

# Learning Koopman eigenfunctions for prediction and control

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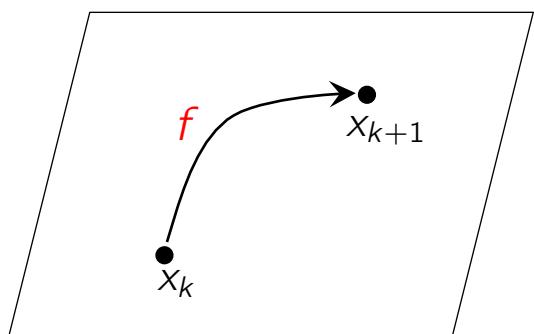
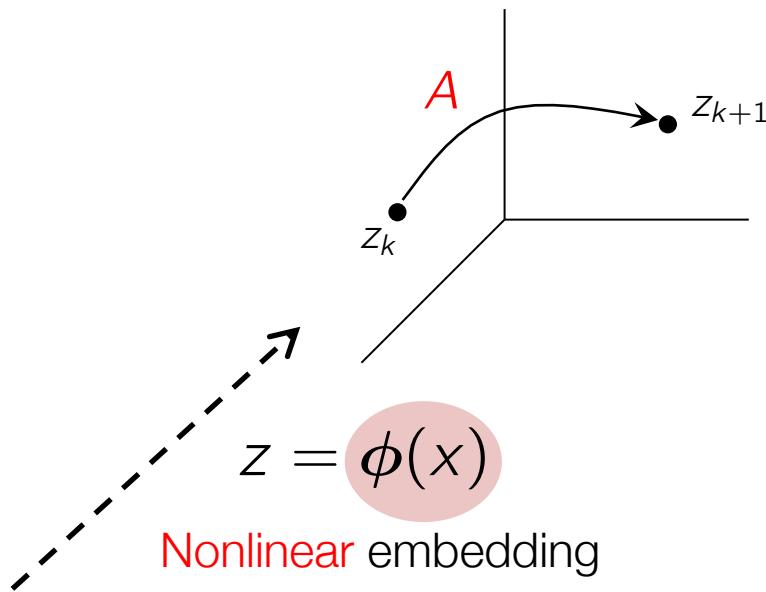
Milan Korda  
(LAAS-CNRS)

Igor Mezić  
(University of California, Santa Barbara)

# Linear prediction

Linear dynamics

$$z_{k+1} = A z_k$$

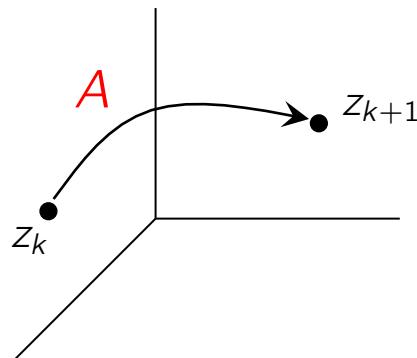


Nonlinear

# Linear prediction

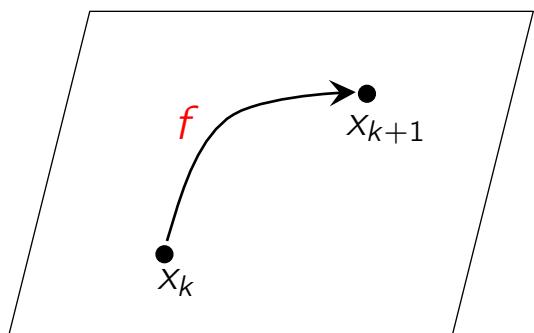
Linear dynamics

$$z_{k+1} = A z_k$$



$$z = \phi(x)$$

Nonlinear embedding



Nonlinear

Linear projection

$$\xi(x_k) \approx C z_k$$

$\xi$  = vector of observables

(e.g.  $\xi(x) = x$ )

# Why linear predictors?

$$z_{k+1} = \mathbf{A} z_k$$

$$z_0 = \phi(x_0)$$

$$\hat{y}_k = \mathbf{C} z_k$$

$$\hat{y}_k \approx \xi(x_k)$$

# Why linear predictors?

$$\begin{aligned}z_{k+1} &= \mathbf{A}z_k \\z_0 &= \phi(x_0) \\\hat{y}_k &= \mathbf{C}z_k\end{aligned}$$

$$\hat{y}_k \approx \xi(x_k)$$

Nonlinear feedback control & estimation using linear techniques

⇒ Model predictive control [Korda & Mezic, 2018]

⇒ State estimation [Surana & Banaszuk, 2016]

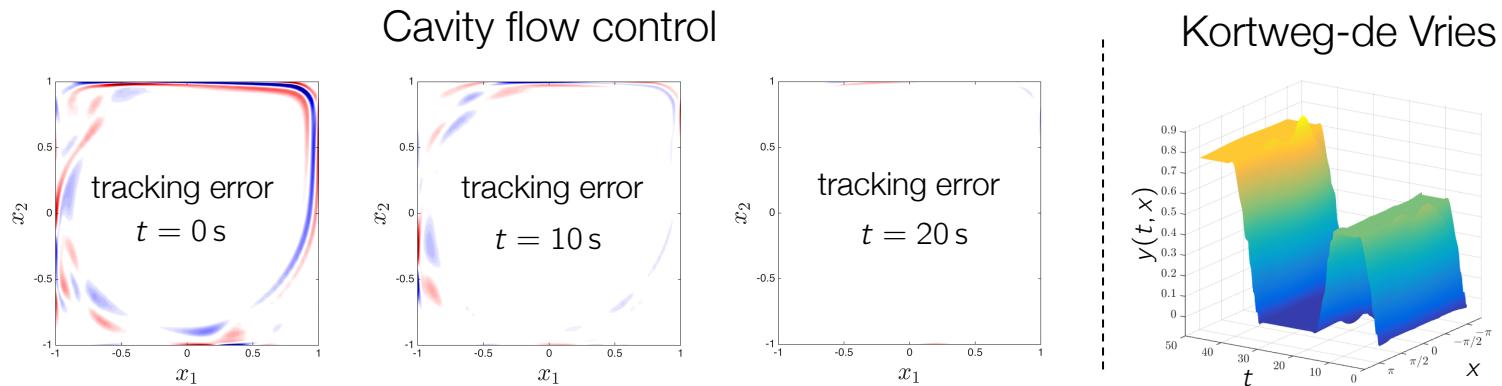
Mature & well understood

Fast computation (linear algebra / convex optimization)

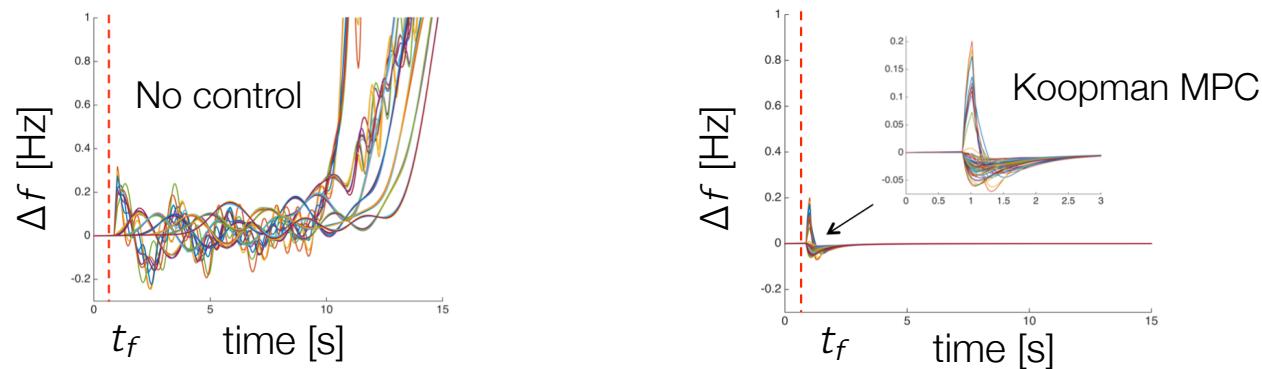
Rapid deployment in applications

# Koopman MPC - applications

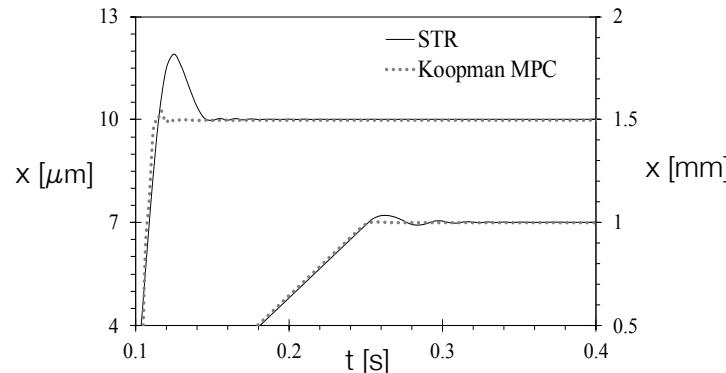
Fluid dynamics  
[Arbabi et al. 2018]



Powergrid  
[Korda et al. 2017]



High-precision positioning  
[Kamenar et al. 2018]



# Choosing the embedding

$$\begin{aligned} z_{k+1} &= \textcolor{red}{A}z_k \\ z_0 &= \textcolor{red}{\phi}(x_0) \\ \hat{y}_k &= \textcolor{red}{C}z_k \end{aligned}$$

When can we predict exactly?

$$\hat{y}_k = \textcolor{red}{\xi}(x_k)$$

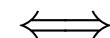
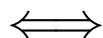
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$$\hat{y}_k = \xi(x_k)$$

if

$\text{span}\{\phi_1, \dots, \phi_N\}$  is Koopman invariant &  $\xi \in \text{span}\{\phi_1, \dots, \phi_N\}$



$\phi_i$ 's are (generalized) Koopman eigenfunctions

(or linear combinations thereof)

