

# ***Model Order Reduction of Large Circuits Using Balanced Truncation Via the Arnoldi Method***

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## **Outline**

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- Introduction**
- State Space Form and System Norm**
- Model Reduction Based on Balanced Truncation**
- Numerical Methods**
- Experimental Results**
- Conclusion**

## Introduction

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- ❑ **In high-frequency range, circuits should be modeled as distributed elements**
- ❑ **Extracted circuits are huge and cannot be simulated without order reduction**
- ❑ **AWE is a method for order reduction based on Pade approximation of the system transfer function**
- ❑ **Improvements of AWE include RICE, PVL and PRIMA (guarantees passivity)**

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## Overview

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- ❑ **Model order reduction has extensively been studied in the control engineering**
- ❑ **Both frequency- and time-domain model reduction techniques have been proposed**
- ❑ **An effective method is based on the balanced realization of the system**
  - ◆ **Guaranteed stability**
  - ◆ **Bound on the error for the reduced system**
  - ◆ **Provably optimal solution**

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## State Space Form

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- Any linear, time-invariant circuit can be written in *standard state space form* as :

$$\dot{\vec{x}} = A\vec{x} + B\vec{u}$$

$$\vec{y} = C\vec{x} + D\vec{u}$$

- Given the state space matrix (A,B,C,D), the transfer function of the system is:

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

- Without loss of generality, assume  $D=0$

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## Singular Value Decomposition

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- Any  $l \times m$  matrix  $A$  may be factorized into a singular value decomposition  $A = U\Sigma V^H$  where the  $l \times l$  matrix  $U$  and the  $m \times m$  matrix  $V$  are unitary and the  $l \times m$  matrix  $\Sigma$  contains a diagonal matrix  $\Sigma_1$  of real, non-negative singular values  $\sigma_i$  arranged as  $[\Sigma_1 \ 0]$  if  $l < m$ , as  $\Sigma_1$  if  $l = m$  and as  $[\Sigma_1 \ 0]^T$  otherwise.

**Note that**  $\Sigma_1 = \text{diag}\{\sigma_1, \sigma_2, \dots, \sigma_k\}$  ;  $k = \min(l, m)$   
and  $\sigma_1 \geq \sigma_2 \geq \dots \geq \sigma_k$  the singular values  $\sigma_i$  are positive square roots of the  $k$  largest eigenvalues of both  $AA^H$  and  $A^H A$ . Matrices  $U$  and  $V$  are unit eigen-vectors of  $AA^H$  and  $A^H A$  .

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## Balancing the System

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- The idea is to find a state vector realization of the system that results in equal coupling of energy from the inputs to states and from states to the outputs
- The *Reachability and Observability Gramians* are measures of such energy couplings:

$$W_r = \int_0^{\infty} e^{A\tau} B B^T e^{A^T \tau} d\tau \quad W_o = \int_0^{\infty} e^{A^T \tau} C^T C e^{A\tau} d\tau$$

- The Gramians are obtained by solving the Lyapunov equations

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## Lyapunov Equations

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- The Lyapunov equations can be stated as follows:

$$A W_r + W_r A^T + B B^T = 0$$

$$A^T W_o + W_o A + C^T C = 0$$

- The Hankel singular values and the Hankel norm are then calculated as:

$$\sigma_i(G(s)) = \sqrt{\lambda_i(W_r W_o)}, i = 1, 2, \dots, n \quad \sigma_1 \geq \sigma_2 \geq \dots \geq \sigma_n \geq 0$$

$$\|G(s)\|_H = \max \sigma_i = \sigma_1$$

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## Time-Domain View of Hankel Norm

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- It can be shown that Hankel norm is also given by:

$$\|G(s)\|_H = \max_{\bar{u}(t)} \frac{\sqrt{\int_0^{\infty} \|\bar{y}(t)\|_2^2}}{\sqrt{\int_{-\infty}^0 \|\bar{u}(t)\|_2^2}}$$

