

Reliable Machine Learning for Dynamical Systems: From Operator Theory to Sea-Ice Forecasting

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Papers and talks:

[http://www.damtp.cam.ac.uk/
user/mjc249/home.html](http://www.damtp.cam.ac.uk/user/mjc249/home.html)

What is a Koopman operator?

- $\mathcal{X}$ – *the state space*
- $\mathcal{X} \ni x$ – *the state*

cts $F: \mathcal{X} \rightarrow \mathcal{X}$ – *the dynamics*: $x_{n+1} = F(x_n)$

Henri Poincaré
(Sorbonne)



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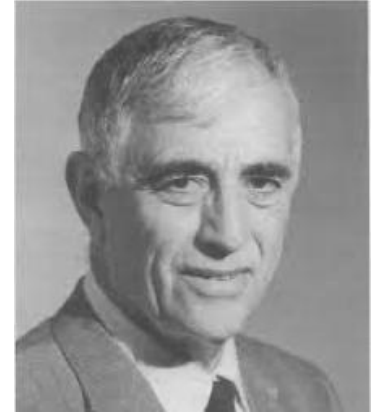
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cts $F: \mathcal{X} \rightarrow \mathcal{X}$ – the dynamics: $x_{n+1} = F(x_n)$

- Functions $g: \mathcal{X} \rightarrow \mathbb{C}$ a.k.a “observables”, $g \in L^2(\mathcal{X}, \omega)$
- Koopman operator $\mathcal{K}_F: [\mathcal{K}_F g](x) = g(F(x))$ LINEAR!


 Observe g one time step forward

Bernard Koopman
(Columbia)



John von Neumann
(IAS)



- Koopman, “Hamiltonian systems and transformation in Hilbert space,” **Proc. Natl. Acad. Sci. USA**, 1931.
- Koopman, v. Neumann, “Dynamical systems of continuous spectra,” **Proc. Natl. Acad. Sci. USA**, 1932.

What is a Koopman operator?

- $\mathcal{X}$ – the state space
- $\mathcal{X} \ni x$ – the state
- Unknown cts $F: \mathcal{X} \rightarrow \mathcal{X}$ – the dynamics: $x_{n+1} = F(x_n)$
- Functions $g: \mathcal{X} \rightarrow \mathbb{C}$ a.k.a “observables”, $g \in L^2(\mathcal{X}, \omega)$
- Koopman operator $\mathcal{K}_F: [\mathcal{K}_F g](x) = g(F(x))$ **LINEAR!**
- Available snapshot data: $\left\{ \left(x^{(m)}, y^{(m)} = F(x^{(m)}) \right) : m = 1, \dots, M \right\}$

Can we compute spectral properties from trajectory data?

$$g(x_n) = [\mathcal{K}^n g](x_0)$$

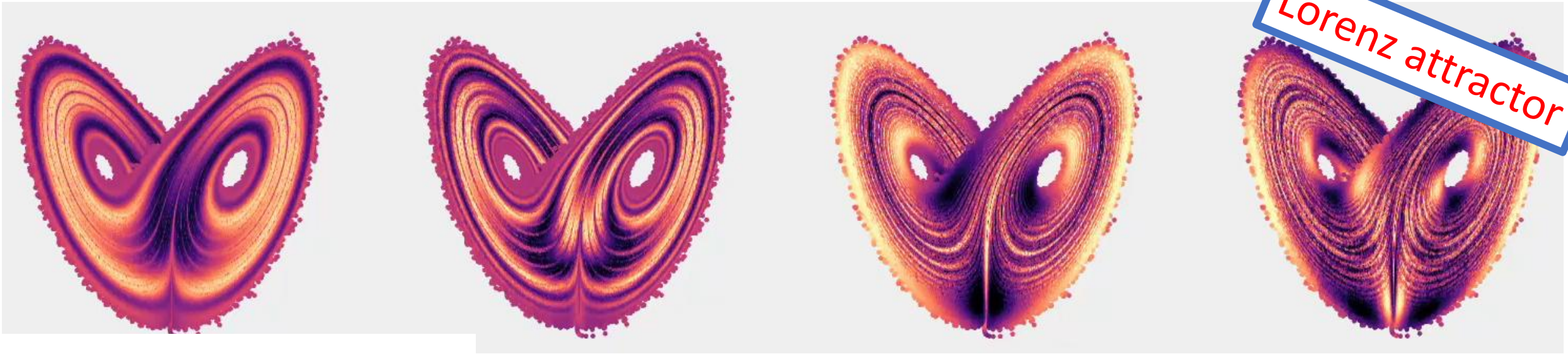
Why?