Span of  $\phi_i$ 's is rich enough

# Eigenfunction construction

# Eigenfunction construction

$$\dot{x} = f(x)$$

Eigenfunction

$$\phi(S_t(x)) = e^{\lambda t} \phi(x)$$

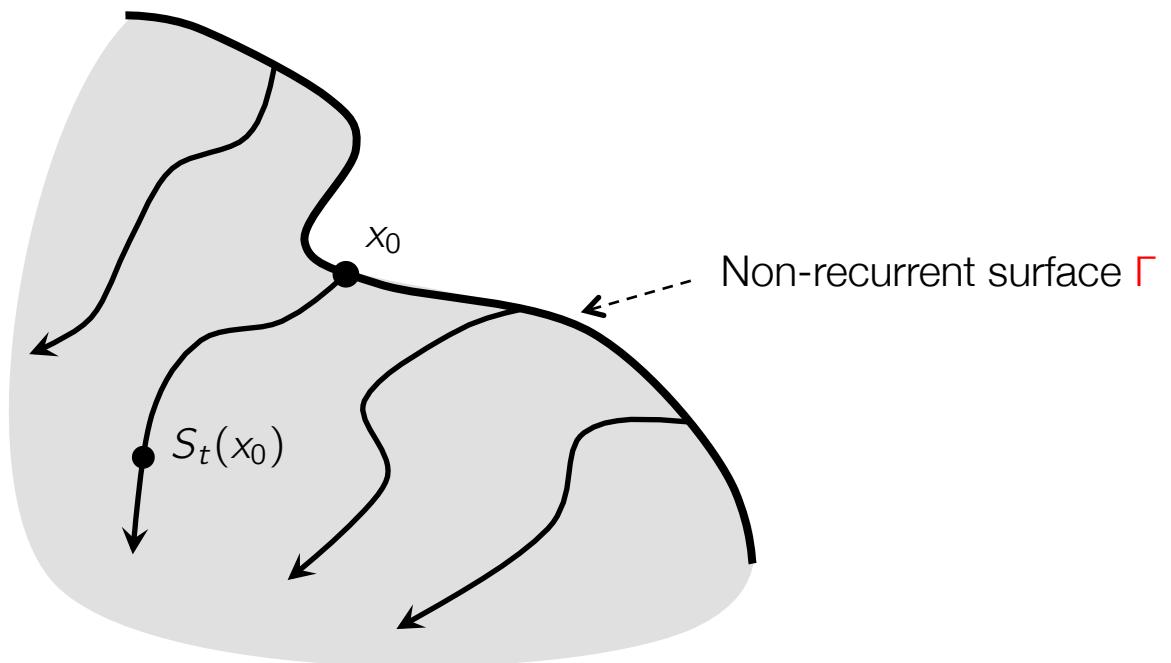
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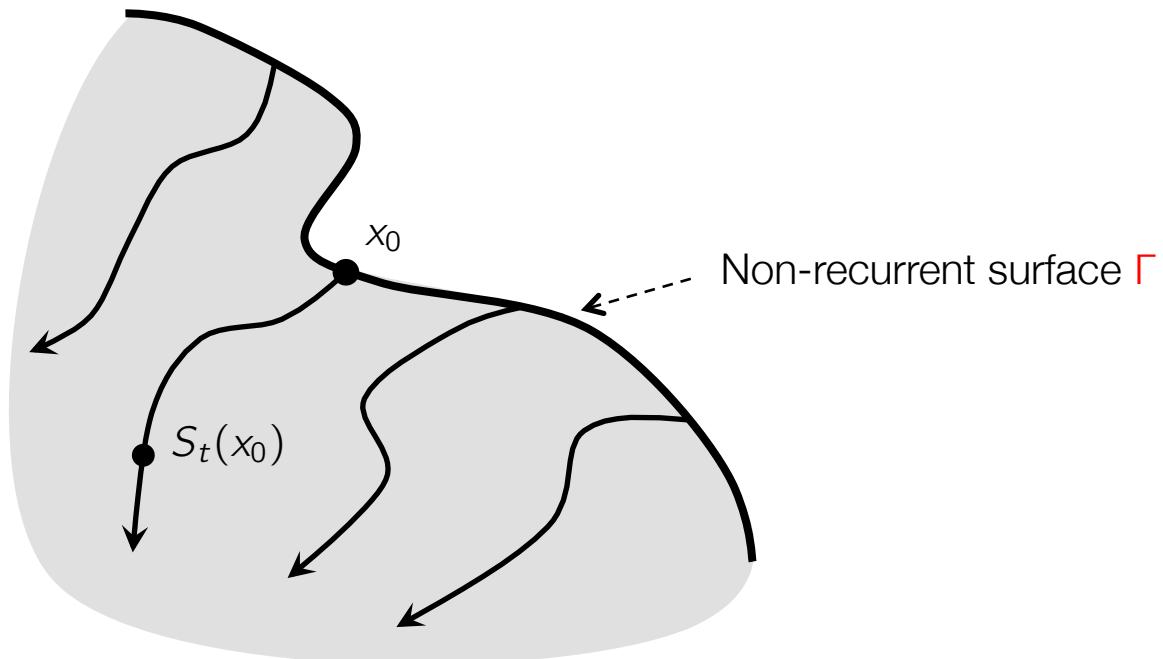
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$\lambda = \text{arbitrary}$  complex number

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$$\phi_{\lambda,g}(S_t(x_0)) = e^{\lambda t} g(x_0) \quad x_0 \in \Gamma$$

$$\phi_{\lambda,g} = g \quad \text{on } \Gamma$$



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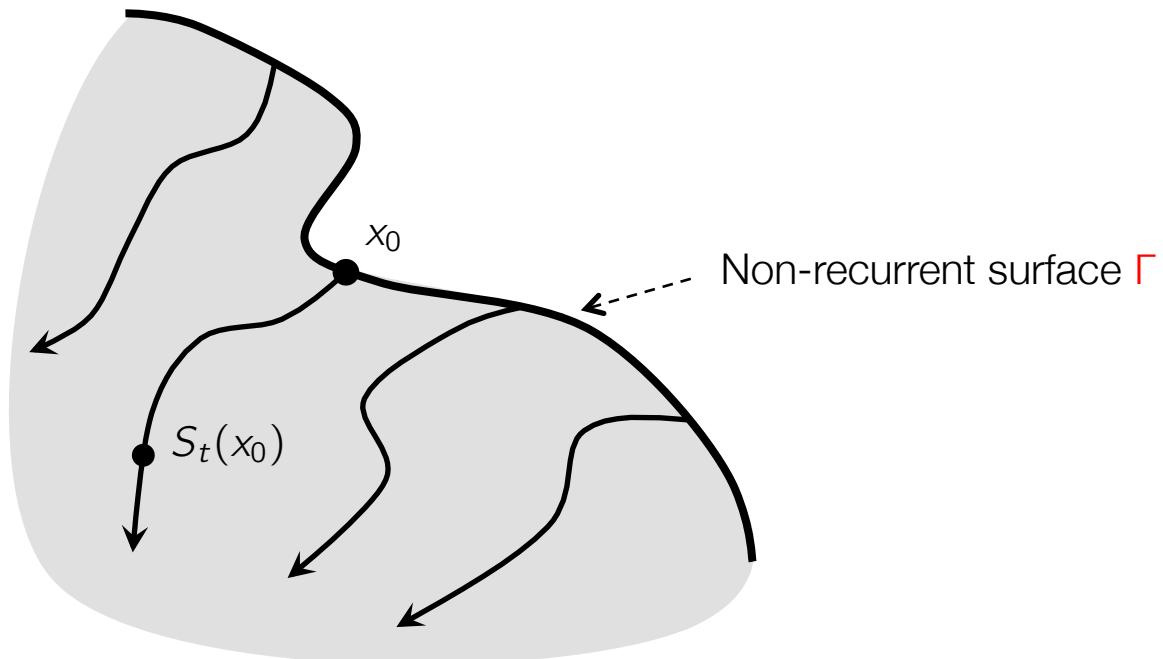
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**Lemma:**  $\Gamma$  non-recurrent &  $g$  continuous  $\Rightarrow \phi_{\lambda,g}$  is a continuous eigenfunction

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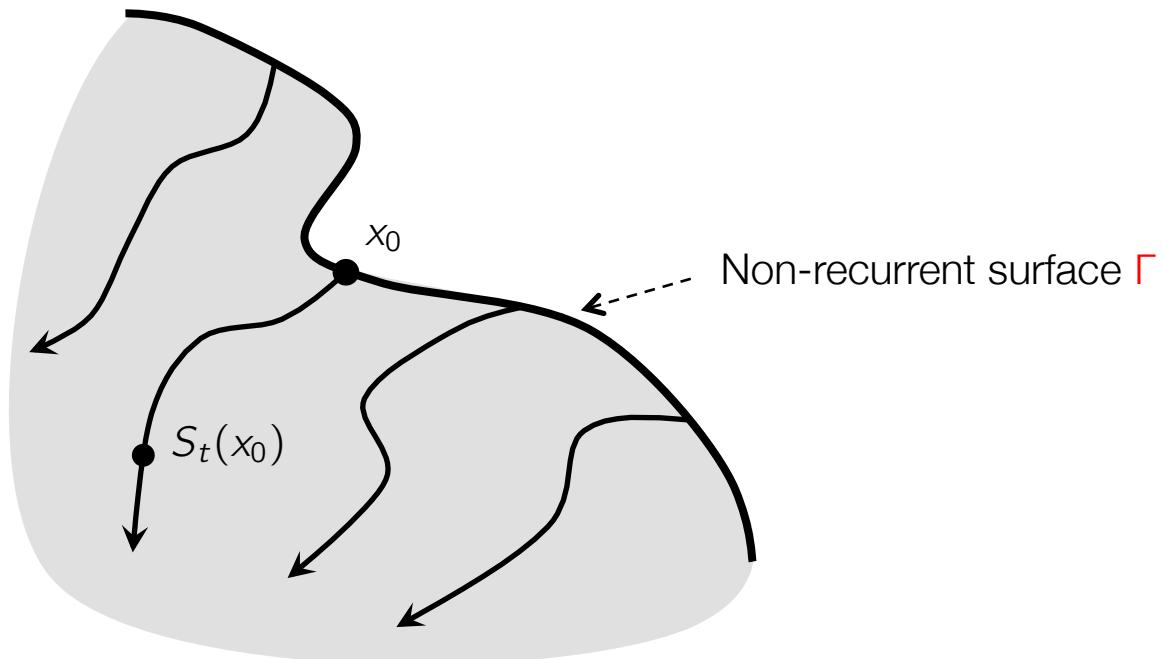
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cf. Open eigenfunctions [Mezic 2017]

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$G = \{g_i\}_{i=1}^{\infty}$  with  $\text{span}\{G\}$  dense in  $\mathcal{C}(\Gamma)$

**Theorem:**  $\Gamma$  non-recurrent,  $\Lambda_0 = \bar{\Lambda}_0$  &  $\exists \lambda \in \Lambda_0$  with  $\text{Re}(\lambda) \neq 0$

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For every continuous function  $\xi$  and every  $\epsilon > 0$  there exists  $\phi_1, \dots, \phi_N \in \Phi_{\Lambda, G}$  such that

$$\sup_x \left| \xi(x) - \sum_{i=1}^N c_i \phi_i(x) \right| < \epsilon$$

for some coefficients  $c_1, \dots, c_N$

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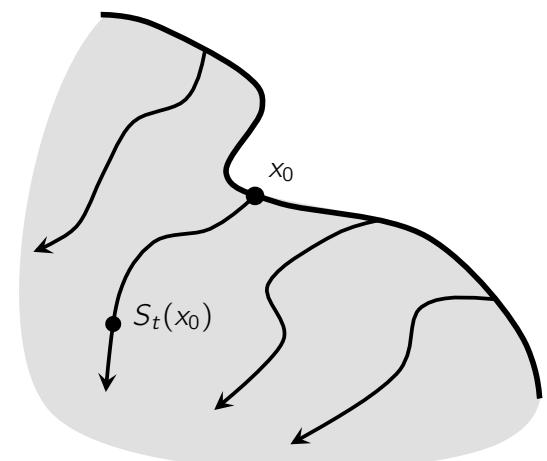
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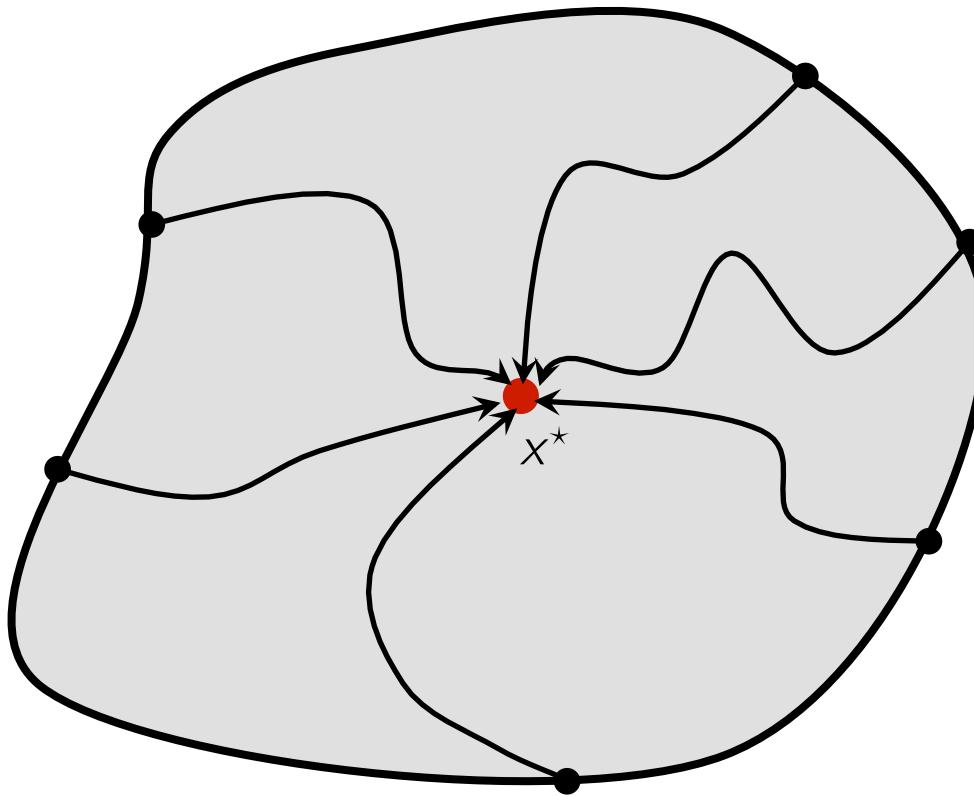
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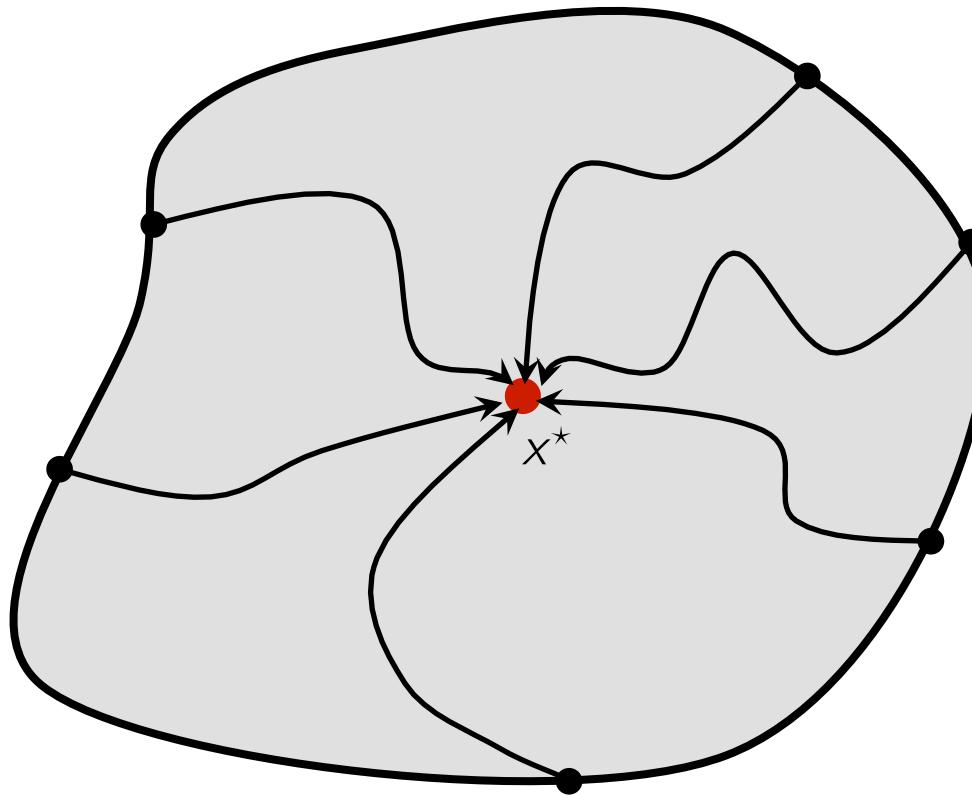
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Singularities don't matter in  $\mathcal{C}$



