



# Model Reduction of Linear and Nonlinear Control Systems

# Arthur J. Krener

Mathematics University of California, Davis and Applied Mathematics Naval Postgraduate School, Monterey

ajkrener@ucdavis.edu

Research supported in part by NSF and AFOSR

# Model Reduction for Control Systems

#### • Full Order Model

$$egin{array}{rcl} \dot{x}&=&f(x,u)\ &y&=&h(x)\ &u\in{I\!\!R}^m, &y\in&{I\!\!R}^p, &x\in{I\!\!R}^n, &n>>1 \end{array}$$

# Model Reduction for Control Systems

#### • Full Order Model

$$egin{array}{rcl} \dot{x}&=&f(x,u)\ &y&=&h(x)\ u\in{I\!\!R}^m, &y\in&{I\!\!R}^p, &x\in{I\!\!R}^n, &n>>1 \end{array}$$

#### • Reduced Order Model

$$egin{array}{rcl} \dot{z}&=&a(z,u)\ &y&=&c(z)\ &u\in{I\!\!R}^m, &y\in&{I\!\!R}^p, &z\in{I\!\!R}^k, &k<< n \end{array}$$

• The reduced order model should have essentially the same input output behaviour as the full order model.

- The reduced order model should have essentially the same input output behaviour as the full order model.
- A compensator that achieves a desired performance for the reduced order model should also do so for the full order model.

- The reduced order model should have essentially the same input output behaviour as the full order model.
- A compensator that achieves a desired performance for the reduced order model should also do so for the full order model.
- The full order model is a compensator that achieves a desired performance for another system and we seek a reduced order compensator that does also.

- The reduced order model should have essentially the same input output behaviour as the full order model.
- A compensator that achieves a desired performance for the reduced order model should also do so for the full order model.
- The full order model is a compensator that achieves a desired performance for another system and we seek a reduced order compensator that does also.
- We shall focus on the first goal.

#### • Full Order Model

$$egin{array}{rcl} \dot{x}&=&f(x)\ x\in{I\!\!R}^n,&n>>1 \end{array}$$

#### • Full Order Model

$$egin{array}{rcl} \dot{x}&=&f(x)\ x\in{I\!\!R}^n,\qquad n>>1 \end{array}$$

#### • Reduced Order Model

$$egin{array}{rcl} \dot{z}&=&a(z)\ z\in{I\!\!R}^k, \qquad k<< n \end{array}$$

#### • Full Order Model

$$egin{array}{rcl} \dot{x}&=&f(x)\ x\in{I\!\!R}^n,\qquad n>>1 \end{array}$$

#### • Reduced Order Model

$$egin{array}{rcl} \dot{z}&=&a(z)\ z\in{I\!\!R}^k, \qquad k<< n \end{array}$$

• The reduced order model should display the "essential" behaviour of the full order model.

• The model reduction problem for dynamical systems can be viewed as one for control systems by adding an input and output,

$$egin{array}{rcl} \dot{x}&=&f(x)+u\ y&=&x\ u\in{I\!\!R}^n,&y\in{I\!\!R}^n \end{array}$$

 The model reduction problem for dynamical systems can be viewed as one for control systems by adding an input and output,

$$egin{array}{rcl} \dot{x}&=&f(x)+u\ y&=&x\ u\in{I\!\!R}^n,&y\in{I\!\!R}^n \end{array}$$

• But because the input and output dimensions are now large, it may be difficult to reduce the model.

• Separation into slow and fast modes.

$$\left[ egin{array}{cc} \dot{x}_1 \ dots \ \dot{x}_n \end{array} 
ight] \;\; = \;\; \left[ egin{array}{cc} \lambda_1 & & 0 \ & \ddots & \ 0 & & \lambda_n \end{array} 
ight] \left[ egin{array}{cc} x_1 \ dots \ x_n \end{array} 
ight] + ar{f}(x_1,\ldots,x_n)$$

• Separation into slow and fast modes.

$$\left[ egin{array}{cc} \dot{x}_1 \ dots \ \dot{x}_n \end{array} 
ight] \;\; = \;\; \left[ egin{array}{cc} \lambda_1 & & 0 \ & \ddots & \ 0 & & \lambda_n \end{array} 
ight] \left[ egin{array}{cc} x_1 \ dots \ x_n \end{array} 
ight] + ar{f}(x_1,\ldots,x_n)$$

• Spectral Gap:

 $0 \geq \lambda_1 \geq \ldots \geq \lambda_k \;\; >> \;\; \lambda_{k+1} \geq \ldots \geq \lambda_n$ 

• Separation into slow and fast modes.

$$\left[ egin{array}{cc} \dot{x}_1 \ dots \ \dot{x}_n \end{array} 
ight] \;\; = \;\; \left[ egin{array}{cc} \lambda_1 & & 0 \ & \ddots & \ 0 & & \lambda_n \end{array} 
ight] \left[ egin{array}{cc} x_1 \ dots \ x_n \end{array} 
ight] + ar{f}(x_1,\ldots,x_n)$$

• Spectral Gap:

$$0 \geq \lambda_1 \geq \ldots \geq \lambda_k \;\; >> \;\; \lambda_{k+1} \geq \ldots \geq \lambda_n$$

• Galerkin projection onto  $x_{k+1} = \cdots = x_n = 0$ ,

$$\left[ egin{array}{cc} \dot{x}_1 \ dots \ \dot{x}_k \end{array} 
ight] \;\;=\;\; \left[ egin{array}{cc} \lambda_1 & & 0 \ & \ddots & \ 0 & & \lambda_k \end{array} 
ight] \left[ egin{array}{cc} x_1 \ dots \ x_k \end{array} 
ight] + ar{f}_1(x_1 \dots, x_k, 0, \dots, 0)$$

• Separation into slow and fast modes.

$$\left[ egin{array}{cc} \dot{x}_1 \ dots \ \dot{x}_n \end{array} 
ight] \;\; = \;\; \left[ egin{array}{cc} \lambda_1 & & 0 \ & \ddots & \ 0 & & \lambda_n \end{array} 
ight] \left[ egin{array}{cc} x_1 \ dots \ x_n \end{array} 
ight] + ar{f}(x_1,\ldots,x_n)$$

• Spectral Gap:

$$0 \geq \lambda_1 \geq \ldots \geq \lambda_k \;\; >> \;\; \lambda_{k+1} \geq \ldots \geq \lambda_n$$

• Galerkin projection onto  $x_{k+1} = \cdots = x_n = 0,$ 

 $\begin{bmatrix} \dot{x}_1 \\ \vdots \\ \dot{x}_k \end{bmatrix} = \begin{bmatrix} \lambda_1 & 0 \\ & \ddots & \\ 0 & & \lambda_k \end{bmatrix} \begin{bmatrix} x_1 \\ \vdots \\ x_k \end{bmatrix} + \bar{f}_1(x_1 \dots, x_k, 0, \dots, 0)$ 

• Other approaches: Petrov Galerkin, nonlinear Galerkin, singular perturbations, center manifolds, inertial manifolds..

Galerkin projection onto the slow modes of the unforced dynamics may not be a satisfactory method of model reduction for control systems.

Galerkin projection onto the slow modes of the unforced dynamics may not be a satisfactory method of model reduction for control systems.

The control may not directly excite the slow modes.

Galerkin projection onto the slow modes of the unforced dynamics may not be a satisfactory method of model reduction for control systems.

The control may not directly excite the slow modes.

The output may not be sensitive to changes in the slow modes.

• We shall focus on state space methods which generalize to nonlinear control systems.

$$egin{array}{rcl} \dot{x}&=&Fx+Gu\ y&=&Hx\ x(0)&=&0\ x\in{I\!\!R}^n,&u&\in{I\!\!R}^m,&y\in{I\!\!R}^p\ \lambda(F)&<&0 \end{array}$$

• We shall focus on state space methods which generalize to nonlinear control systems.

$$egin{array}{rcl} \dot{x}&=&Fx+Gu\ y&=&Hx\ x(0)&=&0\ x\in{I\!\!R}^n, & u&\in&{I\!\!R}^m, & y\in{I\!\!R}^p\ \lambda(F)&<&0 \end{array}$$

• Since the unforced system is Hurwitz, it defines an input-output map

$$egin{array}{rcl} \mathcal{IO}_n: L^2(-\infty,\infty;{I\!\!R}^m) & o & L^2(-\infty,\infty;{I\!\!R}^p) \ \mathcal{IO}_n: u(-\infty:\infty) & \mapsto & y(-\infty:\infty) \ y(t) & = & \int_{-\infty}^t He^{F(t-s)}Gu(s) \; ds \end{array}$$

• What is the smallest state dimension necessary to realize a given input-output map?

- What is the smallest state dimension necessary to realize a given input-output map?
- The state dimension is minimal if the system is controllable,

$$\operatorname{rank}\left[ egin{array}{cccc} G & FG & \ldots & F^{n-1}G \end{array} 
ight] = n,$$

- What is the smallest state dimension necessary to realize a given input-output map?
- The state dimension is minimal if the system is controllable,

$$\operatorname{rank} \left[ \begin{array}{ccc} G & FG & \dots & F^{n-1}G \end{array} \right] = n,$$

and observable,

$$\operatorname{rank} \left[ egin{array}{c} H \ HF \ dots \ HF^{n-1} \end{array} 
ight] = n.$$

- What is the smallest state dimension necessary to realize a given input-output map?
- The state dimension is minimal if the system is controllable,

$$\operatorname{rank} \left[ \begin{array}{ccc} G & FG & \dots & F^{n-1}G \end{array} 
ight] = n,$$

and observable,

$$\operatorname{rank} \begin{bmatrix} H \\ HF \\ \vdots \\ HF^{n-1} \end{bmatrix} = n.$$

• Any system can be reduced to one that is minimal.

• The input-output map is not a compact operator.