If $\|\mathcal{K}g - \lambda g\| \leq \varepsilon$, then $g(x_n) = [\mathcal{K}^n g](x_0) = \lambda^n g(x_0) + \mathcal{O}(n\varepsilon)$

$$g(x_n) = [\mathcal{K}^n g](x_0)$$

Why?

If $\|\mathcal{K}g - \lambda g\| \leq \varepsilon$, then $g(x_n) = [\mathcal{K}^n g](x_0) = \lambda^n g(x_0) + \mathcal{O}(n\varepsilon)$



Coherent features!

$$\text{Sp}_\varepsilon(\mathcal{K}) = \{z \in \mathbb{C} : \exists g, \|g\| = 1, \|\mathcal{K}g - zg\| \leq \varepsilon\}$$

Trades: Nonlinear, finite-dimensional $\Rightarrow$ Linear, infinite-dimensional.

Koopman Mode Decomposition

- Find (g_j, λ_j) with $\|\mathcal{K}g_j - \lambda_j g_j\| \leq \varepsilon$
- Expand observable:

Verified Eigenfunctions

$$h(x) \approx \sum_j c_j g_j(x)$$

coefficients, called
"Koopman modes"

- Forecasts:

$$h(x_n) = \sum_j \lambda_j^n c_j g_j(x) + \mathcal{O}(n\varepsilon)$$

Intuition: A nonlinear separation of variables through a linear operator!

Perils of discretization: Warmup on $\ell^2(\mathbb{Z})$

$$\begin{pmatrix} \ddots & & & & \\ & \ddots & & & \\ & 0 & 1 & & \\ & & 0 & 1 & \\ & & & 0 & 1 \\ & & & & 0 & \ddots \end{pmatrix} \xrightarrow{\text{Two-way infinite}} \begin{pmatrix} 0 & 1 & & & \\ & \ddots & \ddots & & \\ & & \ddots & 1 & \\ & & & \ddots & 0 \end{pmatrix} \in \mathbb{C}^{N \times N}$$

- Spectrum is unit circle.
- Spectrum is stable.
- Continuous spectra.
- Unitary evolution.

- Spectrum is $\{0\}$.
- Spectrum is unstable.
- Discrete spectra.
- Nilpotent evolution.

Lots of Koopman operators are built up from operators like these!

Matrix approximation of $\mathcal{K}$ (EDMD)

Observables $\psi_j: \mathcal{X} \rightarrow \mathbb{C}, j = 1, \dots, N$

$$\{x^{(m)}, y^{(m)} = F(x^{(m)})\}_{m=1}^M$$

quadrature points

$$\langle \psi_k, \psi_j \rangle \approx \sum_{m=1}^M w_m \overline{\psi_j(x^{(m)})} \psi_k(x^{(m)}) = \left[\underbrace{\begin{pmatrix} \psi_1(x^{(1)}) & \dots & \psi_N(x^{(1)}) \\ \vdots & \ddots & \vdots \\ \psi_1(x^{(M)}) & \dots & \psi_N(x^{(M)}) \end{pmatrix}}_{\Psi_X} \right]^* \underbrace{\begin{pmatrix} w_1 & & \\ & \ddots & \\ & & w_M \end{pmatrix}}_W \underbrace{\begin{pmatrix} \psi_1(x^{(1)}) & \dots & \psi_N(x^{(1)}) \\ \vdots & \ddots & \vdots \\ \psi_1(x^{(M)}) & \dots & \psi_N(x^{(M)}) \end{pmatrix}}_{\Psi_X} \right]_{jk}$$

quadrature weights

$$\langle \mathcal{K}\psi_k, \psi_j \rangle \approx \sum_{m=1}^M w_m \overline{\psi_j(x^{(m)})} \underbrace{\psi_k(y^{(m)})}_{[\mathcal{K}\psi_k](x^{(m)})} = \left[\underbrace{\begin{pmatrix} \psi_1(x^{(1)}) & \dots & \psi_N(x^{(1)}) \\ \vdots & \ddots & \vdots \\ \psi_1(x^{(M)}) & \dots & \psi_N(x^{(M)}) \end{pmatrix}}_{\Psi_X} \right]^* \underbrace{\begin{pmatrix} w_1 & & \\ & \ddots & \\ & & w_M \end{pmatrix}}_W \underbrace{\begin{pmatrix} \psi_1(y^{(1)}) & \dots & \psi_N(y^{(1)}) \\ \vdots & \ddots & \vdots \\ \psi_1(y^{(M)}) & \dots & \psi_N(y^{(M)}) \end{pmatrix}}_{\Psi_Y} \right]_{jk}$$