But they do in higher regularity spaces



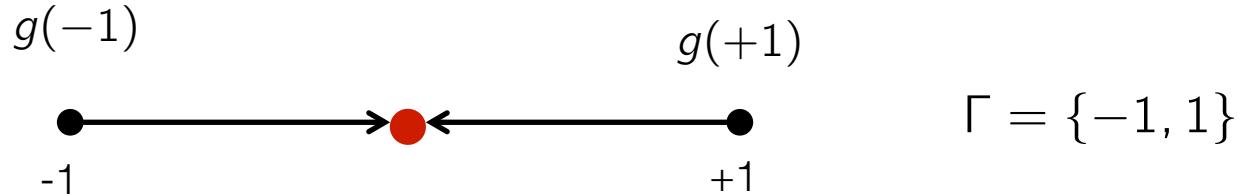
[Mezić 2017]

Stable equilibrium  $x^*$  + conditions on eigenvalues of  $Df(x^*) \Rightarrow$  real **analytic** eigenfunctions **dense in  $\mathcal{C}$**  exist

M. Kvalheim's poster: regular eigenfunctions  $\Rightarrow$  principal eigenvalues

# Example

$$\dot{x} = ax, \quad a < 0$$



$$\phi_{\lambda, g}(x) = \begin{cases} g(-1)|x|^{\frac{\lambda}{a}} & x < 0 \\ g(+1)|x|^{\frac{\lambda}{a}} & x > 0 \end{cases}$$

## Observations:

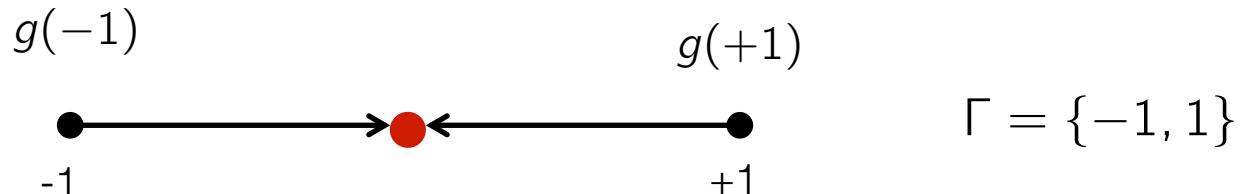
Continuous on  $[-1, 1]$  if  $\lambda < 0$  (and on  $[-1, 1] \setminus \{0\}$  for any  $\lambda$ )

Analytic if  $\lambda = k \cdot a$ ,  $k \in \mathbb{N}$  and  $g(-1) = (-1)^k g(+1)$

$\text{span}\{|x|^{\frac{k\lambda}{a}} : k \in \mathbb{N}\}$  dense in  $\mathcal{C}([0, 1])$  if  $\lambda < 0$

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**Optimal** choice of  $\lambda$  and  $g$  depends on  $\xi$

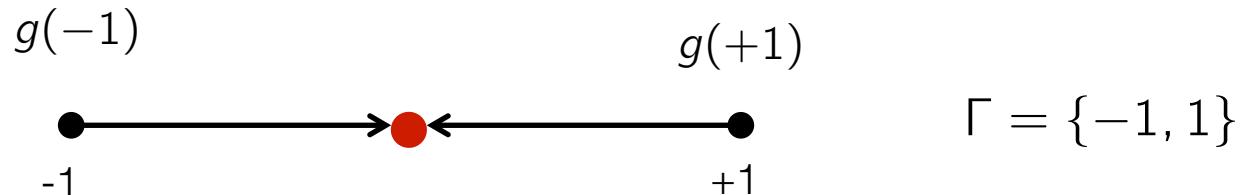
$$\xi = x \quad \Rightarrow \quad \lambda = a \quad \text{and} \quad g(+1) = 1, g(-1) = -1$$

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$$\xi = \max(0, x) \quad \Rightarrow \quad \lambda = a \quad \text{and} \quad g(+1) = 1, g(-1) = 0$$

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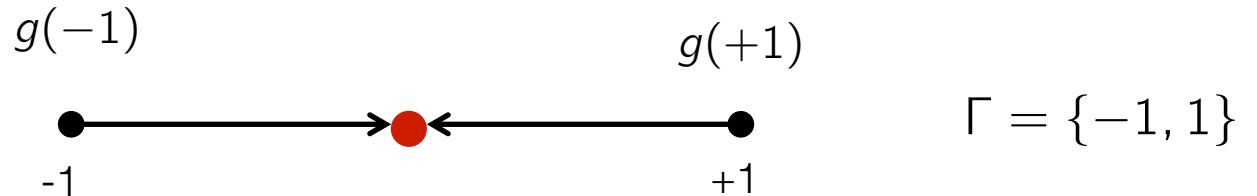
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But  $\text{span}\{\phi_{k\lambda,g} : k \in \mathbb{N}, g \in \mathcal{C}(\Gamma)\}$  dense in  $\mathcal{C}([-1, 1])$  for any  $\lambda < 0$

# Data-driven construction

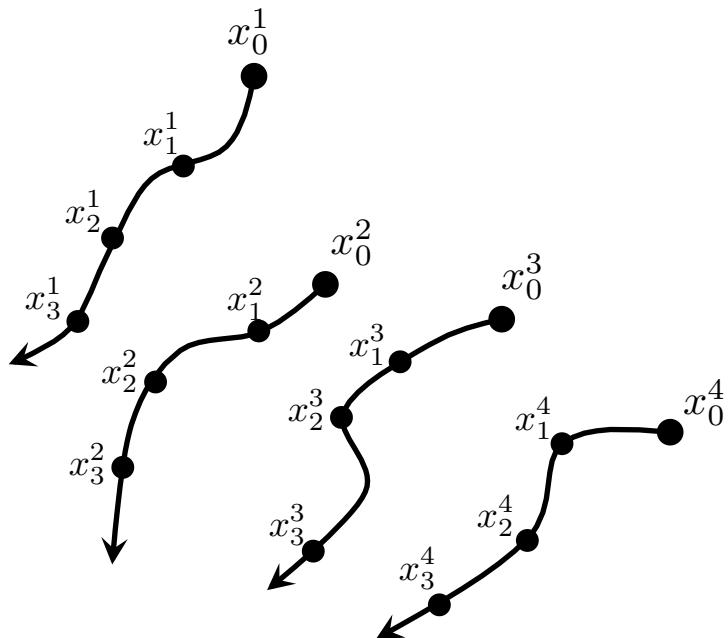
# Data-driven construction

$g = \text{arbitrary}$  continuous function

$\lambda = \text{arbitrary}$  complex number

eigenfunction  $\phi_{\lambda,g}$  defined on data

$$\phi_{\lambda,g}(x_k^j) := e^{\lambda k T_s} g(x_0^j)$$



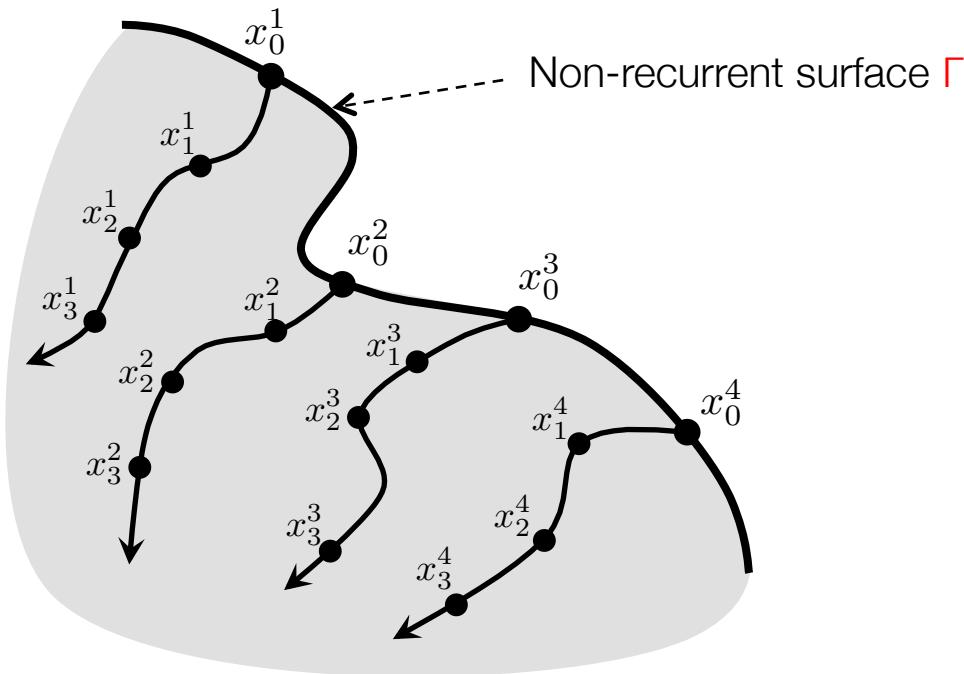
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**Lemma:** Flow rectifiable & initial conditions on distinct trajectories

$\Rightarrow \exists$  non-recurrent surface  $\Gamma$  passing through initial conditions