- The input-output map is not a compact operator.
- The Hankel map takes past inputs to future outputs

$$egin{array}{rcl} \mathcal{H}_n: L^2(-\infty,0;{I\!\!R}^m) &
ightarrow L^2(0,\infty;{I\!\!R}^p) \ \mathcal{H}_n: u(-\infty:0) &\mapsto y(0:\infty) \end{array}$$

- The input-output map is not a compact operator.
- The Hankel map takes past inputs to future outputs

$$egin{array}{lll} \mathcal{H}_n: L^2(-\infty,0;{I\!\!R}^m) &
ightarrow L^2(0,\infty;{I\!\!R}^p) \ \mathcal{H}_n: u(-\infty:0) &\mapsto y(0:\infty) \end{array}$$

• It factors through the current state x(0) and so it is of finite rank hence compact.

$$u(-\infty:0) \mapsto x(0) = \int_{-\infty}^{0} e^{-Fs} Gu(s) ds$$
  
 $y(t) = He^{Ft} x(0)$ 

$$\dot{x} = Fx + Gu$$
  
 $y = Hx$ 

• Find a reduced order linear system with approximately the same input-output map.

$$\dot{x} = Fx + Gu$$
  
 $y = Hx$ 

- Find a reduced order linear system with approximately the same input-output map.
- Moore assumed that

$$\dot{x} = Fx + Gu$$
  
 $y = Hx$ 

- Find a reduced order linear system with approximately the same input-output map.
- Moore assumed that
- F, G is a controllable pair,

$$egin{array}{rcl} \dot{x}&=&Fx+Gu\ y&=&Hx \end{array}$$

- Find a reduced order linear system with approximately the same input-output map.
- Moore assumed that
- F, G is a controllable pair,
- *H*, *F* is a observable pair,

$$egin{array}{rcl} \dot{x}&=&Fx+Gu\ y&=&Hx \end{array}$$

- Find a reduced order linear system with approximately the same input-output map.
- Moore assumed that
- F, G is a controllable pair,
- *H*, *F* is a observable pair,
- F is Hurwitz,  $\lambda(F) < 0$  .

$$\dot{x} = Fx + Gu$$
  
 $y = Hx$ 

- Find a reduced order linear system with approximately the same input-output map.
- Moore assumed that
- F, G is a controllable pair,
- *H*, *F* is a observable pair,
- F is Hurwitz,  $\lambda(F) < 0$  .
- If the system is uncontrollable and/or unobservable, we can make it so by passing to a minimal realization.

$$egin{array}{rcl} \dot{x}&=&Fx+Gu\ y&=&Hx \end{array}$$

- Find a reduced order linear system with approximately the same input-output map.
- Moore assumed that
- F, G is a controllable pair,
- *H*, *F* is a observable pair,
- F is Hurwitz,  $\lambda(F) < 0$  .
- If the system is uncontrollable and/or unobservable, we can make it so by passing to a minimal realization.
- Hurwitz is needed to insure the existence of the input-output map.

## Balanced Realization Theory.

• Moore's insight was that we should restrict to the directions that are easy to excite and ignore directions where changes don't affect the output very much.
- Moore's insight was that we should restrict to the directions that are easy to excite and ignore directions where changes don't affect the output very much.
- Controllablity Function

$$\pi_c(x^0) \;\; = \;\; \inf_{u(-\infty:0)} rac{1}{2} \int_{-\infty}^0 |u(t)|^2 \; dt$$

subject to the system dynamics and  $x(-\infty)=0,\;x(0)=x^0.$ 

- Moore's insight was that we should restrict to the directions that are easy to excite and ignore directions where changes don't affect the output very much.
- Controllablity Function

$$\pi_c(x^0) \;\; = \;\; \inf_{u(-\infty:0)} rac{1}{2} \int_{-\infty}^0 |u(t)|^2 \; dt$$

subject to the system dynamics and  $x(-\infty)=0,\;x(0)=x^0.$ 

• Observability Function

$$\pi_o(x^0) \;\;=\;\; rac{1}{2} \int_0^\infty |y(t)|^2 \; dt$$

subject to the system dynamics and  $x(0) = x^0, \ u(t) = 0, \ t \ge 0.$ 

•  $\pi_c(x)$  is the minimal input energy needed to excite the system from the zero state to x.

- $\pi_c(x)$  is the minimal input energy needed to excite the system from the zero state to x.
- F, G controllable implies  $\pi_c(x)$  is bounded.

- $\pi_c(x)$  is the minimal input energy needed to excite the system from the zero state to x.
- F, G controllable implies  $\pi_c(x)$  is bounded.
- F Hurwitz implies  $\pi_c(x)$  is positive definite.

- $\pi_c(x)$  is the minimal input energy needed to excite the system from the zero state to x.
- F, G controllable implies  $\pi_c(x)$  is bounded.
- *F* Hurwitz implies  $\pi_c(x)$  is positive definite.
- $\pi_o(x)$  is the output energy released by the system as it decays from x back to the zero state.

- $\pi_c(x)$  is the minimal input energy needed to excite the system from the zero state to x.
- F, G controllable implies  $\pi_c(x)$  is bounded.
- *F* Hurwitz implies  $\pi_c(x)$  is positive definite.
- $\pi_o(x)$  is the output energy released by the system as it decays from x back to the zero state.
- *F* Hurwitz implies  $\pi_o(x)$  is bounded.

- $\pi_c(x)$  is the minimal input energy needed to excite the system from the zero state to x.
- F, G controllable implies  $\pi_c(x)$  is bounded.
- *F* Hurwitz implies  $\pi_c(x)$  is positive definite.
- $\pi_o(x)$  is the output energy released by the system as it decays from x back to the zero state.
- *F* Hurwitz implies  $\pi_o(x)$  is bounded.
- H, F observable implies  $\pi_o(x)$  is positive definite.

- $\pi_c(x)$  is the minimal input energy needed to excite the system from the zero state to x.
- F, G controllable implies  $\pi_c(x)$  is bounded.
- *F* Hurwitz implies  $\pi_c(x)$  is positive definite.
- $\pi_o(x)$  is the output energy released by the system as it decays from x back to the zero state.
- *F* Hurwitz implies  $\pi_o(x)$  is bounded.
- H, F observable implies  $\pi_o(x)$  is positive definite.
- $\pi_c(x)$  and  $\pi_o(x)$  are quadratic functions because the system is linear and the energies are quadratic,

$$\pi_c(x) = rac{1}{2} x' P_c^{-1} x, \qquad \pi_o(x) = rac{1}{2} x' P_o x$$

#### • P<sub>c</sub>, P<sub>o</sub> are the unique positive definite solutions of

$$0 = FP_c + P_cF' + GG'$$
  
$$0 = F'P_o + P_oF + H'H.$$

#### • P<sub>c</sub>, P<sub>o</sub> are the unique positive definite solutions of

$$0 = FP_c + P_cF' + GG'$$
  
$$0 = F'P_o + P_oF + H'H.$$

• They are called the controllablity and observability gramians.

• P<sub>c</sub>, P<sub>o</sub> are the unique positive definite solutions of

$$0 = FP_c + P_cF' + GG'$$
  
$$0 = F'P_o + P_oF + H'H.$$

- They are called the controllablity and observability gramians.
- $\pi_c(x)$  large implies that it takes a lot of input energy to excite the system in the direction x and so this direction might be ignored in a reduced order model.

• P<sub>c</sub>, P<sub>o</sub> are the unique positive definite solutions of

$$0 = FP_c + P_cF' + GG'$$
  
$$0 = F'P_o + P_oF + H'H.$$

- They are called the controllablity and observability gramians.
- $\pi_c(x)$  large implies that it takes a lot of input energy to excite the system in the direction x and so this direction might be ignored in a reduced order model.
- $\pi_o(x)$  small implies that changes in the direction x lead to small changes in the output energy and so this direction might be ignored in a reduced order model.

•  $P_c$ ,  $P_o$  transform differently under a linear change of states coordinates x = Tz

$$\begin{array}{rccc} P_c & \mapsto & T^{-1}P_cT'^{-1} \\ P_o & \mapsto & T'P_oT \end{array}$$

•  $P_c$ ,  $P_o$  transform differently under a linear change of states coordinates x = Tz

$$\begin{array}{rccc} P_c & \mapsto & T^{-1}P_cT'^{-1} \\ P_o & \mapsto & T'P_oT \end{array}$$

• P<sub>o</sub>P<sub>c</sub> is a similarity invariant

$$P_oP_c \mapsto T'P_oP_cT'^{-1}$$

•  $P_c$ ,  $P_o$  transform differently under a linear change of states coordinates x = Tz

$$\begin{array}{rccc} P_c & \mapsto & T^{-1}P_cT'^{-1} \\ P_o & \mapsto & T'P_oT \end{array}$$

• P<sub>o</sub>P<sub>c</sub> is a similarity invariant

$$P_oP_c \mapsto T'P_oP_cT'^{-1}$$

• Its eigenvalues are the squares of the singular values of the Hankel map.

• There is a linear change of state coordinates so that the controllability and observability gramians are diagonal and equal,

$$P_c = P_o = \left[egin{array}{cc} \sigma_1 & 0 \ & \ddots & \ 0 & \sigma_n \end{array}
ight]$$

• There is a linear change of state coordinates so that the controllability and observability gramians are diagonal and equal,

$$P_c = P_o = \left[ egin{array}{cc} \sigma_1 & 0 \ & \ddots & \ 0 & \sigma_n \end{array} 
ight]$$

• The Hankel singular values can be ordered

$$\sigma_1 \geq \sigma_2 \geq \ldots \geq \sigma_n > 0$$

• There is a linear change of state coordinates so that the controllability and observability gramians are diagonal and equal,

$$P_c = P_o = \left[egin{array}{cc} \sigma_1 & 0 \ & \ddots & \ 0 & \sigma_n \end{array}
ight]$$

• The Hankel singular values can be ordered

$$\sigma_1 \geq \sigma_2 \geq \ldots \geq \sigma_n > 0$$

• In these new state coordinates the system is said to be balanced.