- Hankel norm can be interpreted as a kind of induced norm from past inputs to future outputs

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## Balanced Realization (BR)

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- By applying a transformation  $T$  to the system, we can change  $W_r$  and  $W_o$  as follows:

$$\hat{W}_r = T^{-1}W_rT^{-T} \quad \hat{W}_o = T^TW_oT$$

- It can be shown that for any system, there is a transformation which makes

$$\hat{W}_r = \hat{W}_o$$

- Using such a transformation, the new system is called a *balanced realization*

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## Reduced System

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- **To reduce order of the system, we simply ignore states with small Hankel singular values:**

$$A = \begin{bmatrix} A_{11} & A_{12} \\ A_{21} & A_{22} \end{bmatrix}, B = \begin{bmatrix} B_1 \\ B_2 \end{bmatrix}, C = [C_1 \quad C_2] \quad \Sigma = \begin{bmatrix} \Sigma_1 & 0 \\ 0 & \Sigma_2 \end{bmatrix}$$

- **Original System**       $G(s) = C(sI - A)^{-1}B$

- **Reduced System**       $G_a^k(s) = C_1(sI - A_{11})^{-1}B_1$

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## H<sub>∞</sub> System Norm

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- **H<sub>∞</sub> norm of the system is defined as:**

$$\|G(s)\|_{\infty} = \max_{\omega} \sigma_1(G(j\omega))$$

- **Note that given any matrix A,  $\sigma_i$  is the defined as:**  $\sigma_i(A) = \sqrt{\lambda_i(A^H A)} = \sqrt{\lambda_i(AA^H)}$

- **$\sigma_1$  is the largest  $\sigma_i$**

- **It can be shown that:**  $\|G(s)\|_{\infty} = \max_{\vec{u}(t) \neq 0} \frac{\|\vec{y}(t)\|_2}{\|\vec{u}(t)\|_2}$

- **So when  $\|G_1(s) - G_2(s)\|_{\infty} \approx 0$ , the two systems are almost identical**

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## Main Theorems

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- Let  $G(s)$  denote a stable rational transfer function of degree  $n$  with Hankel singular values  $\sigma_1 \geq \sigma_2 \geq \dots \geq \sigma_n \geq 0$ . Let  $G_a^k(s)$  denote the order  $k$  reduction of this transfer function as defined previously. We have:

$$\|G(s) - G_a^k(s)\|_\infty \leq 2(\sigma_{k+1} + \sigma_{k+2} + \dots + \sigma_n)$$

- The minimum error for approximating  $G(s)$  with an arbitrary transfer function  $H(s)$  of degree  $r < n$  is given by:

$$\|G(s) - H(s)\|_\infty \geq \sigma_{k+1}$$

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## Numerical Methods

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- We can directly obtain the reduced order system without calculating the BR

- Procedure (Safanov'89):

- ◆ Compute matrices  $V_{L,k}$  and  $V_{R,k}$  whose columns form bases for the right and left eigen-spaces of  $W_r W_o$  associated with the big eigen-values  $\sigma_1^2, \dots, \sigma_k^2$

- ◆ Set  $E = V_{L,k}^T V_{R,k}$

- ◆ Compute singular value decomposition  $U_E \Sigma_k V_E^T = E$

- ◆ Set  $S_L = V_{L,k} U_E \Sigma^{-1/2} \in \mathfrak{R}^{n \times k}$

$$S_R = V_{R,k} V_E \Sigma^{-1/2} \in \mathfrak{R}^{n \times k}$$

- ◆ The reduced order system is given by:

$$\hat{A} = S_L^T A S_R \quad \hat{B} = S_L^T B \quad \hat{C} = C S_R \quad \hat{D} = D$$