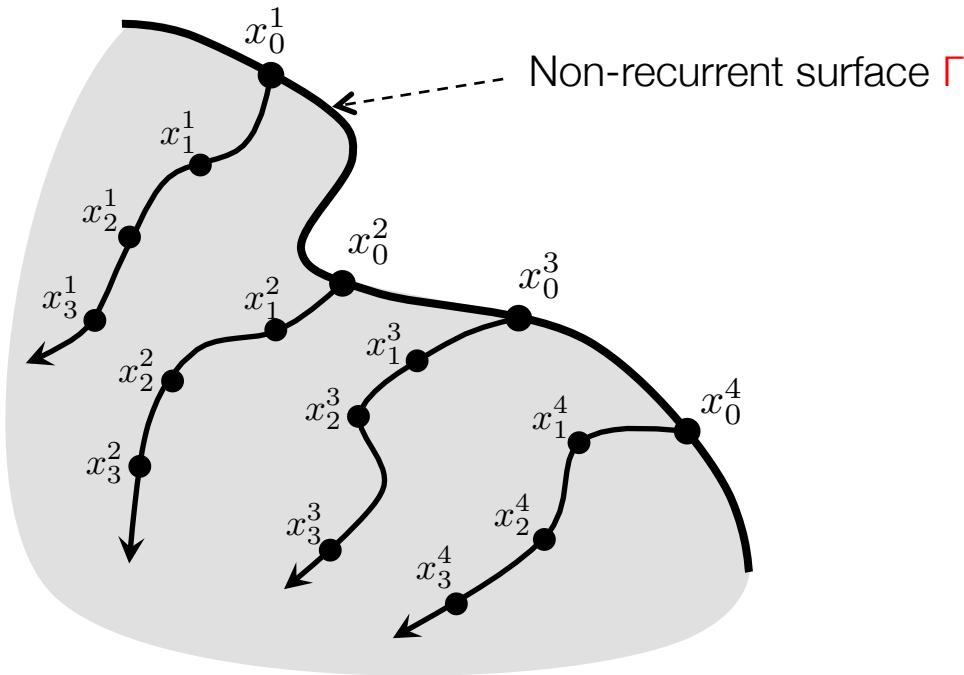
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**Lemma:** Flow rectifiable & initial conditions on distinct trajectories

$\Rightarrow \exists$  non-recurrent surface  $\Gamma$  passing through initial conditions

$\Rightarrow \{\phi_{\lambda,g}(x_k^j)\}_{j,k}$  samples of a **continuous** eigenfunction  $\Rightarrow$  can **interpolate**

# Algorithm summary

Eigenfunction construction

**Given** trajectory data  $(x_k^j)_{j,k}$

**Choose**  $\lambda_1, \dots, \lambda_{N_\lambda}$  complex numbers

**Choose**  $g_1, \dots, g_{N_g}$  continuous functions

**Construct**  $N := N_\lambda N_g$  eigenfunctions by

**Set**  $\phi_{\lambda,g}(x_k^j) := e^{\lambda k T_s} g(x_0^j)$  for each  $\lambda$  and  $g$

**Interpolate**  $\phi_{\lambda,g}(x_k^j)$  to get  $\hat{\phi}_{\lambda,g}$

**Output**  $\hat{\phi} = [\hat{\phi}_1, \dots, \hat{\phi}_N]$

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## Predictor matrices

**Set**  $A = \text{diag}(\lambda_1, \dots, \lambda_N)$

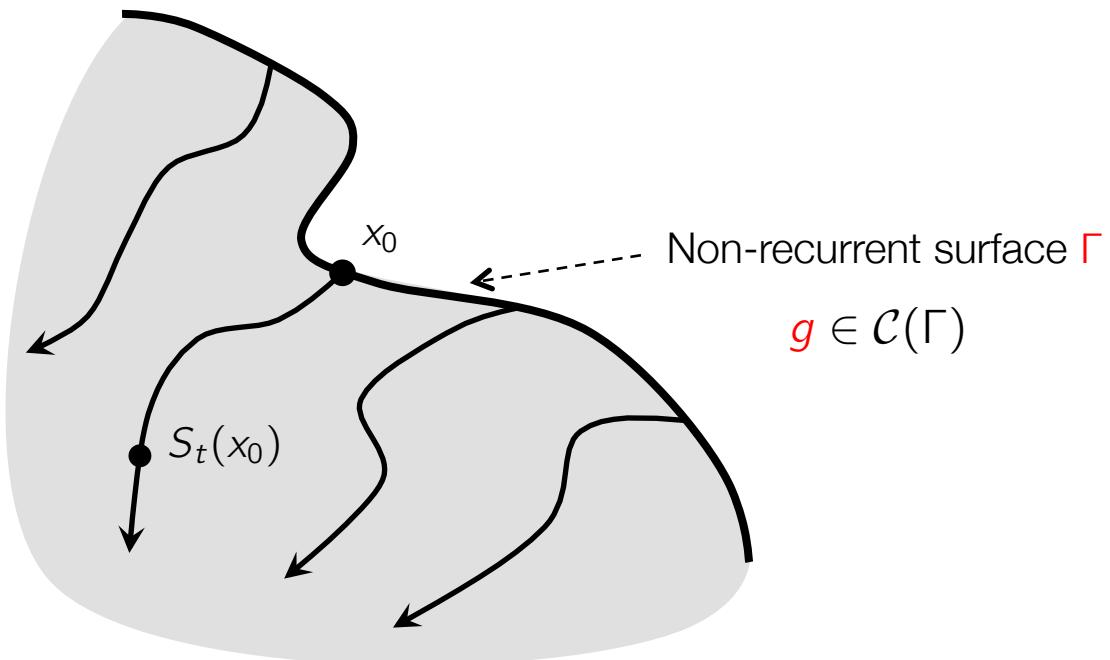
**Get**  $C$  by minimizing  $\sum_{i=1}^M \|\xi(\bar{x}_i) - C\hat{\phi}(\bar{x}_i)\|^2$   
(Linear least-squares)

$$\begin{aligned} z_{k+1} &= \textcolor{red}{A}z_k \\ z_0 &= \hat{\phi}(x_0) \\ \hat{y}_k &= \textcolor{red}{C}z_k \end{aligned}$$

# Optimal choice of boundary functions

# Choice of boundary functions

$$\phi_{\lambda,g}(S_t(x_0)) = e^{\lambda t} g(x_0)$$



**Observation:**  $\phi_{\lambda,g}$  depends linearly on  $g$   $\Rightarrow$  maybe can choose  $g$  using convex optimization

# Choice of boundary functions

$$\phi_{\lambda, \textcolor{red}{g}}(S_t(x_0)) = e^{\lambda t} \textcolor{red}{g}(x_0)$$

Given  $\lambda_1, \dots, \lambda_{N_\lambda}$  there exist **linear** operators  $\mathcal{L}_{\lambda_i}$  such that

$$\mathcal{L}_{\lambda_i} g = \phi_{\lambda_i, g}$$

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$$\mathcal{L}_{\lambda_i} g = \phi_{\lambda, g}$$

Given  $\xi$  we want to find  $\textcolor{red}{g}_1, \dots, \textcolor{red}{g}_{N_\lambda}$  such that

$$\|\xi - \text{Proj}_{\text{span}\{\mathcal{L}_{\lambda_1} \textcolor{red}{g}_1, \dots, \mathcal{L}_{\lambda_{N_\lambda}} \textcolor{red}{g}_{N_\lambda}\}} \xi\| \text{ is minimized}$$

Is this convex in  $\textcolor{red}{g}$ 's?

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Is this convex in  $\textcolor{red}{g}$ 's?

**Answer:** probably not, but a **convex reformulation** exists if each component of  $\xi$  is considered separately

# Convex reformulation

$$\underset{\substack{\boldsymbol{g}_i \in \mathcal{C}(\Gamma)}}{\text{minimize}} \|\boldsymbol{\xi} - \text{Proj}_{\text{span}\{\mathcal{L}_{\lambda_1} \boldsymbol{g}_1, \dots, \mathcal{L}_{\lambda_{N_\lambda}} \boldsymbol{g}_{N_\lambda}\}} \boldsymbol{\xi}\|$$

↔

$$\underset{\substack{\boldsymbol{g}_i \in \mathcal{C}(\Gamma), \ c_i \in \mathbb{C}^{N_\xi}}} \text{minimize} \left\| \boldsymbol{\xi} - \sum_{i=1}^{N_\lambda} c_i \mathcal{L}_{\lambda_i} \boldsymbol{g}_i \right\|$$

↔  $\boldsymbol{\xi}$  scalar & substitution  $\tilde{\boldsymbol{g}}_i = c_i \boldsymbol{g}_i$

$$\underset{\substack{\tilde{\boldsymbol{g}}_i \in \mathcal{C}(\Gamma)}}{\text{minimize}} \left\| \boldsymbol{\xi} - \sum_{i=1}^{N_\lambda} \mathcal{L}_{\lambda_i} \tilde{\boldsymbol{g}}_i \right\|$$

Convex

# Regularization

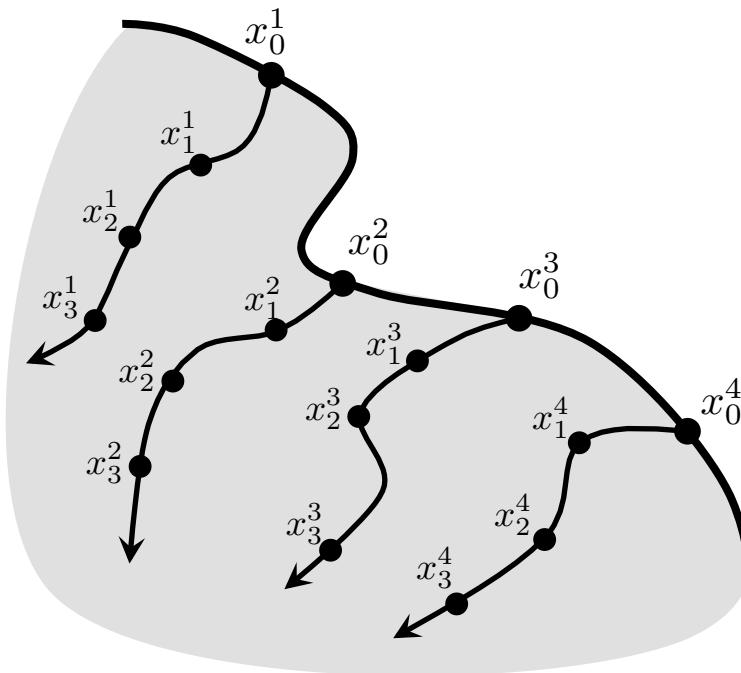
$$\underset{\boldsymbol{g}_1, \dots, \boldsymbol{g}_{N_\lambda}}{\text{minimize}} \left\| \boldsymbol{\xi} - \sum_{i=1}^{N_\lambda} \mathcal{L}_{\lambda_i} \boldsymbol{g}_i \right\| + \text{regularizer}(\boldsymbol{g}_1, \dots, \boldsymbol{g}_{N_\lambda})$$

## Examples

$$\text{regularizer} = \sum_i \text{Lipschitz}(\mathcal{L}_{\lambda_i} \boldsymbol{g}_i)$$

$$\text{regularizer} = \sum_i \int \|\nabla \mathcal{L}_{\lambda_i} \boldsymbol{g}_i\|^2$$

# Data-driven construction



When restricted to the available data set, the operators  $\mathcal{L}_i$  become matrices  
the functions  $g$  become vectors

⇒ Finite-dimensional **convex** optimization problem

$\ell_2$  norm squared + quadratic regularization ⇒ **least-squares**

# Adding control

# Adding control

$$z_{k+1} = Az_k + \textcolor{red}{B}u_k$$

$$z_0 = \hat{\phi}(x_0)$$

$$\hat{y}_k = Cz_k$$

$A, C, \hat{\phi}$  known

Minimize **multi-step** prediction error

$$\underset{\textcolor{red}{B} \in \mathbb{R}^{N \times m}}{\text{minimize}} \quad \sum_{j=1}^{\#\text{traj}} \sum_{k=1}^{\text{trajLen}} \|\xi(x_k^j) - \hat{y}_k(x_0^j)\|_2^2,$$

$\hat{y}_k$  is **linear** in  $\textcolor{red}{B}$

$$\hat{y}_k(x_0^j) = CA^k z_0^j + \sum_{i=0}^{k-1} CA^{k-i-1} \textcolor{red}{B} u_i^j$$

# Adding control

$$z_{k+1} = Az_k + \mathcal{B}u_k$$

$$z_0 = \hat{\phi}(x_0)$$

$$\hat{y}_k = Cz_k$$

$A, C, \hat{\phi}$  known

Minimize **multi-step** prediction error

$$\underset{\mathcal{B} \in \mathbb{R}^{N \times m}}{\text{minimize}} \quad \sum_{j=1}^{\#\text{traj}} \sum_{k=1}^{\text{trajLen}} \|\xi(x_k^j) - \hat{y}_k(x_0^j)\|_2^2,$$

$\hat{y}_k$  is **linear** in  $\mathcal{B}$        $\hat{y}_k(x_0^j) = CA^k z_0^j + \sum_{i=0}^{k-1} CA^{k-i-1} \mathcal{B} u_i^j$