• There is a linear change of state coordinates so that the controllability and observability gramians are diagonal and equal,

$$P_c = P_o = \left[egin{array}{cc} \sigma_1 & 0 \ & \ddots & \ 0 & \sigma_n \end{array}
ight]$$

• The Hankel singular values can be ordered

$$\sigma_1 \geq \sigma_2 \geq \ldots \geq \sigma_n > 0$$

- In these new state coordinates the system is said to be balanced.
- If the Hankel singular values are distinct then the balanced coordinates are unique up to changes of signs  $x_i\mapsto -x_i$ .

• The reduced model is obtained by Galerkin projection onto the states corresponding to large Hankel singular values because they can be reached with relatively small input energy and they produce relatively large output energy.

- The reduced model is obtained by Galerkin projection onto the states corresponding to large Hankel singular values because they can be reached with relatively small input energy and they produce relatively large output energy.
- Suppose  $\sigma_k >> \sigma_{k+1}$ . Let  $x_1$  denote the first k components of xLet  $x_2$  denote the last n - k components.

- The reduced model is obtained by Galerkin projection onto the states corresponding to large Hankel singular values because they can be reached with relatively small input energy and they produce relatively large output energy.
- Suppose  $\sigma_k >> \sigma_{k+1}$ . Let  $x_1$  denote the first k components of xLet  $x_2$  denote the last n - k components.
- Full Order Model

$$egin{array}{rcl} \dot{x}_1 \ \dot{x}_2 \end{array} &=& \left[ egin{array}{cc} F_{11} & F_{12} \ F_{21} & F_{22} \end{array} 
ight] \left[ egin{array}{cc} x_1 \ x_2 \end{array} 
ight] + \left[ egin{array}{cc} G_1 \ G_2 \end{array} 
ight] u \ y &=& \left[ egin{array}{cc} H_1 & H_2 \end{array} 
ight] \left[ egin{array}{cc} x_1 \ x_2 \end{array} 
ight] \end{array}$$

- The reduced model is obtained by Galerkin projection onto the states corresponding to large Hankel singular values because they can be reached with relatively small input energy and they produce relatively large output energy.
- Suppose  $\sigma_k >> \sigma_{k+1}$ . Let  $x_1$  denote the first k components of xLet  $x_2$  denote the last n - k components.
- Full Order Model

$$egin{array}{ccc} \left[ egin{array}{c} \dot{x}_1 \ \dot{x}_2 \end{array} 
ight] &=& \left[ egin{array}{ccc} F_{11} & F_{12} \ F_{21} & F_{22} \end{array} 
ight] \left[ egin{array}{c} x_1 \ x_2 \end{array} 
ight] + \left[ egin{array}{c} G_1 \ G_2 \end{array} 
ight] u \ y &=& \left[ egin{array}{c} H_1 & H_2 \end{array} 
ight] \left[ egin{array}{c} x_1 \ x_2 \end{array} 
ight] \end{array}$$

• Balanced Truncation obtained by Galerkin projection.

$$\dot{z} = F_{11}z + G_1u$$
  
 $y = H_1z$ 

• The reduced model is obtained by Galerkin (orthogonal) projection in balanced coordinates.

- The reduced model is obtained by Galerkin (orthogonal) projection in balanced coordinates.
- In the original coordinates, it is a Petrov Galerkin (oblique) projection.

- The reduced model is obtained by Galerkin (orthogonal) projection in balanced coordinates.
- In the original coordinates, it is a Petrov Galerkin (oblique) projection.
- The Hankel singular values of the reduced model are  $\sigma_1, \ldots, \sigma_k$ .

- The reduced model is obtained by Galerkin (orthogonal) projection in balanced coordinates.
- In the original coordinates, it is a Petrov Galerkin (oblique) projection.
- The Hankel singular values of the reduced model are  $\sigma_1, \ldots, \sigma_k$ .
- But the reduced model is not an optimal Hankel norm approximation of the full model because the singular vectors are different. Typically

$$\|\mathcal{H}_n - \mathcal{H}_k\| > \sigma_{k+1}$$

Adamjan-Arov-Krein, Glover

- The reduced model is obtained by Galerkin (orthogonal) projection in balanced coordinates.
- In the original coordinates, it is a Petrov Galerkin (oblique) projection.
- The Hankel singular values of the reduced model are  $\sigma_1, \ldots, \sigma_k$ .
- But the reduced model is not an optimal Hankel norm approximation of the full model because the singular vectors are different. Typically

$$\|\mathcal{H}_n - \mathcal{H}_k\| > \sigma_{k+1}$$

Adamjan-Arov-Krein, Glover

• Glover has shown that for balanced truncation

$$\left\|\mathcal{IO}_n-\mathcal{IO}_k
ight\|\leq 2\sum_{j=k+1}^n\sigma_j$$

• Large and small are relative terms and we need one quadratic form to normalize the other.

- Large and small are relative terms and we need one quadratic form to normalize the other.
- The eigenvalues of the dynamics play an indirect role. It is very hard to excite the system in a direction corresponding to a very stable eigenvalue and so the controllability function tends to be large in such a direction.

- Large and small are relative terms and we need one quadratic form to normalize the other.
- The eigenvalues of the dynamics play an indirect role. It is very hard to excite the system in a direction corresponding to a very stable eigenvalue and so the controllability function tends to be large in such a direction.
- Moreover, a very stable state direction damps out quickly and so the observability function tends to be small in such a direction.

- Large and small are relative terms and we need one quadratic form to normalize the other.
- The eigenvalues of the dynamics play an indirect role. It is very hard to excite the system in a direction corresponding to a very stable eigenvalue and so the controllability function tends to be large in such a direction.
- Moreover, a very stable state direction damps out quickly and so the observability function tends to be small in such a direction.
- Hence the very stable directions of the dynamics tend to be ignored in the reduction process.

#### **Balanced Reduction**

Suppose we have a linear dyamical system in modal coordinates

$$\dot{x} \;=\; \left[egin{array}{ccc} \lambda_1 & & 0 \ & \ddots & \ 0 & & \lambda_n \end{array}
ight]x$$

 $0>\lambda_1\geq\ldots\geq\lambda_n$  .

#### **Balanced Reduction**

#### Suppose we have a linear dyamical system in modal coordinates

$$\dot{x} \;=\; \left[egin{array}{ccc} \lambda_1 & & 0 \ & \ddots & \ 0 & & \lambda_n \end{array}
ight]x$$

 $0>\lambda_1\geq\ldots\geq\lambda_n$  .

#### As before we add a dummy input and output,

$$\dot{x} = \left[egin{array}{ccc} \lambda_1 & & 0 \ & \ddots & \ 0 & & \lambda_n \end{array}
ight]x+u, \quad y=x$$

# **Balanced Reduction**

#### Then

$$P_c=P_o=\left[egin{array}{cc} -rac{1}{2\lambda_1}&0\ &\ &\ddots\ &\ &\ &0&-rac{1}{2\lambda_n}\end{array}
ight], \quad \sigma_i=-rac{1}{2\lambda_i}$$
# **Balanced Reduction**

#### Then

$$P_c=P_o=\left[egin{array}{cc} -rac{1}{2\lambda_1}&0\ &\ &\ddots\ &\ &\ &0&-rac{1}{2\lambda_n}\end{array}
ight], \quad \sigma_i=-rac{1}{2\lambda_i}$$

So balanced truncation is the usual truncation onto the slow modes.

$$-rac{1}{2\lambda_1}\geq -rac{1}{2\lambda_2}\geq \ldots \geq -rac{1}{2\lambda_n}>0$$

Chain of three masses connected by springs and dashpots attached to a wall at one end. The input is a force applied to the mass next to the wall and the output is the displacement of the mass at the other end. Assume that each mass is  $\mu$ , each spring constant is c and each dampening constant is b.



#### The system is linear,

$$F = \begin{bmatrix} 0 & 0 & 0 & 1 & 0 & 0 \\ 0 & 0 & 0 & 0 & 1 & 0 \\ 0 & 0 & 0 & 0 & 0 & 1 \\ -\frac{2c}{m} & \frac{c}{\mu} & 0 & -\frac{2b}{\mu} & \frac{b}{\mu} & 0 \\ \frac{c}{\mu} & -\frac{2c}{\mu} & \frac{c}{\mu} & \frac{b}{\mu} & -\frac{2b}{\mu} & \frac{b}{\mu} \\ 0 & \frac{c}{\mu} & -\frac{c}{\mu} & 0 & \frac{b}{\mu} & -\frac{b}{\mu} \end{bmatrix} \qquad G = \begin{bmatrix} 0 \\ 0 \\ 0 \\ \frac{1}{\mu} \\ 0 \\ 0 \end{bmatrix}$$

 $H = \left[\begin{array}{rrrrr} 0 & 0 & 1 & 0 & 0 \end{array}\right]$ 





$$\sigma_2 = 1.4901 >> 0.1404 = \sigma_3$$

This suggests taking a reduced order model of dimension k=2 .



#### Glover has shown that for balanced truncation

$$\left\|\mathcal{IO}_n-\mathcal{IO}_k
ight\|\leq 2\sum_{j=k+1}^n\sigma_j$$

#### Glover has shown that for balanced truncation

$$\left\|\mathcal{IO}_n-\mathcal{IO}_k
ight\|\leq 2\sum_{j=k+1}^n\sigma_j$$

We know that

$$\sigma_{k+1} \leq \|\mathcal{H}_n - \mathcal{H}_k\| \leq \|\mathcal{IO}_n - \mathcal{IO}_k\| \leq 2\sum_{j=k+1}^n \sigma_j$$

#### Glover has shown that for balanced truncation

$$\left\|\mathcal{IO}_n-\mathcal{IO}_k
ight\|\leq 2\sum_{j=k+1}^n\sigma_j$$

We know that

$$\sigma_{k+1} \leq \|\mathcal{H}_n - \mathcal{H}_k\| \leq \|\mathcal{IO}_n - \mathcal{IO}_k\| \leq 2\sum_{j=k+1}^n \sigma_j$$

For the spring mass example with k = 2 this yields

 $\sigma_3 = 0.1404 \leq \|\mathcal{H}_n - \mathcal{H}_k\| \leq \|\mathcal{IO}_n - \mathcal{IO}_k\| \leq 0.5672$ 

If we restrict the Hankel maps to optimal inputs of the full system then

$$\|\mathcal{H}_n - \mathcal{H}_k\| \leq 0.1432$$

Notice how close this is to  $\sigma_3 = 0.1404$ .