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## Numerical Methods

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- In Safanov's algorithm, we need  $W_r$  and  $W_o$  and then the Schur decomposition of  $W_r W_o$  to obtain  $V_{L,k}$  and  $V_{R,k}$
- Large Lyapunov equations can be solved directly using Krylov-subspace methods (based on the Arnoldi algorithm) as shown in [Saad'89]
- For the decomposition of  $W_r W_o$ , we resort again to the Arnoldi algorithm

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## Balanced Truncation Via Arnoldi

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- Procedure BTVA
  - ◆ Use Krylov-subspace method to calculate  $W_r, W_o$
  - ◆ Use Arnoldi algorithm to calculate big eigenvalues and corresponding left and right eigenvectors ( $V_{r,k}, V_{l,k}$ ) of  $W_r W_o$
  - ◆ Choose the degree for reduced order system based on calculated eigenvalues and the desired error bounds
  - ◆ Compute
 
$$E = V_{L,k}^T V_{R,k} \qquad U_E \Sigma_k V_E^T = E$$

$$S_L = V_{L,k} U_E \Sigma^{-1/2} \in \mathfrak{R}^{n \times k} \qquad S_R = V_{R,k} V_E \Sigma^{-1/2} \in \mathfrak{R}^{n \times k}$$
  - ◆ Compute the reduced order system

$$\hat{A} = S_L^T A S_R \qquad \hat{B} = S_L^T B \qquad \hat{C} = C S_R \qquad \hat{D} = D$$

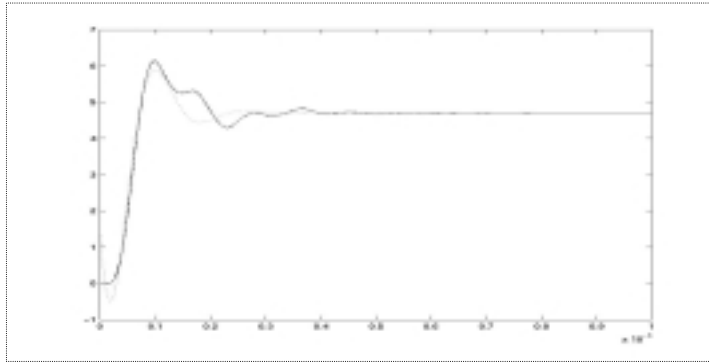
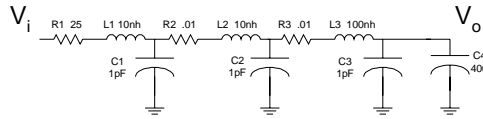
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## Experimental Results

- Small Circuit:
- 8->4 identical
- 8->2 Some error

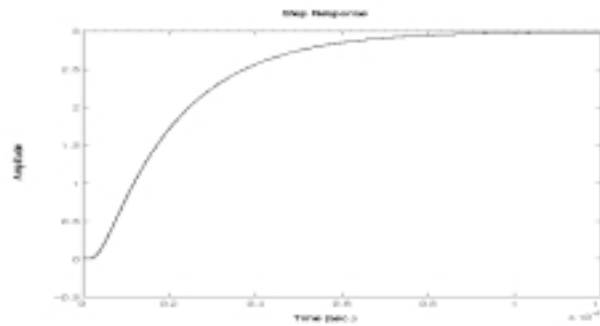
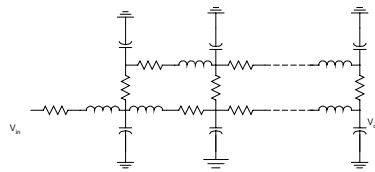


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## Experimental Result

- Big System
- 320-> 4 : Identical



Pc.....

## Conclusions

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- ❑ **Balanced realization is a provably optimal solution to order reduction of LTI systems**
- ❑ **It results in better reduced system compared to the Pade-based techniques**
- ❑ **The computational complexity may however limit the application of this method**
- ❑ **Better methods for solving Lyapunov equations are required to handle higher order systems**