&

$A$  and  $C$  **known**       $\Rightarrow \underset{\mathcal{B} \in \mathbb{R}^{Nm}}{\text{minimize}} \quad \|\Theta \mathcal{B} - \theta\|^2$       where       $\mathcal{b} = \text{vec}(\mathcal{B})$

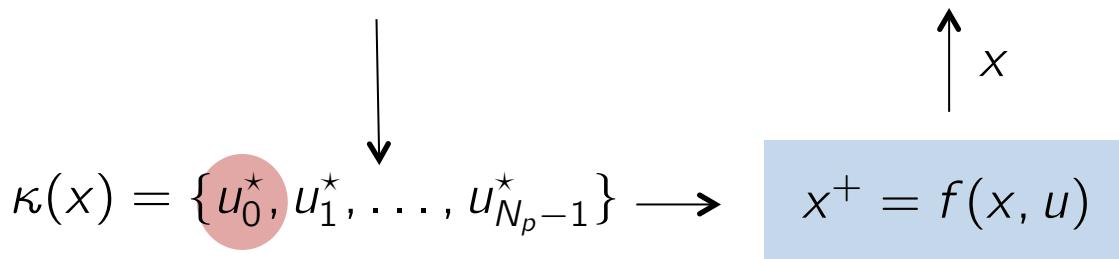
**Linear least-squares** problem

$\Rightarrow \mathcal{B} = \text{vec}^{-1}(\Theta^\dagger \theta)$

# Koopman MPC [Korda, Mezić 2018]

## Nonlinear MPC

$$\begin{array}{ll}\text{minimize}_{u_i, x_i} & \sum_{i=0}^{N_p-1} l_x(x_i) + u_i^\top R u_i + r^\top u_i \\ \text{subject to} & x_{i+1} = f(x_i, u_i), \quad i = 0, \dots, N_p - 1 \\ & c_x(x_i) + C_u u_i \leq b, \quad i = 0, \dots, N_p - 1 \\ \text{parameter} & x_0 = x\end{array}$$



# Koopman MPC [Korda, Mezić 2018]

## Koopman MPC

$$\underset{u_i, z_i, \hat{y}_i}{\text{minimize}} \quad \sum_{i=0}^{N_p-1} \hat{y}_i^\top Q \hat{y}_i + u_i^\top R u_i + q^\top \hat{y}_i + r^\top u_i$$

$$\text{subject to} \quad z_{i+1} = \mathbf{A}z_i + \mathbf{B}u_i, \quad i = 0, \dots, N_p - 1$$

$$\hat{y}_i = \mathbf{C}z_i \quad i = 0, \dots, N_p - 1$$

$$Ez_i + Fu_i \leq b, \quad i = 0, \dots, N_p - 1$$

$$\text{parameter} \quad z_0 = \hat{\phi}(x)$$

$$\kappa(x) = \{u_0^*, u_1^*, \dots, u_{N_p-1}^*\} \longrightarrow \begin{array}{c} \uparrow x \\ x^+ = f(x, u) \end{array}$$

Can handle **nonlinear constraints** and **costs** in a linear fashion

# Koopman MPC [Korda, Mezić 2018]

## Dense-form Koopman MPC

$$\underset{\mathbf{u} \in \mathbb{R}^{mN_p}}{\text{minimize}} \quad \mathbf{u}^\top H \mathbf{u} + h^\top \mathbf{u} + z_0^\top G \mathbf{u}$$

$$\text{subject to} \quad L\mathbf{u} + Mz_0 \leq c$$

$$\text{parameter} \quad z_0 = \hat{\phi}(x)$$

Convex QP!

$$\kappa(x) = \begin{bmatrix} u_0^* \\ \vdots \\ u_{N_p-1}^* \end{bmatrix} \longrightarrow \begin{array}{c} \uparrow x \\ x^+ = f(x, u) \end{array}$$

Computation cost **independent** of the size of the lift!

# Koopman MPC summary

At each step of closed-loop operation

- Set  $z_0 = \hat{\phi}(x_{\text{current}})$

- Solve

$$\begin{array}{ll}\text{minimize}_{\mathbf{u} \in \mathbb{R}^{mN_p}} & \mathbf{u}^\top H \mathbf{u}^\top + h^\top \mathbf{u} + z_0^\top G \mathbf{u} \\ \text{subject to} & L \mathbf{u} + M z_0 \leq c\end{array}$$

$$\Rightarrow \mathbf{u}^* = \begin{bmatrix} u_0^* \\ \vdots \\ u_{N_p-1}^* \end{bmatrix}$$

- Apply  $u_0^*$  to the system

Main benefits

**Data-driven:** No model required

**Fast & simple:** only small **convex quadratic program** solved online

**Nonlinear constraints** and **costs** handled in a linear fashion

# Numerical examples

# Numerical examples – damped Duffing

## Dynamics

$$\dot{x}_1 = x_2$$

$$\dot{x}_2 = -0.5x_2 - x_1(4x_1^2 - 1) + 0.5\textcolor{red}{u}$$

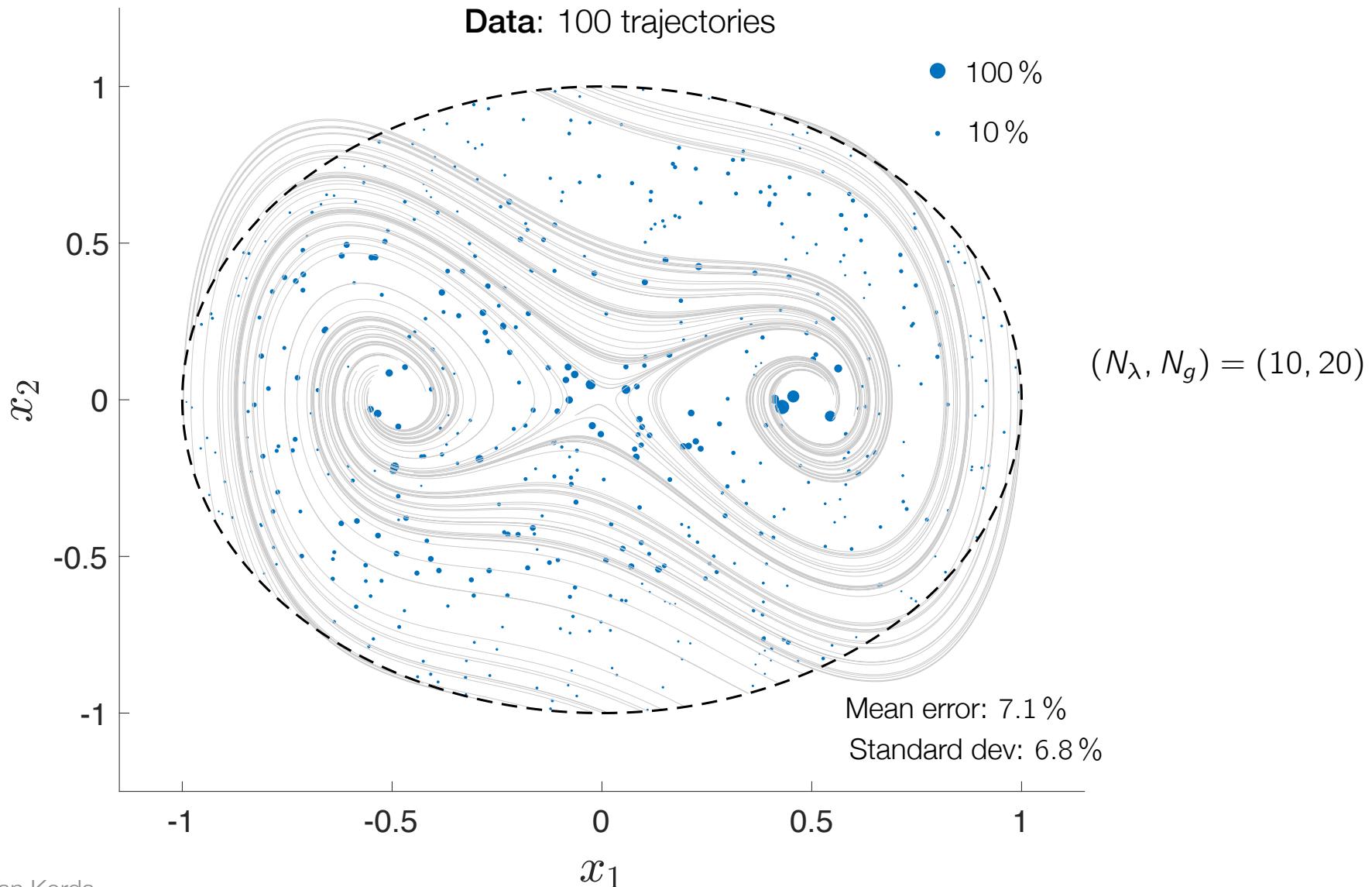
**Data:** 100 trajectories, 8 second long

**Eigenvalues:** Lattice from DMD eigenvalues

**Boundary functions:** Thin plate spline RBFs

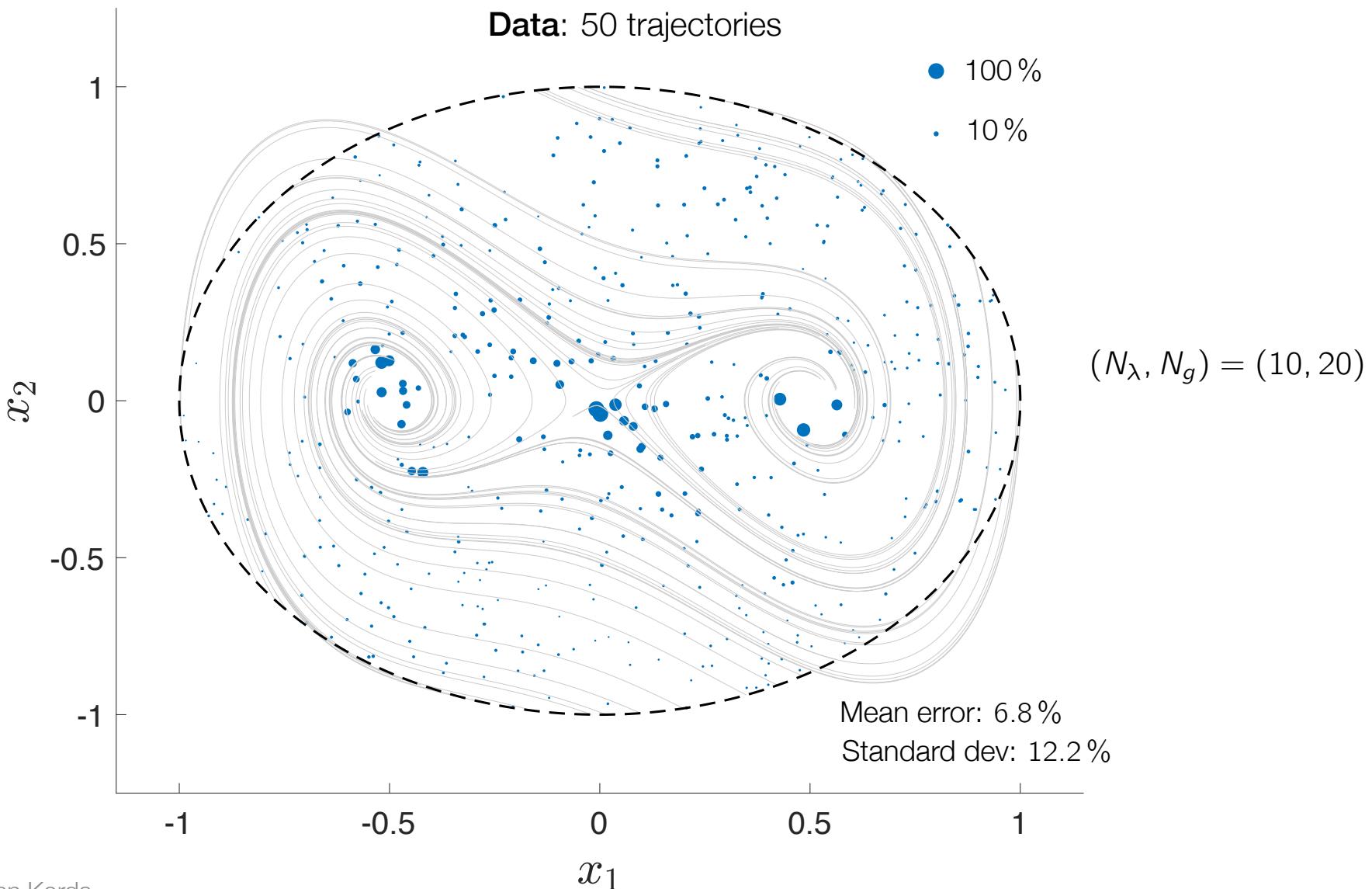
# Numerical examples – damped Duffing

Spatial distribution of one-second prediction error (with control)