If we restrict the Hankel maps to optimal inputs of the reduced system then

$$\|\mathcal{H}_n-\mathcal{H}_k\|\leq 0.0867$$

Notice how much smaller this is than  $\sigma_3 = 0.1404$ .



• Balanced truncation does not minimize the difference between the input-output maps of the full and reduced models.

- Balanced truncation does not minimize the difference between the input-output maps of the full and reduced models.
- Balanced truncation does not minimize the difference between the Hankel maps of the full and reduced models.

- Balanced truncation does not minimize the difference between the input-output maps of the full and reduced models.
- Balanced truncation does not minimize the difference between the Hankel maps of the full and reduced models.
- So what problem does it solve? And how can it be generalized to nonlinear systems?

- Balanced truncation does not minimize the difference between the input-output maps of the full and reduced models.
- Balanced truncation does not minimize the difference between the Hankel maps of the full and reduced models.
- So what problem does it solve? And how can it be generalized to nonlinear systems?
- Nonlinear balanced truncation of Scherpen.

- Balanced truncation does not minimize the difference between the input-output maps of the full and reduced models.
- Balanced truncation does not minimize the difference between the Hankel maps of the full and reduced models.
- So what problem does it solve? And how can it be generalized to nonlinear systems?
- Nonlinear balanced truncation of Scherpen.
- Stochastic interpretation of Newman and Krishnaprasad.

- Balanced truncation does not minimize the difference between the input-output maps of the full and reduced models.
- Balanced truncation does not minimize the difference between the Hankel maps of the full and reduced models.
- So what problem does it solve? And how can it be generalized to nonlinear systems?
- Nonlinear balanced truncation of Scherpen.
- Stochastic interpretation of Newman and Krishnaprasad.
- Differential eigenstructure of nonlinear Hankel maps, Fujimoto and Scherpen.

- Balanced truncation does not minimize the difference between the input-output maps of the full and reduced models.
- Balanced truncation does not minimize the difference between the Hankel maps of the full and reduced models.
- So what problem does it solve? And how can it be generalized to nonlinear systems?
- Nonlinear balanced truncation of Scherpen.
- Stochastic interpretation of Newman and Krishnaprasad.
- Differential eigenstructure of nonlinear Hankel maps, Fujimoto and Scherpen.
- Here is a new way of viewing and generalizing linear balanced truncation.

**Full Order Model** 

$$\dot{x} = Fx + Gu$$
  
 $y = Hx$ 

Full Order Model

$$\dot{x} = Fx + Gu$$
  
 $y = Hx$ 

We restrict to those reduced order models that can be obtained by Petrov Galerkin projection. For this we need an injection  $\Psi$  and a surjection  $\Phi$ 

$$egin{array}{rcl} \Psi:I\!\!R^k& o&I\!\!R^n\ \Psi:z&\mapsto &x=\Psi z\ \Phi:I\!\!R^n& o&I\!\!R^k\ \Phi:x&\mapsto &z=\Phi x\ \Phi\Psi z=z, &(\Psi\Phi)^2x=\Psi\Phi x \end{array}$$

Full Order Model

$$\dot{x} = Fx + Gu$$
  
 $y = Hx$ 

We restrict to those reduced order models that can be obtained by Petrov Galerkin projection. For this we need an injection  $\Psi$  and a surjection  $\Phi$ 

$$\begin{split} \Psi : I\!\!R^k &\to I\!\!R^n \\ \Psi : z &\mapsto x = \Psi z \\ \Phi : I\!\!R^n &\to I\!\!R^k \\ \Phi : x &\mapsto z = \Phi x \\ \Phi \Psi z = z, \qquad (\Psi \Phi)^2 x = \Psi \Phi x \end{split}$$

**Reduced Order Model** 

$$\dot{z} = \Phi F \Psi z + \Phi G u$$
  
 $y = H \Psi z$ 

How should we choose the injection  $\Psi$  ?

How should we choose the injection  $\Psi$  ?

We choose  $\Psi : \mathbb{R}^k \to \mathbb{R}^n$  so that its k dimensional range maximizes output energy  $\pi_o(x)$  for given input energy  $\pi_c(x)$ .

How should we choose the injection  $\Psi$  ?

We choose  $\Psi : \mathbb{R}^k \to \mathbb{R}^n$  so that its k dimensional range maximizes output energy  $\pi_o(x)$  for given input energy  $\pi_c(x)$ .

To do this it is convenient to put the system in input normal form, that is, make a linear change of state coordinates so that

$$\pi_c(x) = rac{1}{2} \sum_i x_i^2, \quad \pi_o(x) = rac{1}{2} \sum_i au_i x_i^2$$

How should we choose the injection  $\Psi$  ?

We choose  $\Psi : \mathbb{R}^k \to \mathbb{R}^n$  so that its k dimensional range maximizes output energy  $\pi_o(x)$  for given input energy  $\pi_c(x)$ .

To do this it is convenient to put the system in input normal form, that is, make a linear change of state coordinates so that

$$\pi_c(x)=rac{1}{2}\sum_i x_i^2, \quad \ \pi_o(x)=rac{1}{2}\sum_i au_i x_i^2$$

This is just a diagonal change from balanced coordinates and

 $\tau_i = \sigma_i^2$  are the squared Hankel singular values.

How should we choose the injection  $\Psi$  ?

We choose  $\Psi : \mathbb{R}^k \to \mathbb{R}^n$  so that its k dimensional range maximizes output energy  $\pi_o(x)$  for given input energy  $\pi_c(x)$ .

To do this it is convenient to put the system in input normal form, that is, make a linear change of state coordinates so that

$$\pi_c(x)=rac{1}{2}\sum_i x_i^2, \quad \ \pi_o(x)=rac{1}{2}\sum_i au_i x_i^2$$

This is just a diagonal change from balanced coordinates and

 $\tau_i = \sigma_i^2$  are the squared Hankel singular values.

If  $au_k>> au_{k+1}$  then we should take the range of  $\Psi(z)$  to be  $x_{k+1}=\dots=x_n=0$  , e.g.,

$$\Psi(z_1,\ldots,z_k)= egin{array}{cc} x &=(z_1,\ldots,z_k,0,\ldots,0) \end{array}$$

How should we choose the surjection  $\Phi$  ?

How should we choose the surjection  $\Phi$  ?

We choose  $\Phi : \mathbb{R}^n \to \mathbb{R}^k$  to minimize the  $L^2$  norm of the difference between the outputs from x and  $\Psi(\Phi(x))$ .

How should we choose the surjection  $\Phi$  ?

We choose  $\Phi : \mathbb{R}^n \to \mathbb{R}^k$  to minimize the  $L^2$  norm of the difference between the outputs from x and  $\Psi(\Phi(x))$ .

Define the co-observability function

$$\pi_{oo}(x,ar{x}) = rac{1}{2}\int_0^\infty |y(t)-ar{y}(t)|^2 \; dt$$

where  $y(t), \ ar{y}(t)$  are the outputs from  $x(0)=x, \ x(0)=ar{x}.$ 

How should we choose the surjection  $\Phi$  ?

We choose  $\Phi : \mathbb{R}^n \to \mathbb{R}^k$  to minimize the  $L^2$  norm of the difference between the outputs from x and  $\Psi(\Phi(x))$ .

Define the co-observability function

$$\pi_{oo}(x,ar{x}) = rac{1}{2}\int_0^\infty |y(t)-ar{y}(t)|^2 \; dt$$

where  $y(t), \ \bar{y}(t)$  are the outputs from  $x(0) = x, \ x(0) = \bar{x}$ . Because the system is linear

$$\pi_{oo}(x,\bar{x}) = \pi_o(x-\bar{x})$$

How should we choose the surjection  $\Phi$  ?

We choose  $\Phi : \mathbb{R}^n \to \mathbb{R}^k$  to minimize the  $L^2$  norm of the difference between the outputs from x and  $\Psi(\Phi(x))$ .