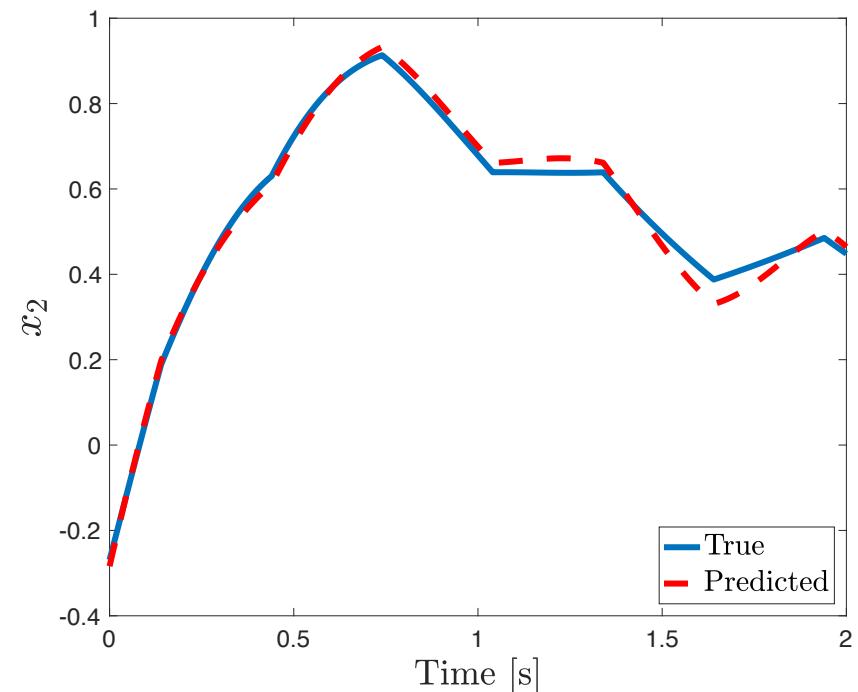
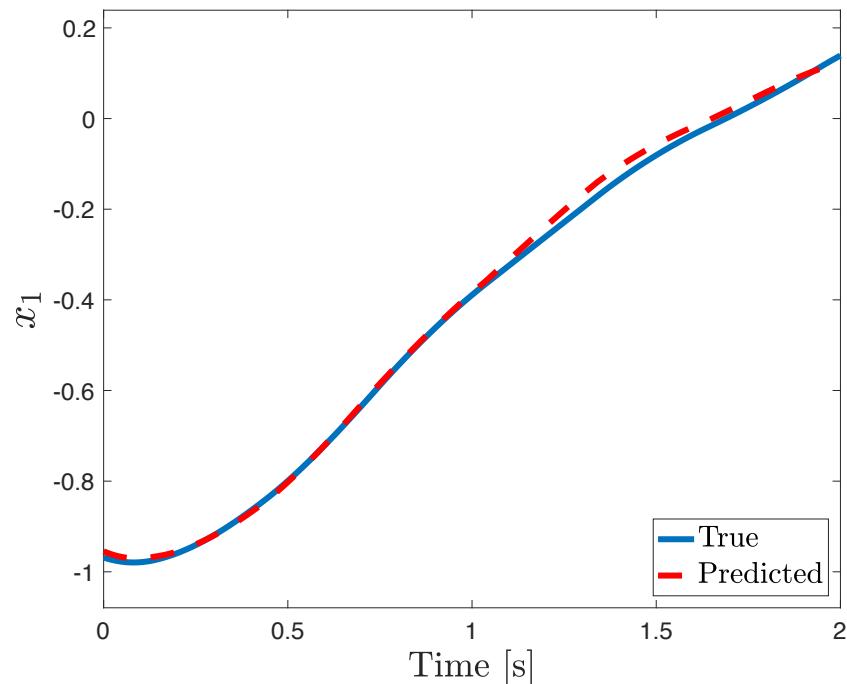


# Numerical examples – damped Duffing

Spatial distribution of one-second prediction error (with control)



# Numerical examples – damped Duffing



$$(N_\lambda, N_g) = (10, 20)$$

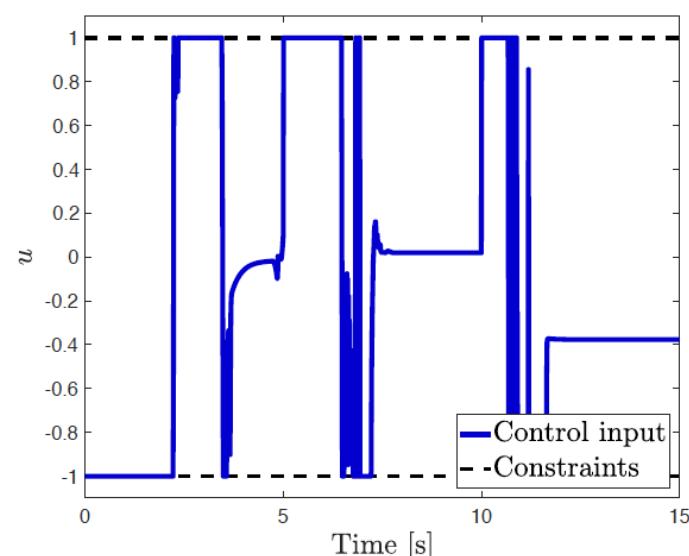
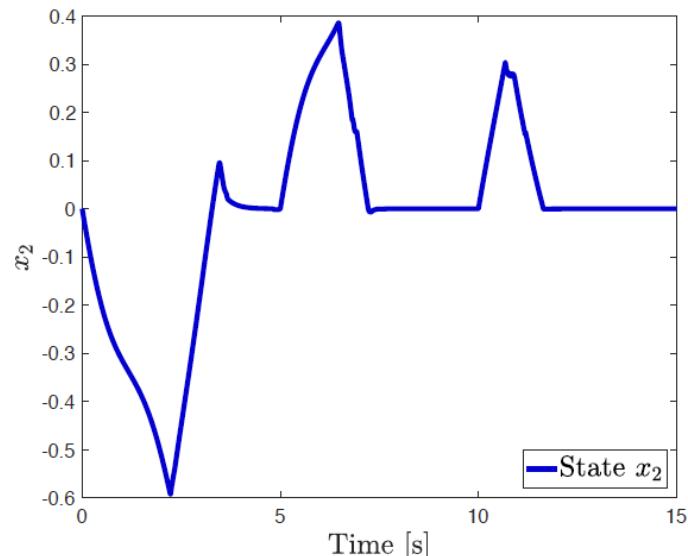
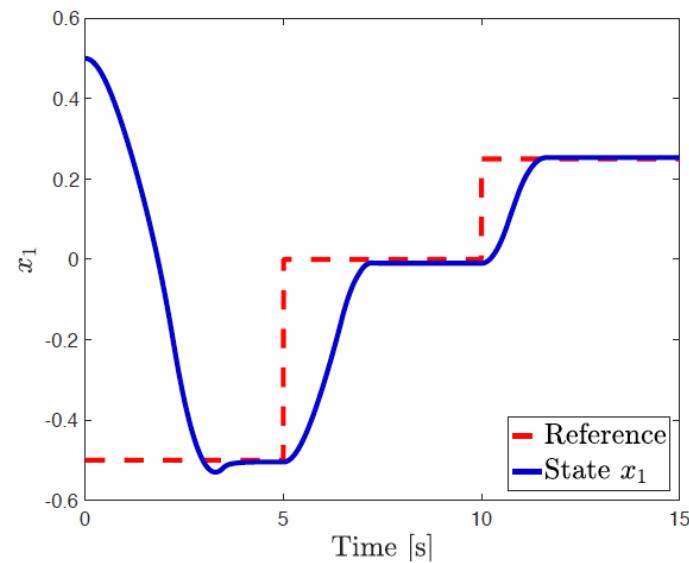
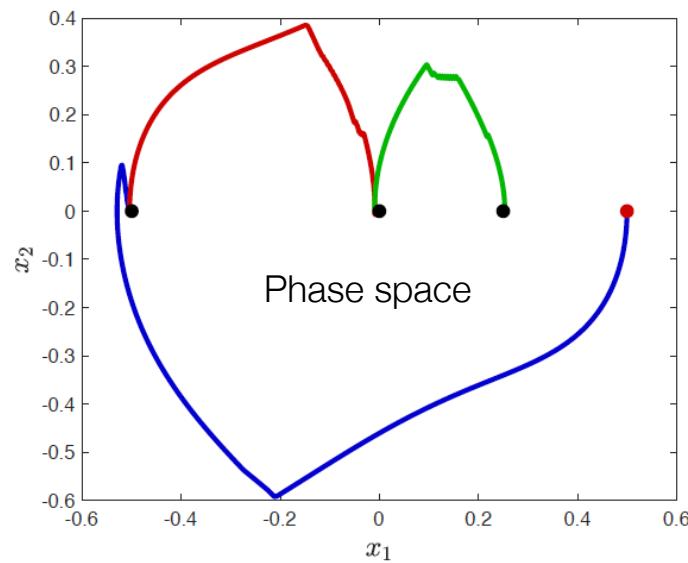
# Numerical examples – damped Duffing

$(N_A, N_G)$	(10, 30)	(10, 20)	(6, 20)	(10, 10)	(10, 5)	(10, 3)
Mean error [uncontrolled]	6.9 %	8.9 %	17.4 %	19.9 %	38.8 %	56.2 %
Mean error [controlled]	4.6 %	6.7 %	15.8 %	15.7 %	35.6 %	53.5 %

EDMD error (200 RBF basis functions) = 25.1 %

# Numerical examples – damped Duffing

Feedback control – Koopman MPC



# Numerical examples – Van der Pol

## Dynamics

$$\dot{x}_1 = 2x_2$$

$$\dot{x}_2 = -0.8x_1 + 2x_2 - 10x_1^2x_2 + \textcolor{red}{u}$$

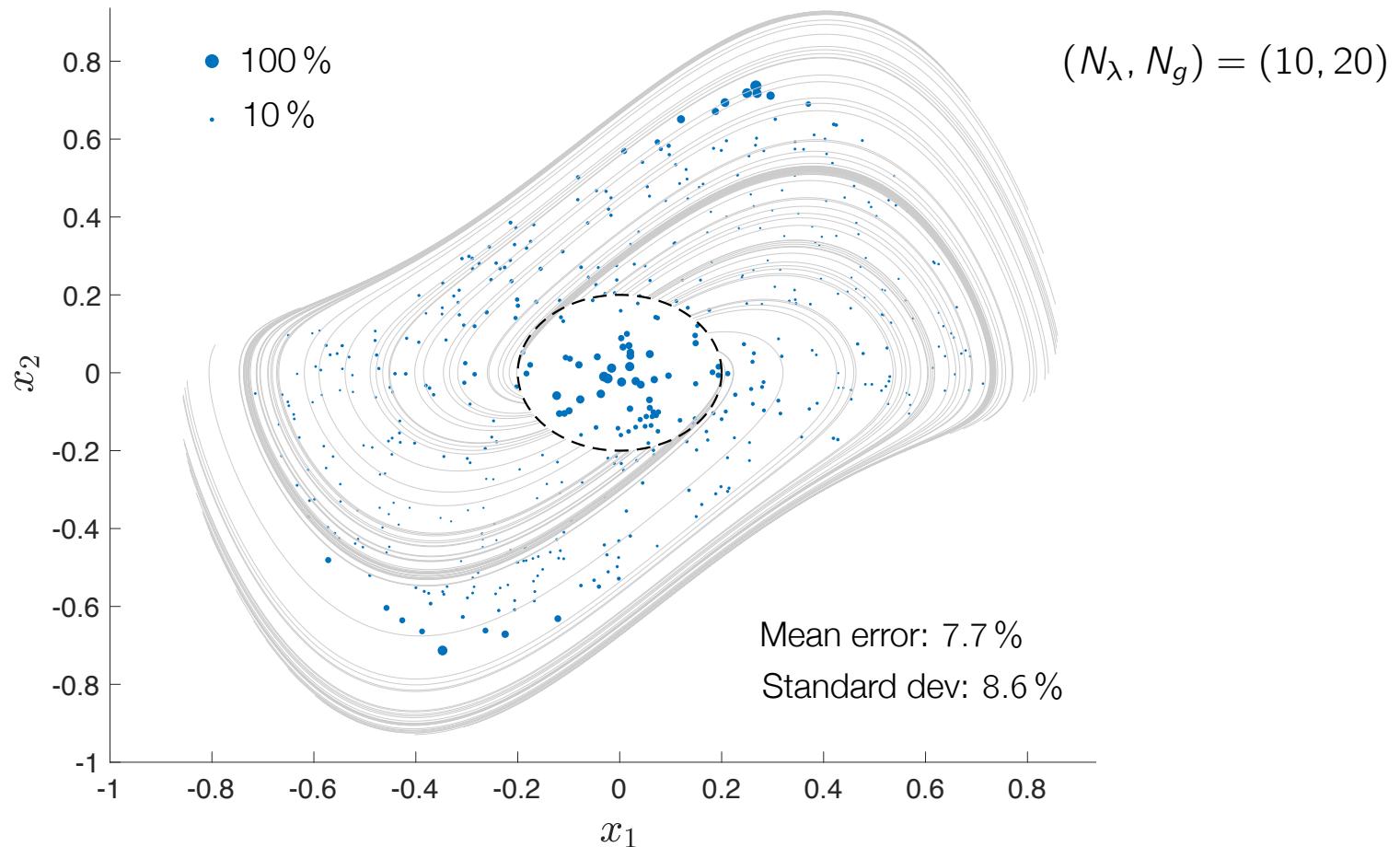
**Data:** 100 trajectories, 3 second long