Define the co-observability function

$$\pi_{oo}(x,ar{x}) = rac{1}{2}\int_0^\infty |y(t)-ar{y}(t)|^2 \; dt$$

where  $y(t), \ \bar{y}(t)$  are the outputs from  $x(0) = x, \ x(0) = \bar{x}$ . Because the system is linear

$$\pi_{oo}(x, \bar{x}) = \pi_o(x - \bar{x})$$

If the system is in input normal form then the optimal  $\Phi$  is

$$\Phi(x_1,\ldots,x_n) = z = (x_1,\ldots,x_k)$$

#### Scherpen generalized Moore to nonlinear systems

$$\dot{x} = f(x,u)$$
  
 $y = h(x)$ 

#### Scherpen generalized Moore to nonlinear systems

$$\dot{x} = f(x,u)$$
  
 $y = h(x)$ 

#### She defined the controllability function,

$$\pi_c(x^0) \;\;=\;\; \inf rac{1}{2} \int_{-\infty}^0 |u(t)|^2 \; dt$$

subject to the system dynamics and  $x(-\infty)=0, \ x(0)=x^0.$ 

#### Scherpen generalized Moore to nonlinear systems

$$\dot{x} = f(x,u)$$
  
 $y = h(x)$ 

#### She defined the controllability function,

$$\pi_c(x^0) \;\; = \;\; \inf rac{1}{2} \int_{-\infty}^0 |u(t)|^2 \; dt$$

subject to the system dynamics and  $x(-\infty) = 0$ ,  $x(0) = x^0$ . And the observability function,

$$\pi_o(x^0) \;\;=\;\; rac{1}{2} \int_0^\infty |y(t)|^2 \; dt$$

subject to the system dynamics and  $x(0)=x^0,\ u(t)=0.$ 

The controllability function  $\pi_c(x)$  and the optimal control  $u = \kappa(x)$  satisfy the HJB PDE

$$\begin{array}{lcl} 0 & = & \displaystyle \frac{\partial \pi_c}{\partial x}(x) f(x,\kappa(x)) - \displaystyle \frac{1}{2} |\kappa(x)|^2 \\ 0 & = & \displaystyle \frac{\partial \pi_c}{\partial x}(x) \displaystyle \frac{\partial f}{\partial u}(x,\kappa(x)) - \kappa'(x) \end{array}$$

The controllability function  $\pi_c(x)$  and the optimal control  $u = \kappa(x)$  satisfy the HJB PDE

$$egin{array}{rcl} 0&=&rac{\partial\pi_c}{\partial x}(x)f(x,\kappa(x))-rac{1}{2}|\kappa(x)|^2\ 0&=&rac{\partial\pi_c}{\partial x}(x)rac{\partial f}{\partial u}(x,\kappa(x))-\kappa'(x) \end{array}$$

The observability function  $\pi_o(x)$  satisfies the Lyapunov PDE

$$0 = \frac{\partial \pi_o}{\partial x}(x)f(x,0) + \frac{1}{2}h'(x)h(x).$$
#### Suppose

• The system is smooth with Taylor expansion

$$\dot{x} \;\;=\;\; f(x,u) \;\;=\;\; Fx + Gu \;\;+ f^{[2]}(x,u) \;\;+ \ldots$$

$$y = h(x) = Hx + h^{[2]}(x) + \dots$$

where [d] denotes a vector field that is a homogeneous polynomial of degree d.

#### Suppose

• The system is smooth with Taylor expansion

$$\dot{x} \;\;=\;\; f(x,u) \;\;=\;\; Fx + Gu \;\;+ f^{[2]}(x,u) \;\;+ \dots$$

where [d] denotes a vector field that is a homogeneous polynomial of degree d.

• The linear part of the system is Hurwitz, controllable and observable

#### Then

• there exist smooth, positive definite local solutions to the above PDEs around x = 0

#### Then

- there exist smooth, positive definite local solutions to the above PDEs around x=0

$$\begin{aligned} \pi_c(x) &= \frac{1}{2} x' P_c^{-1} x &+ \pi_c^{[3]}(x) &+ \dots \\ \pi_o(x) &= \frac{1}{2} x' P_o x &+ \pi_o^{[3]}(x) &+ \dots \end{aligned}$$

#### Then

- there exist smooth, positive definite local solutions to the above PDEs around x=0

$$\begin{aligned} \pi_c(x) &= \frac{1}{2}x'P_c^{-1}x + \pi_c^{[3]}(x) + \dots \\ \pi_o(x) &= \frac{1}{2}x'P_ox + \pi_o^{[3]}(x) + \dots \end{aligned}$$

•  $P_c$ ,  $P_o$  are the controllability and observability gramians of the linear part of the system.

Scherpen showed that there is a local change of coordinates that brings the system into the form

$$\pi_c(x)=rac{1}{2}x'x, \qquad \pi_o(x)=rac{1}{2}x' \left[ egin{array}{cc} au_1(x) & 0 \ & \ddots & \ 0 & au_n(x) \end{array} 
ight]x$$

Scherpen showed that there is a local change of coordinates that brings the system into the form

$$\pi_c(x) = rac{1}{2}x'x, \qquad \pi_o(x) = rac{1}{2}x' \left[ egin{array}{cc} au_1(x) & 0 \ & \ddots & \ 0 & au_n(x) \end{array} 
ight] x$$

The  $\tau_i(x)$  are the squared singular value functions,

$$au_i(0) = au_i = \sigma_i^2$$

where  $\sigma_i$  are the Hankel singular values of the linear part of the system.

Scherpen showed that there is a local change of coordinates that brings the system into the form

$$\pi_c(x)=rac{1}{2}x'x, \qquad \pi_o(x)=rac{1}{2}x'\left[egin{array}{cc} au_1(x) & 0 \ & \ddots & \ 0 & au_n(x) \end{array}
ight]x$$

The  $\tau_i(x)$  are the squared singular value functions,

$$\tau_i(0) = \tau_i = \sigma_i^2$$

where  $\sigma_i$  are the Hankel singular values of the linear part of the system.

Unfortunately neither these coordinates nor the squared singular value functions are unique.

Scherpen showed that there is a local change of coordinates that brings the system into the form

$$\pi_c(x)=rac{1}{2}x'x, \qquad \pi_o(x)=rac{1}{2}x'\left[egin{array}{cc} au_1(x) & 0 \ & \ddots & \ 0 & au_n(x) \end{array}
ight]x$$

The  $\tau_i(x)$  are the squared singular value functions,

$$\tau_i(0) = \tau_i = \sigma_i^2$$

where  $\sigma_i$  are the Hankel singular values of the linear part of the system.

Unfortunately neither these coordinates nor the squared singular value functions are unique.

She obtained a reduced order model by Galerkin projection onto the states with large  $\tau_i(x)$ .

The functions  $\pi_c(x)$  and  $\pi_o(x)$  have power series expansions

$$\pi_c(x) = \frac{1}{2}x'P_c^{-1}x + \pi_c^{[3]}(x) + \pi_c^{[4]}(x) + \dots$$

$$\pi_o(x) = \frac{1}{2}x'P_ox + \pi_o^{[3]}(x) + \pi_o^{[4]}(x) + \dots$$

The functions  $\pi_c(x)$  and  $\pi_o(x)$  have power series expansions

$$\pi_c(x) = rac{1}{2}x'P_c^{-1}x + \pi_c^{[3]}(x) + \pi_c^{[4]}(x) + \dots$$

$$\pi_o(x) = \frac{1}{2}x'P_ox + \pi_o^{[3]}(x) + \pi_o^{[4]}(x) + \dots$$

Following Moore and Scherpen we can make a linear change of coordinates so that

$$P_c = \left[ egin{array}{cc} 1 & 0 \ & \ddots & \ 0 & 1 \end{array} 
ight], \qquad P_o = \left[ egin{array}{cc} au_1 & 0 \ & \ddots & \ 0 & au_n \end{array} 
ight]$$

where  $\tau_i = \sigma_i^2$  and  $\tau_1 \ge \tau_2 \ge \ldots \ge \tau_n > 0$ . After this linear change of coordinates the system is said to be in input normal form of degree one.

For simplicity of exposition we shall assume that the  $\tau_i$  are distinct,  $\tau_1 > \tau_2 > \ldots > \tau_n > 0$ .

For simplicity of exposition we shall assume that the  $\tau_i$  are distinct,  $\tau_1 > \tau_2 > \ldots > \tau_n > 0$ .

Now

$$egin{array}{rcl} \pi^{[3]}_c(x) &=& \displaystyle\sum_{i\leq j\leq k} \gamma^{ijk}_c x_i x_j x_k \ \pi^{[3]}_o(x) &=& \displaystyle\sum_{i\leq j\leq k} \gamma^{ijk}_o x_i x_j x_k \end{array}$$

For simplicity of exposition we shall assume that the  $\tau_i$  are distinct,  $\tau_1 > \tau_2 > \ldots > \tau_n > 0$ .

Now

$$egin{array}{rcl} \pi^{[3]}_c(x) &=& \displaystyle\sum_{i\leq j\leq k}\gamma^{ijk}_cx_ix_jx_k \ \pi^{[3]}_o(x) &=& \displaystyle\sum_{i\leq j\leq k}\gamma^{ijk}_ox_ix_jx_k \end{array}$$

Choose three indices  $1 \leq r \leq s \leq t \leq n$  , at least two indices

are different, r < t . Consider the change of coordinates

$$egin{array}{rcl} x_r &=& \xi_r + eta_r \xi_s \xi_t \ x_t &=& \xi_t + eta_t \xi_r \xi_s \ x_l &=& \xi_l & ext{otherwise} \end{array}$$

Then the quadratic parts of  $\pi_c$ ,  $\pi_o$  are left unchanged but the cubic parts each pick up an extra term,

$$\begin{aligned} \pi_c^{[3]}(\xi) &= \sum_{i \leq j \leq k} \gamma_c^{ijk} \xi_i \xi_j \xi_k + (\beta_r + \beta_t) \xi_r \xi_s \xi_t \\ \pi_o^{[3]}(\xi) &= \sum_{i \leq j \leq k} \gamma_o^{ijk} \xi_i \xi_j \xi_k + (\tau_r \beta_r + \tau_t \beta_t) \xi_r \xi_s \xi_t \end{aligned}$$

Then the quadratic parts of  $\pi_c$ ,  $\pi_o$  are left unchanged but the cubic parts each pick up an extra term,

$$\begin{aligned} \pi_c^{[3]}(\xi) &= \sum_{i \leq j \leq k} \gamma_c^{ijk} \xi_i \xi_j \xi_k + (\beta_r + \beta_t) \xi_r \xi_s \xi_t \\ \pi_o^{[3]}(\xi) &= \sum_{i \leq j \leq k} \gamma_o^{ijk} \xi_i \xi_j \xi_k + (\tau_r \beta_r + \tau_t \beta_t) \xi_r \xi_s \xi_t \end{aligned}$$

Since  $au_r > au_t$  we can solve the linear system

$$\left[\begin{array}{cc}1 & 1\\\tau_r & \tau_t\end{array}\right]\left[\begin{array}{c}\beta_r\\\beta_t\end{array}\right] = -\left[\begin{array}{c}\gamma_c^{rst}\\\gamma_o^{rst}\end{array}\right]$$

Then the quadratic parts of  $\pi_c$ ,  $\pi_o$  are left unchanged but the cubic parts each pick up an extra term,

$$\begin{aligned} \pi_c^{[3]}(\xi) &= \sum_{i \leq j \leq k} \gamma_c^{ijk} \xi_i \xi_j \xi_k + (\beta_r + \beta_t) \xi_r \xi_s \xi_t \\ \pi_o^{[3]}(\xi) &= \sum_{i < j \leq k} \gamma_o^{ijk} \xi_i \xi_j \xi_k + (\tau_r \beta_r + \tau_t \beta_t) \xi_r \xi_s \xi_t \end{aligned}$$

Since  $au_r > au_t$  we can solve the linear system

$$\left[ egin{array}{c} 1 & 1 \ au_r & au_t \end{array} 
ight] \left[ egin{array}{c} eta_r \ eta_t \end{array} 
ight] = - \left[ egin{array}{c} \gamma_c^{rst} \ \gamma_o^{rst} \end{array} 
ight]$$

This change of coordinates cancels the monomials  $\xi_r \xi_s \xi_t$  from  $\pi_c^{[3]}(\xi), \ \pi_o^{[3]}(\xi)$  .

But if r = s = t then we can cancel the monomial  $x_r^3$  from only one of  $\pi_c^{[3]}(\xi)$ ,  $\pi_o^{[3]}(\xi)$  by a change of coordinates of the form

$$egin{array}{rcl} x_r &=& \xi_r + eta_r \xi_r^2 \ x_l &=& \xi_l & ext{otherwise} \end{array}$$

But if r = s = t then we can cancel the monomial  $x_r^3$  from only one of  $\pi_c^{[3]}(\xi), \ \pi_o^{[3]}(\xi)$  by a change of coordinates of the form

$$egin{array}{rcl} x_r &=& \xi_r + eta_r \xi_r^2 \ x_l &=& \xi_l & ext{otherwise} \end{array}$$

We do so in  $\pi_c^{[3]}(\xi)$  to obtain input normal form of degree two,

$$egin{array}{rcl} \pi^{[3]}_{c}(\xi) &=& 0 \ \pi^{[3]}_{o}(\xi) &=& \sum_{i} \gamma^{iii}_{o} \xi_{i} \xi_{i} \xi_{i} \end{array}$$

But if r = s = t then we can cancel the monomial  $x_r^3$  from only one of  $\pi_c^{[3]}(\xi), \ \pi_o^{[3]}(\xi)$  by a change of coordinates of the form

$$egin{array}{rcl} x_r &=& \xi_r + eta_r \xi_r^2 \ x_l &=& \xi_l & ext{otherwise} \end{array}$$

We do so in  $\pi_c^{[3]}(\xi)$  to obtain input normal form of degree two,

$$egin{array}{rcl} \pi^{[3]}_{c}(\xi) &=& 0 \ \pi^{[3]}_{o}(\xi) &=& \sum_{i} \gamma^{iii}_{o} \xi_{i} \xi_{i} \xi_{i} \end{array}$$

If the three indices are distinct r < s < t then there are many ways to cancel  $x_r x_s x_t$  from  $\pi_c$ ,  $\pi_o$ .

We can do similarly for higher degrees and in this way bring the system into input normal form of degree d,

$$\begin{aligned} \pi_c(x) &= \frac{1}{2} \sum_{i=1}^n x_i^2 + O(x)^{d+2} \\ \pi_o(x) &= \frac{1}{2} \sum_{i=1}^n \tau_i^{[0:d-1]}(x_i) x_i^2 + O(x)^{d+2} \end{aligned}$$

where the squared singular value polynomials  $au_i^{[0:d-1]}(x_i)$  are of degrees 0 through d-1

$$au_i^{[0:d-1]}(x_i) \;\; = \;\; au_i + au_{i,1} x_i + \ldots + au_{i,d-1} x_i^{d-1}$$

We can do similarly for higher degrees and in this way bring the system into input normal form of degree d,

$$\begin{aligned} \pi_c(x) &= \frac{1}{2} \sum_{i=1}^n x_i^2 + O(x)^{d+2} \\ \pi_o(x) &= \frac{1}{2} \sum_{i=1}^n \tau_i^{[0:d-1]}(x_i) x_i^2 + O(x)^{d+2} \end{aligned}$$

where the squared singular value polynomials  $au_i^{[0:d-1]}(x_i)$  are of degrees 0 through d-1

$$au_i^{[0:d-1]}(x_i) \;\; = \;\; au_i + au_{i,1} x_i + \ldots + au_{i,d-1} x_i^{d-1}$$

We have "simultaneously diagonalized"  $\pi_c(x)$ ,  $\pi_o(x)$  through terms of degree  $\leq d + 1$ . There are no cross terms,  $x_i x_j \dots$ 

The squared singular value polynomial  $\tau_i^{[0:d-1]}(x_i)$  measures the relative importance of the state coordinate  $x_i$ .

The squared singular value polynomial  $\tau_i^{[0:d-1]}(x_i)$  measures the relative importance of the state coordinate  $x_i$ .

There may be several ways to bring the system to input normal form of degree d but if  $d \leq 6$  then the  $\tau_i^{[0:d-1]}(x_i)$  are unique.

The squared singular value polynomial  $\tau_i^{[0:d-1]}(x_i)$  measures the relative importance of the state coordinate  $x_i$ .

There may be several ways to bring the system to input normal form of degree d but if  $d \le 6$  then the  $\tau_i^{[0:d-1]}(x_i)$  are unique.

If the system is odd

$$egin{array}{rcl} f(-x,-u)&=&-f(x,u)\ h(-x)&=&-h(x) \end{array}$$

then  $\pi_c(x)$ ,  $\pi_o(x)$  are even and the  $\tau_i^{[0:d-1]}(x_i)$  are unique for  $d \leq 12$ .

As with linear balancd truncation we restrict to reduced order models that can be found by nonlinear Galerkin projection.

As with linear balancd truncation we restrict to reduced order models that can be found by nonlinear Galerkin projection.

For this we need an embedding  $\psi$  and a submersion  $\phi$ 

$$egin{aligned} \psi : I\!\!R^k & o & I\!\!R^n \ \psi : z & \mapsto & x = \psi(z) \ \phi : I\!\!R^n & o & I\!\!R^k \ \phi : x & \mapsto & z = \phi(x) \ \phi(\psi(z)) = z, & (\psi \circ \phi)^2(x) = (\psi \circ \phi)(x) \end{aligned}$$

As with linear balancd truncation we restrict to reduced order models that can be found by nonlinear Galerkin projection.

For this we need an embedding  $\psi$  and a submersion  $\phi$ 

$$\begin{split} \psi : I\!\!R^k &\to I\!\!R^n \\ \psi : z &\mapsto x = \psi(z) \\ \phi : I\!\!R^n &\to I\!\!R^k \\ \phi : x &\mapsto z = \phi(x) \\ \phi(\psi(z)) = z, \qquad (\psi \circ \phi)^2(x) = (\psi \circ \phi)(x) \end{split}$$

**Reduced Order Model** 

$$egin{array}{rcl} \dot{z}&=&rac{\partial\phi}{\partial x}(\psi(z))f(\psi(z),u)\ y&=&h(\psi(z)) \end{array}$$

We would like to choose the embedding  $\psi : \mathbb{R}^k \to \mathbb{R}^n$  so the k dimensional submanifold that is its range "maximizes" the output energy  $\pi_o(x)$  for given input energy  $\pi_c(x)$ .

We would like to choose the embedding  $\psi : \mathbb{R}^k \to \mathbb{R}^n$  so the k dimensional submanifold that is its range "maximizes" the output energy  $\pi_o(x)$  for given input energy  $\pi_c(x)$ .

If k > 1 then this prescription is not mathematically well-defined so we shall settle for a submanifold that approximates it.

We would like to choose the embedding  $\psi : \mathbb{R}^k \to \mathbb{R}^n$  so the k dimensional submanifold that is its range "maximizes" the output energy  $\pi_o(x)$  for given input energy  $\pi_c(x)$ .

If k > 1 then this prescription is not mathematically well-defined so we shall settle for a submanifold that approximates it.

Assume the system is in input normal form of degree d, that the range of input energy of interest is

$$\pi_c(x) ~pprox ~rac{1}{2}|x|^2 \leq rac{1}{2}c^2$$

and that

$$au_i^{[0:d-1]}(x_i) >> au_j^{[0:d-1]}(x_j) \ 1 \leq i \leq k < j \leq n ext{ and } |x_i|, \; |x_j| \leq c.$$

for

We would like to choose the embedding  $\psi : \mathbb{R}^k \to \mathbb{R}^n$  so the k dimensional submanifold that is its range "maximizes" the output energy  $\pi_o(x)$  for given input energy  $\pi_c(x)$ .

If k > 1 then this prescription is not mathematically well-defined so we shall settle for a submanifold that approximates it.

Assume the system is in input normal form of degree d, that the range of input energy of interest is

$$\pi_c(x) ~pprox ~rac{1}{2}|x|^2 \leq rac{1}{2}c^2$$

and that

$$\tau_i^{[0:d-1]}(x_i) >> \tau_j^{[0:d-1]}(x_j)$$

for  $1 \leq i \leq k < j \leq n$  and  $|x_i|, \; |x_j| \leq c.$ 

Then a "reasonable" choice of  $\psi(z) = x$  is

$$\psi(z_1,\ldots,z_k)= \ x \ =(z_1,\ldots,z_k,0,\ldots,0)$$

How should we choose the submersion  $\phi$  ?

How should we choose the submersion  $\phi$  ?

We choose  $\phi : \mathbb{R}^n \to \mathbb{R}^k$  to minimize the  $L^2$  norm of the difference of the outputs from x and  $\psi(\phi(x))$ .

How should we choose the submersion  $\phi$  ?