**Eigenvalues:** Lattice from DMD eigenvalues

**Boundary functions:** Thin plate spline RBFs

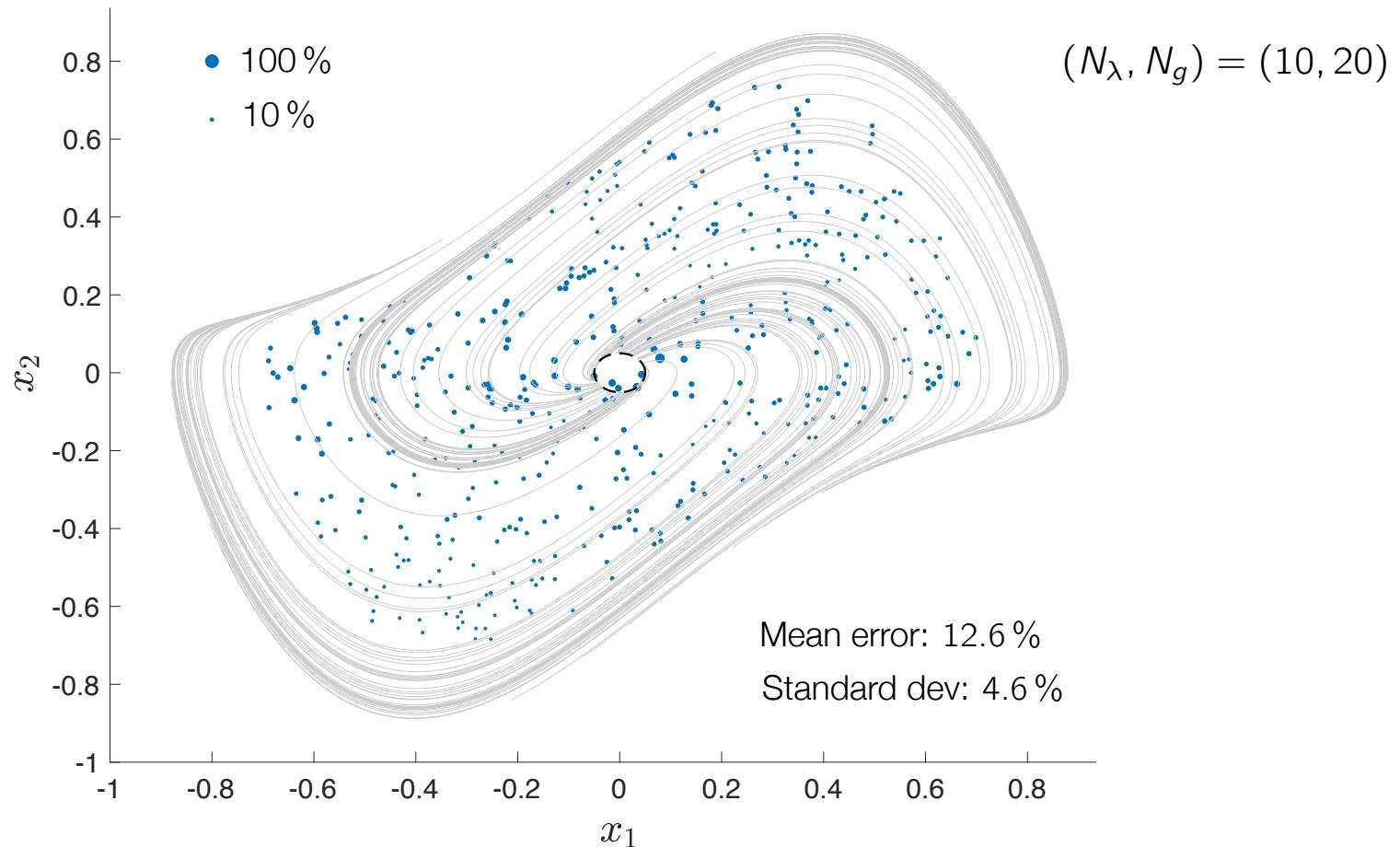
# Numerical examples – Van der Pol

Spatial distribution of one-second prediction error (with control)

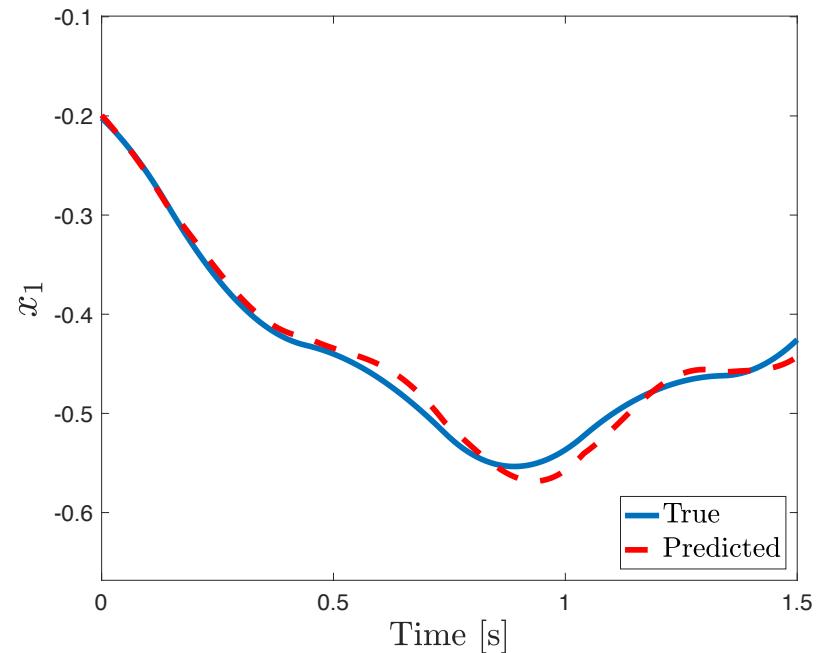
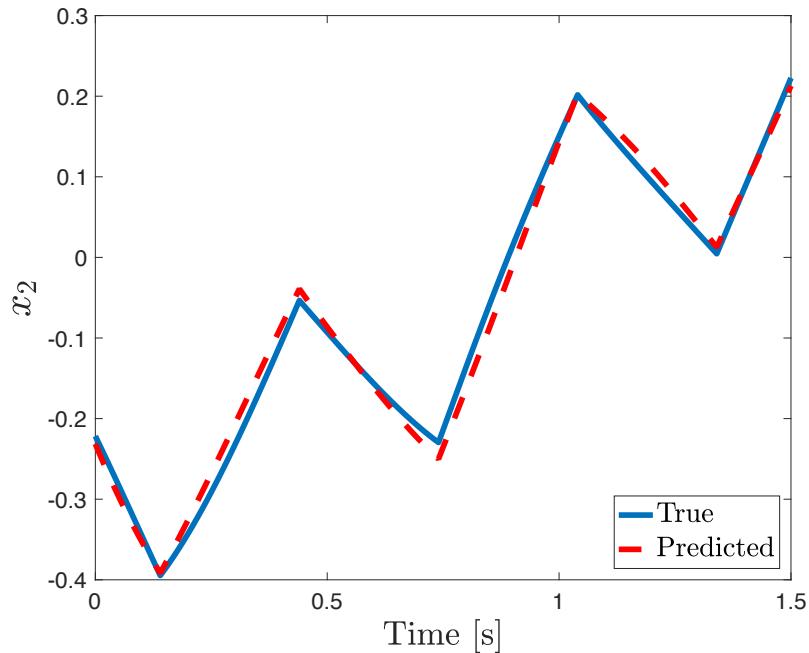


# Numerical examples – Van der Pol

Spatial distribution of one-second prediction error (with control)



# Numerical examples – Van der Pol



$$(N_\lambda, N_g) = (10, 20)$$

# Numerical examples – Van der Pol

Mean prediction error for different number of eigenfunctions

$(N_\lambda, N_g)$	(10, 20)	(6, 20)	(10, 10)	(10, 5)	(10, 3)
Mean error [uncontrolled]	10.4 %	18.5 %	14.0 %	26.4 %	33.4 %
Mean error [controlled]	12.6 %	18.3 %	16.0 %	26.5 %	34.2 %

# Optimized x Not optimized

Total # of eigenfunctions

Without optimization of  $g$

$$N = N_\lambda N_G$$

With optimization of  $g$

$$N = N_\lambda N_\xi$$

# Optimized x Not optimized

Total # of eigenfunctions

Without optimization of  $g$

$$N = N_\lambda N_G$$

With optimization of  $g$

$$N = N_\lambda N_\xi$$

Duffing	Not optimized						Optimized		
	$N$	300	200	120	100	50	30	20	12
Mean error [controlled]		4.6 %	6.7 %	15.8 %	15.7 %	35.6 %	53.5 %	5.03 %	15.4 %

# Optimized x Not optimized

Total # of eigenfunctions

Without optimization of  $g$

$$N = N_\lambda N_G$$

With optimization of  $g$

$$N = N_\lambda N_\xi$$

Duffing	Not optimized						Optimized		
	$N$	300	200	120	100	50	30	20	12
Mean error [controlled]		4.6 %	6.7 %	15.8 %	15.7 %	35.6 %	53.5 %	5.03 %	15.4 %

Van der Pol	Not optimized						Optimized		
	$N$	300	200	120	100	50	30	30	20
Mean error [controlled]		11.6 %	12.6 %	18.3 %	16.0 %	26.5 %	34.2 %	7.6 %	13.6 %

# Open problems

# Open problem 1:

Optimal choice of  $g$  **jointly** for all components of  $\xi$

$$\underset{\mathbf{x} \in \mathbb{R}^n}{\text{maximize}} \quad \mathbf{x}^\top A^\top \left( \sum_i B_i \mathbf{x} \mathbf{x}^\top B_i^\top \right)^{-1} A \mathbf{x}$$

**Remarks:** Homogenous of degree zero

Similar to (generalized) eigenvalue problem

# Open problem 2:

Optimal choice of eigenvalues

$$\phi_{\lambda,g}(S_t(x_0)) = e^{\lambda t} g(x_0)$$

**Remarks:** Log-linear in  $\lambda$

Geometric programming?