We choose  $\phi : \mathbb{R}^n \to \mathbb{R}^k$  to minimize the  $L^2$  norm of the difference of the outputs from x and  $\psi(\phi(x))$ .

Define the co-observability function as before

$$\pi_{oo}(x,ar{x}) = rac{1}{2}\int_0^\infty |y(t)-ar{y}(t)|^2 \; dt$$

where  $y(t), \ \bar{y}(t)$  are the outputs from  $x(0) = x, \ x(0) = \bar{x}$ . when u(t) = 0.

How should we choose the submersion  $\phi$  ?

We choose  $\phi : \mathbb{R}^n \to \mathbb{R}^k$  to minimize the  $L^2$  norm of the difference of the outputs from x and  $\psi(\phi(x))$ .

Define the co-observability function as before

$$\pi_{oo}(x,ar{x}) = rac{1}{2}\int_0^\infty |y(t)-ar{y}(t)|^2 \; dt$$

where  $y(t), \ \bar{y}(t)$  are the outputs from  $x(0) = x, \ x(0) = \bar{x}.$ when u(t) = 0.

Then  $\pi_{oo}$  satisfies the Lyapunov PDE

$$0 \;\; = \;\; rac{\partial \pi_{oo}}{\partial (x,ar{x})} (x,ar{x}) \left[ egin{array}{c} f(x,0) \ f(ar{x},0) \end{array} 
ight] + rac{1}{2} \, |h(x) - h(ar{x})|^2$$
As before  $\pi_{oo}$  has a power series expansion

$$\pi_{oo}(x,\bar{x}) = \frac{1}{2} \sum_{i} \tau_i (x_i - \bar{x}_i)^2 + \pi_{oo}^{[3]}(x,\bar{x}) + \dots$$

that can be computed term by term.

As before  $\pi_{oo}$  has a power series expansion

$$\pi_{oo}(x,\bar{x}) = \frac{1}{2} \sum_{i} \tau_i (x_i - \bar{x}_i)^2 + \pi_{oo}^{[3]}(x,\bar{x}) + \dots$$

that can be computed term by term.

We define  $\phi(x) = z$  as

$$\phi(x) = \operatorname{argmin}_{z} \pi_{oo}(x, \psi(z))$$

As before  $\pi_{oo}$  has a power series expansion

$$\pi_{oo}(x,\bar{x}) = \frac{1}{2} \sum_{i} \tau_i (x_i - \bar{x}_i)^2 + \pi_{oo}^{[3]}(x,\bar{x}) + \dots$$

that can be computed term by term.

We define  $\phi(x) = z$  as

$$\phi(x) = \operatorname{argmin}_{z} \pi_{oo}(x, \psi(z))$$

so  $\phi(x)$  satisfies

$$0 \;\; = \;\; {\partial \pi_{oo} \over \partial {ar x}} (x, \psi(\phi(x))) {\partial \psi \over \partial z}(\phi(x))$$

From the choice of  $\psi$  ,

$$rac{\partial \psi}{\partial z}(\phi(x)) \;\; = \; \left[ egin{array}{c} I \ 0 \end{array} 
ight]$$

From the choice of  $\psi$  ,

$$rac{\partial \psi}{\partial z}(\phi(x)) \;\; = \;\; \left[ egin{array}{c} I \ 0 \end{array} 
ight]$$

so  $\phi_i(x) = z_i$  satisfies for  $1 \leq i \leq k$ 

$$egin{array}{rl} \phi_i(x)&=&x_i\ &+rac{1}{ au_i}\left(rac{\partial\pi^{[3]}_{oo}}{\partialar x_i}(x,(\phi(x),0))+rac{\partial\pi^{[4]}_{oo}}{\partialar x_i}(x,(\phi(x),0))+\ldots
ight) \end{array}$$

From the choice of  $\psi$  ,

$$rac{\partial \psi}{\partial z}(\phi(x)) \;\; = \;\; \left[ egin{array}{c} I \ 0 \end{array} 
ight]$$

so  $\phi_i(x) = z_i$  satisfies for  $1 \leq i \leq k$ 

.

$$egin{array}{rl} \phi_i(x)&=&x_i\ &+rac{1}{ au_i}\left(rac{\partial\pi^{[3]}_{oo}}{\partialar x_i}(x,(\phi(x),0))+rac{\partial\pi^{[4]}_{oo}}{\partialar x_i}(x,(\phi(x),0))+\ldots
ight)$$

This can be solved term by term via repeated substitution.

$$\begin{array}{lll} \phi_i^{(1)}(x) &=& x_i \\ \phi_i^{(2)}(x) &=& x_i + \frac{1}{\tau_i} \frac{\partial \pi_{oo}^{[3]}}{\partial \bar{x}_i}(x,(\phi^{(1)}(x),0)) \end{array}$$

Full Order Model  $x \in I\!\!R^n$ 

$$\dot{x} = f(x,u)$$
  
 $y = h(x)$ 

Full Order Model  $x \in {I\!\!R}^n$ 

$$egin{array}{rcl} \dot{x}&=&f(x,u)\ y&=&h(x) \end{array}$$

Reduced Order Model  $z \in I\!\!R^k$ 

$$egin{array}{rcl} \dot{z}&=&a(z,u)&=&rac{\partial\phi}{\partial x}(\psi(z))f(\psi(z),u) \ y&=&c(z)&=&h(\psi(z)) \end{array}$$

Full Order Model  $x \in {I\!\!R}^n$ 

$$egin{array}{rcl} \dot{x}&=&f(x,u)\ y&=&h(x) \end{array}$$

Reduced Order Model  $z \in I\!\!R^k$ 

$$egin{array}{rcl} \dot{z}&=&a(z,u)&=&rac{\partial\phi}{\partial x}(\psi(z))f(\psi(z),u) \ y&=&c(z)&=&h(\psi(z)) \end{array}$$

What is the error between their input-output maps?

Full Order Model  $x \in I\!\!R^n$ 

$$egin{array}{rcl} \dot{x}&=&f(x,u)\ y&=&h(x) \end{array}$$

Reduced Order Model  $z \in I\!\!R^k$ 

$$egin{array}{rcl} \dot{z}&=&a(z,u)&=&rac{\partial\phi}{\partial x}(\psi(z))f(\psi(z),u) \ y&=&c(z)&=&h(\psi(z)) \end{array}$$

What is the error between their input-output maps? What is the error between their Hankel maps?

Full Order Model  $x \in I\!\!R^n$ 

$$egin{array}{rcl} \dot{x}&=&f(x,u)\ y&=&h(x) \end{array}$$

Reduced Order Model  $z \in I\!\!R^k$ 

$$egin{array}{rcl} \dot{z}&=&a(z,u)&=&rac{\partial\phi}{\partial x}(\psi(z))f(\psi(z),u) \ y&=&c(z)&=&h(\psi(z)) \end{array}$$

What is the error between their input-output maps? What is the error between their Hankel maps? What is the error between their Hankel maps restricted to optimal inputs?

**Full Order Optimal Feedback** 

$$egin{array}{rcl} u&=&\kappa(x)=\left(rac{\partial\pi_c}{\partial x}(x)rac{\partial f}{\partial u}(x,\kappa(x))
ight)'\ &=&G'x+\left(x'rac{\partial f^{[2]}}{\partial u}(x,G'x)
ight)'+\dots \end{array}$$

This can be solved term by term via repeated substitution.

**Full Order Optimal Feedback** 

$$egin{array}{rcl} u&=&\kappa(x)=\left(rac{\partial\pi_c}{\partial x}(x)rac{\partial f}{\partial u}(x,\kappa(x))
ight)'\ &=&G'x+\left(x'rac{\partial f^{[2]}}{\partial u}(x,G'x)
ight)'+\dots \end{array}$$

This can be solved term by term via repeated substitution. Combined closed loop system

$$\left[ egin{array}{c} \dot{x} \ \dot{z} \end{array} 
ight] \;\; = \;\; \left[ egin{array}{c} F+GG' & 0 \ BG' & A \end{array} 
ight] \left[ egin{array}{c} x \ z \end{array} 
ight] + \ldots$$

Full Order Optimal Feedback

$$egin{array}{rcl} u&=&\kappa(x)=\left(rac{\partial\pi_c}{\partial x}(x)rac{\partial f}{\partial u}(x,\kappa(x))
ight)'\ &=&G'x+\left(x'rac{\partial f^{[2]}}{\partial u}(x,G'x)
ight)'+\dots \end{array}$$

This can be solved term by term via repeated substitution. Combined closed loop system

$$\left[ egin{array}{c} \dot{x} \ \dot{z} \end{array} 
ight] \;\; = \;\; \left[ egin{array}{cc} F+GG' & 0 \ BG' & A \end{array} 
ight] \left[ egin{array}{c} x \ z \end{array} 
ight] + \ldots$$

F + GG' is antistable. A is stable.

#### So there is an unstable manifold $z = \theta(x)$ satisfying

$$a( heta(x),\kappa(x)) \;\;=\;\; rac{\partial heta}{\partial x}(x)f(x,\kappa(x))$$

So there is an unstable manifold  $z = \theta(x)$  satisfying

$$a( heta(x),\kappa(x)) \;\;=\;\; rac{\partial heta}{\partial x}(x)f(x,\kappa(x))$$

This PDE can be solved term by term.

#### Define the cross-observability function

$$ho(x^0,z^0) \;\;=\;\; rac{1}{2}\int_0^\infty |y(t)-w(t)|^2 dt$$

where

$$egin{array}{rcl} \dot{x} &=& f(x,0) & \dot{z} &=& a(z,0) \ y &=& h(x) & w &=& c(z) \ x(0) &=& x^0 & z(0) &=& z^0 \end{array}$$

#### Define the cross-observability function

$$ho(x^0,z^0) \;\;=\;\; rac{1}{2}\int_0^\infty |y(t)-w(t)|^2 dt$$

where

$$egin{array}{rcl} \dot{x}&=&f(x,0)&&\dot{z}&=&a(z,0)\ y&=&h(x)&&w&=&c(z)\ x(0)&=&x^0&&z(0)&=&z^0 \end{array}$$

Then  $\rho$  satisfies the Lyapunov PDE

$$0 \hspace{.1in} = \hspace{.1in} rac{\partial 
ho}{\partial (x,z)} (x,z) \left[ egin{array}{c} f(x,0) \ a(z,0) \end{array} 
ight] + rac{1}{2} \left| h(x) - c(z) 
ight|^2$$

which can be solved term by term.