Low-dimensional

# Conclusion

- Data-driven construction of Koopman eigenfunctions

Only linear algebra and/or convex optimization needed

Readily applicable to control and estimation

Seems very robust

## Future work

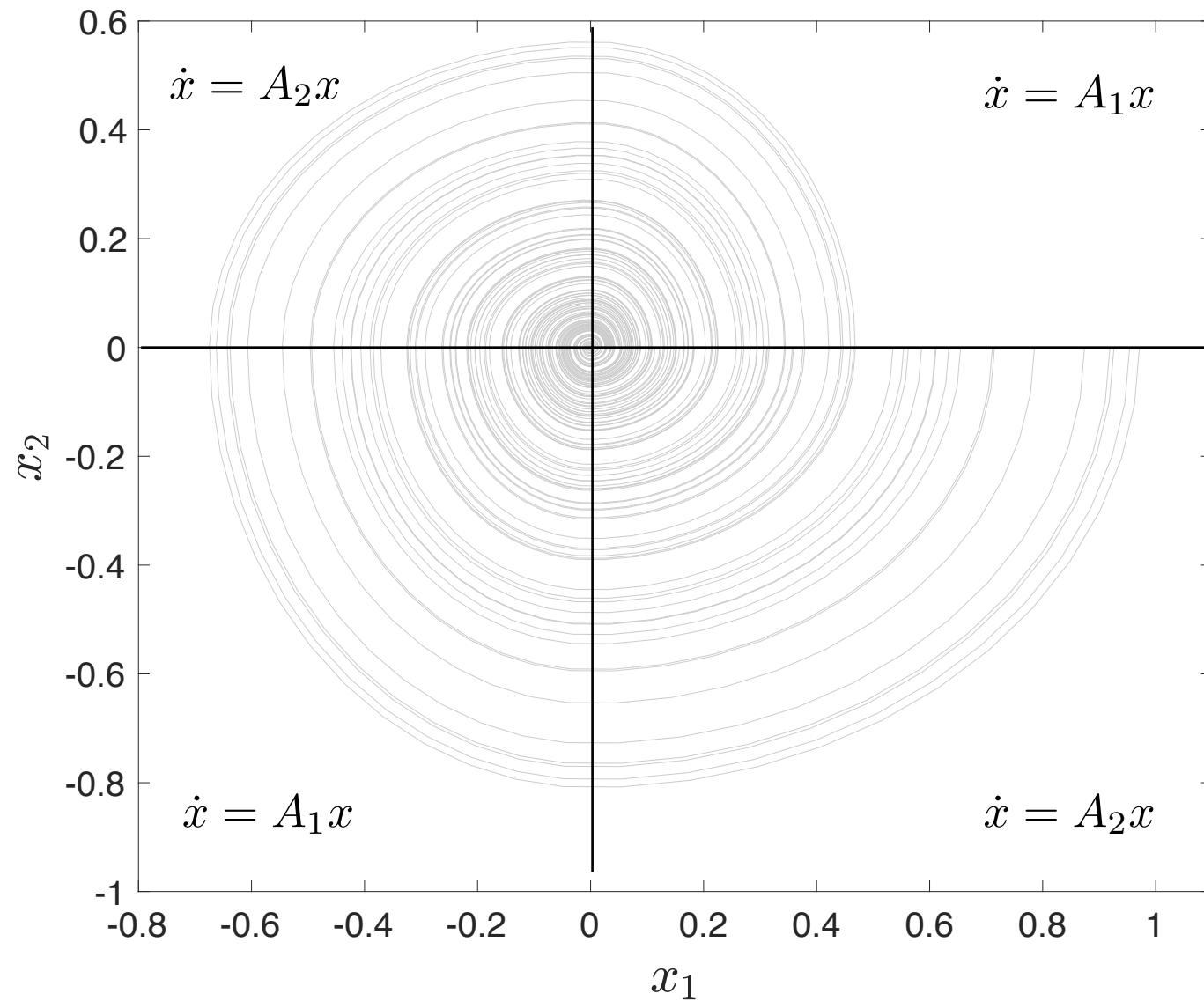
- High dimensional interpolation / approximation
- Exploit **algebraic structure** (products of eigenfunctions)

$\phi_1, \dots, \phi_N$  eigenfunctions  $\Rightarrow \phi_1^{p_1} \cdot \dots \cdot \phi_N^{p_N}$  also an eigenfunction

- Generalized eigenfunctions – Jordan blocks

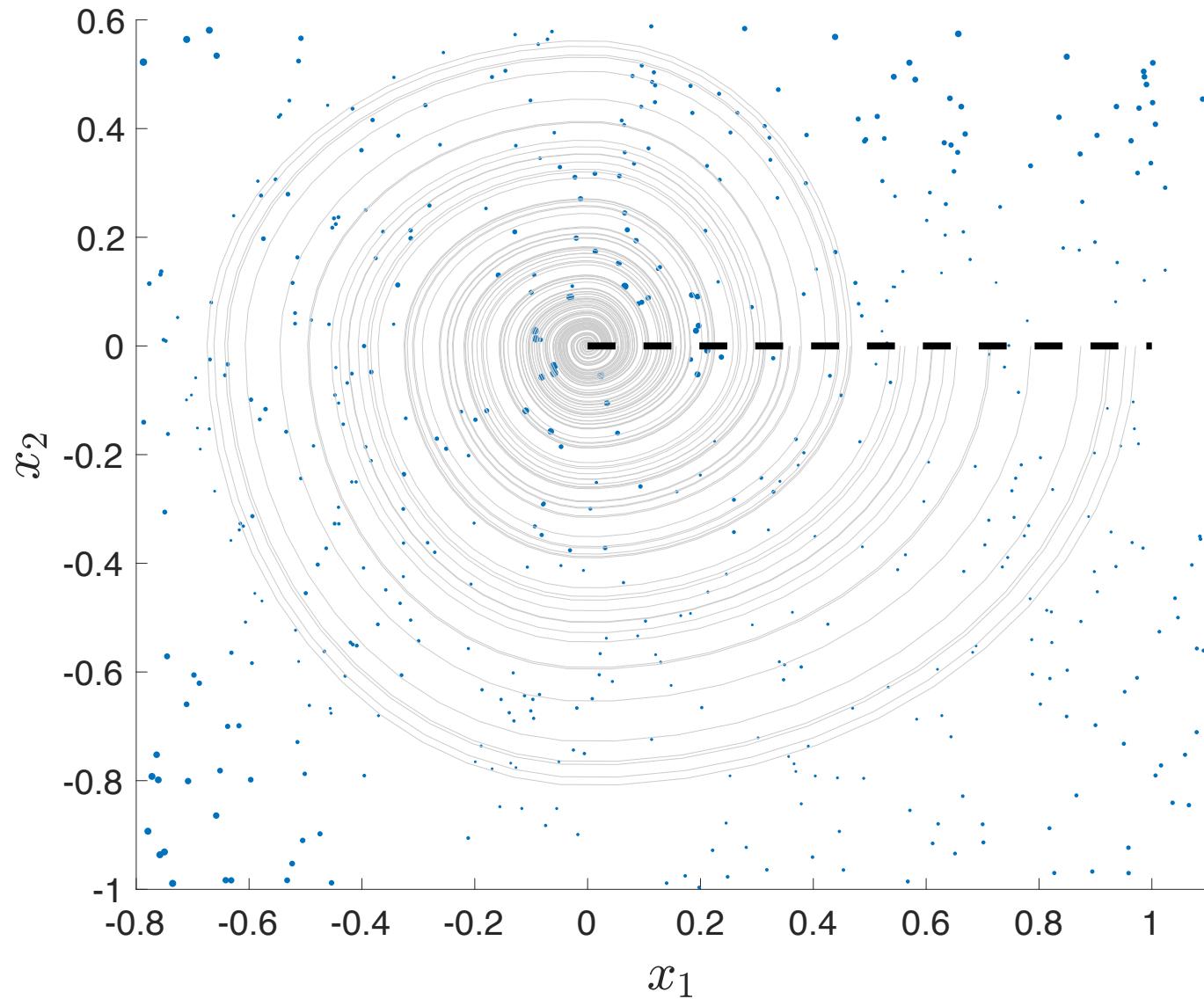
$$\begin{bmatrix} \phi_1(x(t)) \\ \phi_2(x(t)) \end{bmatrix} := \exp\left(t \begin{bmatrix} \lambda & 1 \\ 0 & \lambda \end{bmatrix}\right) \begin{bmatrix} g_1(x_0) \\ g_2(x_0) \end{bmatrix} \quad \Rightarrow \quad \text{span}\{\phi_1, \phi_2\} \text{ is invariant!}$$

# Switched linear system



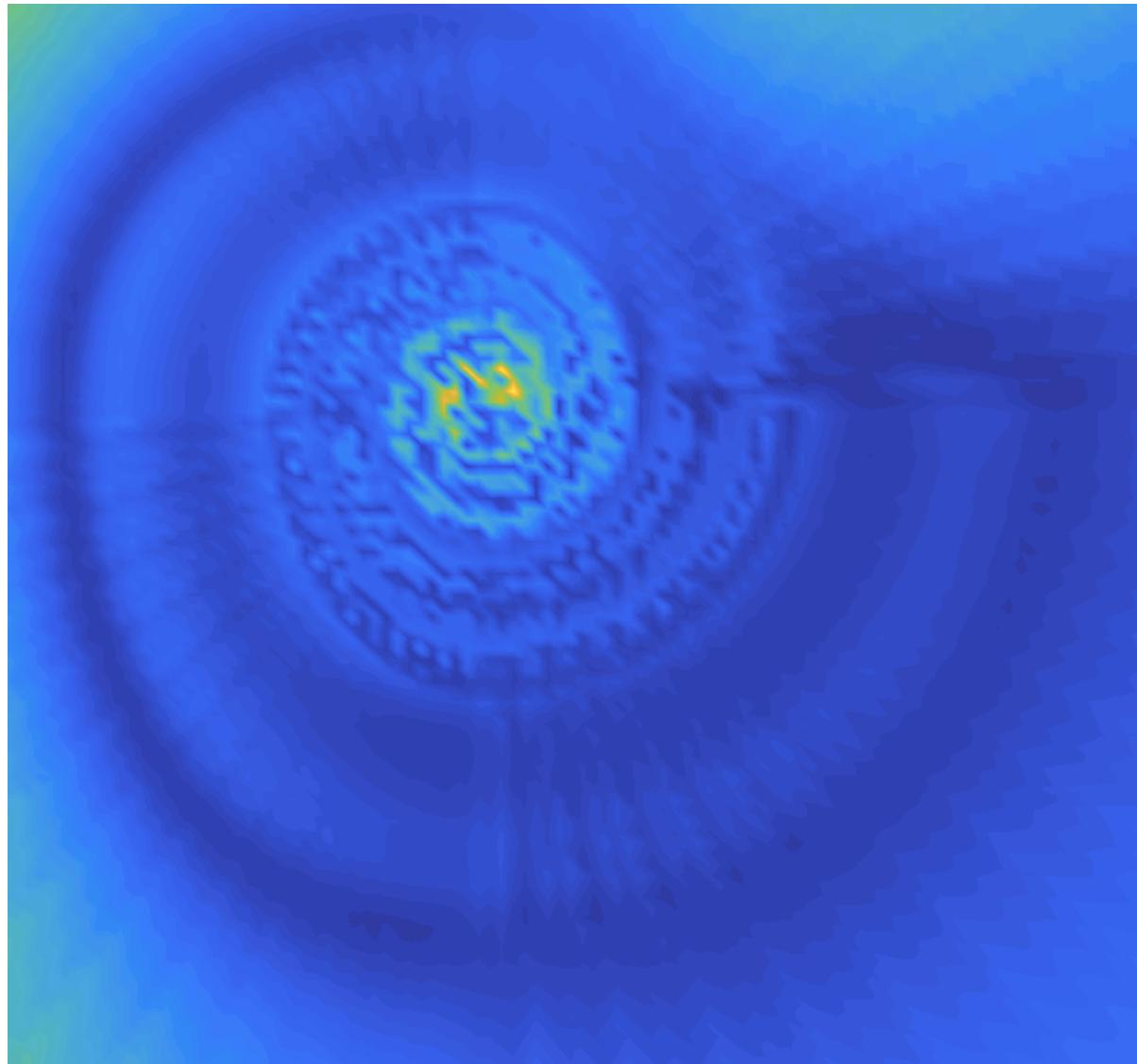
# Switched linear system

Mean error = 5.8 %  
Standard dev. = 3.4 %



# Switched linear system

Mean error = 5.8 %  
Standard dev. = 3.4 %



# Switched linear system

Mean error = 5.8 %  
Standard dev. = 3.4 %