Let  $u_x(-\infty:0)$  be the optimal control that excites the full order system to x(0)=x . Then the nonlinear Hankel maps satisfy

$$|\mathcal{H}_n(u_x(-\infty:0))-\mathcal{H}_k(u_x(-\infty:0))|^2\leq
ho(x, heta(x))$$

Let  $u_x(-\infty:0)$  be the optimal control that excites the full order system to x(0)=x . Then the nonlinear Hankel maps satisfy

$$|\mathcal{H}_n(u_x(-\infty:0))-\mathcal{H}_k(u_x(-\infty:0))|^2\leq
ho(x, heta(x))$$

A good estimate of the error between the nonlinear Hankel maps is

$$\sup\left(rac{
ho(x, heta(x))}{\pi_c(x)}
ight)^{1/2}$$

As with input normal form we can make a change of coordinates so that

$$egin{array}{rcl} \pi_c(x) &=& rac{1}{2} |x|^2 + O(x)^{d+2} \ &
ho(x, heta(x)) &=& rac{1}{2} \sum \epsilon_i^{[0:d-1]}(x_i) x_i^2 + O(x)^{d+2} \end{array}$$

As with input normal form we can make a change of coordinates so that

$$egin{array}{rll} \pi_c(x)&=&rac{1}{2}|x|^2+O(x)^{d+2} \ &
ho(x, heta(x))&=&rac{1}{2}\sum\epsilon_i^{[0:d-1]}(x_i)x_i^2+O(x)^{d+2} \end{array}$$

 $\epsilon_i^{[0:d-1]}(x_i)$  are the error polynomials of degrees 0 through d-1.

As with input normal form we can make a change of coordinates so that

$$egin{array}{rcl} \pi_c(x) &=& rac{1}{2} |x|^2 + O(x)^{d+2} \ &
ho(x, heta(x)) &=& rac{1}{2} \sum \epsilon_i^{[0:d-1]}(x_i) x_i^2 + O(x)^{d+2} \end{array}$$

 $\epsilon_i^{[0:d-1]}(x_i)$  are the error polynomials of degrees 0 through d-1.

The error polynomials are unique for  $d \leq 6$ .

Three linked rods connected by planar rotary joints with springs and dampening hanging from the ceiling. The input is a torque applied to the top joint and the output is the horizontal displacement of the bottom. Each rod is uniform of length l = 2, mass  $\mu = 1$ , with spring constant c = 3, dampening constant b = 0.5 and gravity constant g = 0.5.



We approximated the nonlinear system by its Taylor series through terms of degree 5.

We approximated the nonlinear system by its Taylor series through terms of degree 5.

The Taylor series of controllability and observability functions  $\pi_c(x), \ \pi_o(x)$  were computed through terms of degree 6.

We approximated the nonlinear system by its Taylor series through terms of degree 5.

The Taylor series of controllability and observability functions  $\pi_c(x), \ \pi_o(x)$  were computed through terms of degree 6.

The system was brought into input normal form of degree 5 by a changes of state coordinates of degrees 1 through 5.

We approximated the nonlinear system by its Taylor series through terms of degree 5.

The Taylor series of controllability and observability functions  $\pi_c(x), \ \pi_o(x)$  were computed through terms of degree 6.

The system was brought into input normal form of degree 5 by a changes of state coordinates of degrees 1 through 5.

The Hankel singular values of the linear part of the system are

 $\begin{bmatrix} 15.3437 & 14.9678 & 0.3102 & 0.2470 & 0.0156 & 0.0091 \end{bmatrix}$ 

We approximated the nonlinear system by its Taylor series through terms of degree 5.

The Taylor series of controllability and observability functions  $\pi_c(x), \ \pi_o(x)$  were computed through terms of degree 6.

The system was brought into input normal form of degree 5 by a changes of state coordinates of degrees 1 through 5.

The Hankel singular values of the linear part of the system are

 $\begin{bmatrix} 15.3437 & 14.9678 & 0.3102 & 0.2470 & 0.0156 & 0.0091 \end{bmatrix}$ 

Apparently only two dimensions are linearly significant.

Here are the squared singular value polynomials.

$$au_1^{[0:4]}(x_1) ~=~ 235.4298 - 3.4163 x_1^2 - 0.3104 x_1^4$$

$$au_2^{[0:4]}(x_2) ~=~ 224.0356 - 3.2750 x_2^2 - 0.2941 x_2^4$$

$$au_3^{[0:4]}(x_3) ~=~ 000.0962 + 0.0014 x_3^2 - 0.0001 x_3^4$$

$$au_4^{[0:4]}(x_4) ~=~ 000.0610 + 0.0006 x_4^2 + 0.0000 x_4^4$$

$$au_5^{[0:4]}(x_5) ~=~ 000.0002 + 0.0000 x_5^2 + 0.0000 x_5^4$$

$$au_6^{[0:4]}(x_6) = 000.0001 + 0.0000x_6^2 + 0.0000x_6^4$$

Here are the squared singular value polynomials.

$$au_1^{[0:4]}(x_1) ~=~ 235.4298 - 3.4163 x_1^2 - 0.3104 x_1^4$$

$$au_2^{[0:4]}(x_2) ~=~ 224.0356 - 3.2750 x_2^2 - 0.2941 x_2^4$$

$$au_3^{[0:4]}(x_3) = 000.0962 + 0.0014 x_3^2 - 0.0001 x_3^4$$

$$au_4^{[0:4]}(x_4) ~=~ 000.0610 + 0.0006 x_4^2 + 0.0000 x_4^4$$

$$au_5^{[0:4]}(x_5) ~=~ 000.0002 + 0.0000 x_5^2 + 0.0000 x_5^4$$

$$au_6^{[0:4]}(x_6) ~=~ 000.0001 + 0.0000 x_6^2 + 0.0000 x_6^4$$

#### Apparently only two dimensions are nonlinearly significant.

# Semilog plot of the squared singular value polynomials $au_i^{[0:4]}$



# Semilog plot of the squared singular value polynomials $au_i^{[0:4]}$



Notice the difference in scale and how flat they are.

Let  $u_x(-\infty:0)$  be the optimal input that excites the full system to x.

Let  $u_x(-\infty:0)$  be the optimal input that excites the full system to x.

The error between the nonlinear Hankel maps satisfies

$$egin{aligned} |\mathcal{H}_n(u_x(-\infty:0)) - \mathcal{H}_k(u_x(-\infty:0))|^2 \ &\leq 0.0965 |x|^2 - 0.0009 |x|^4 + \dots \end{aligned}$$

Let  $u_x(-\infty:0)$  be the optimal input that excites the full system to x.

The error between the nonlinear Hankel maps satisfies

$$egin{aligned} |\mathcal{H}_n(u_x(-\infty:0)) - \mathcal{H}_k(u_x(-\infty:0))|^2 \ &\leq 0.0965 |x|^2 - 0.0009 |x|^4 + \dots \end{aligned}$$

By way of comparison, the square of the third Hankel singular value is

0.0962

so this estimate is tight.
#### Nonlinear Example

# Here are the error polynomials $\epsilon_i^{[0:2]}$



#### Nonlinear Example

# Here are outputs of the Hankel maps of the full and reduced systems excited by an optimal control $u_x(-\infty:0)$ for random x.



#### Nonlinear Example

Here are the responses of the full nonlinear model (blue), the reduced nonlinear model (green) and the linear part of the full model (red) to a sinusoidal input.



• We gave a new interpretation of linear balanced truncation that readily generalizes to nonlinear systems.

- We gave a new interpretation of linear balanced truncation that readily generalizes to nonlinear systems.
- The nonlinear method is computionally feasible for moderate state dimensions and moderate degrees of approximation.

- We gave a new interpretation of linear balanced truncation that readily generalizes to nonlinear systems.
- The nonlinear method is computionally feasible for moderate state dimensions and moderate degrees of approximation.
- For both linear and nonlinear model reduction, we gave new computable error bounds for the difference of the full and reduced Hankel maps restricted to optimal inputs.

- We gave a new interpretation of linear balanced truncation that readily generalizes to nonlinear systems.
- The nonlinear method is computionally feasible for moderate state dimensions and moderate degrees of approximation.
- For both linear and nonlinear model reduction, we gave new computable error bounds for the difference of the full and reduced Hankel maps restricted to optimal inputs.
- We believe that these bounds are typical of the errors for inputs that are used in practice.

- We gave a new interpretation of linear balanced truncation that readily generalizes to nonlinear systems.
- The nonlinear method is computionally feasible for moderate state dimensions and moderate degrees of approximation.
- For both linear and nonlinear model reduction, we gave new computable error bounds for the difference of the full and reduced Hankel maps restricted to optimal inputs.
- We believe that these bounds are typical of the errors for inputs that are used in practice.
- The methodology extends directly to other methods of balancing such as  $LQG, \ H^{\infty}$ .

- We gave a new interpretation of linear balanced truncation that readily generalizes to nonlinear systems.
- The nonlinear method is computionally feasible for moderate state dimensions and moderate degrees of approximation.
- For both linear and nonlinear model reduction, we gave new computable error bounds for the difference of the full and reduced Hankel maps restricted to optimal inputs.
- We believe that these bounds are typical of the errors for inputs that are used in practice.
- The methodology extends directly to other methods of balancing such as  $LQG,\ H^\infty$  .
- Thank you for listening.