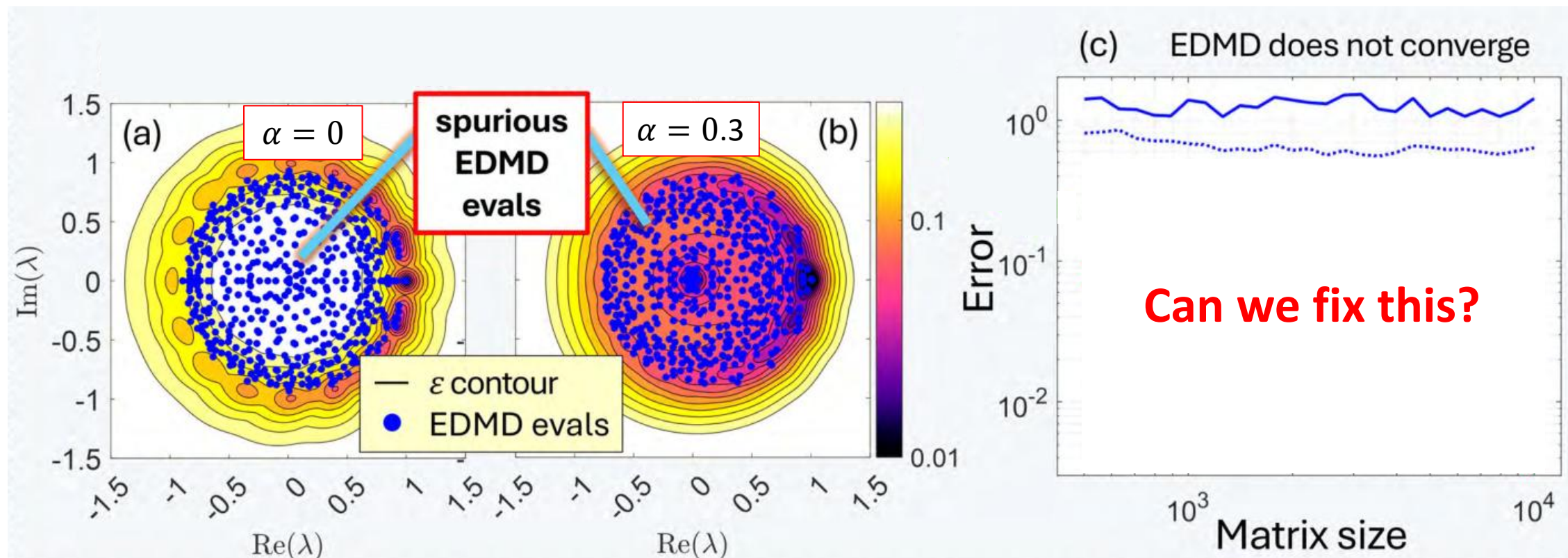
Galerkin
Approximation

$$\mathcal{K} \rightarrow (\Psi_X^* W \Psi_X)^{-1} \Psi_X^* W \Psi_Y \in \mathbb{C}^{N \times N}$$

- Schmid, "Dynamic mode decomposition of numerical and experimental data," **J. Fluid Mech.**, 2010.
- Rowley, Mezić, Bagheri, Schlatter, Henningson, "Spectral analysis of nonlinear flows," **J. Fluid Mech.**, 2009.
- Williams, Kevrekidis, Rowley "A data-driven approximation of the Koopman operator: Extending dynamic mode decomposition," **J. Nonlinear Sci.**, 2015.

EDMD doesn't converge!

- Duffing oscillator: $\dot{x} = y$, $\dot{y} = -\alpha y + x(1 - x^2)$, sampled $\Delta t = 0.3$.
- Gaussian radial basis functions, Monte Carlo integration ($M = 50000$)



The fix: Residual DMD (ResDMD)

$$\langle \psi_k, \psi_j \rangle \approx \sum_{m=1}^M w_m \overline{\psi_j(x^{(m)})} \psi_k(x^{(m)}) = \left[\underbrace{\Psi_X^* W \Psi_X}_G \right]_{jk}$$

$$\langle \mathcal{K}\psi_k, \psi_j \rangle \approx \sum_{m=1}^M w_m \overline{\psi_j(x^{(m)})} \underbrace{\psi_k(y^{(m)})}_{[\mathcal{K}\psi_k](x^{(m)})} = \left[\underbrace{\Psi_X^* W \Psi_Y}_A \right]_{jk}$$

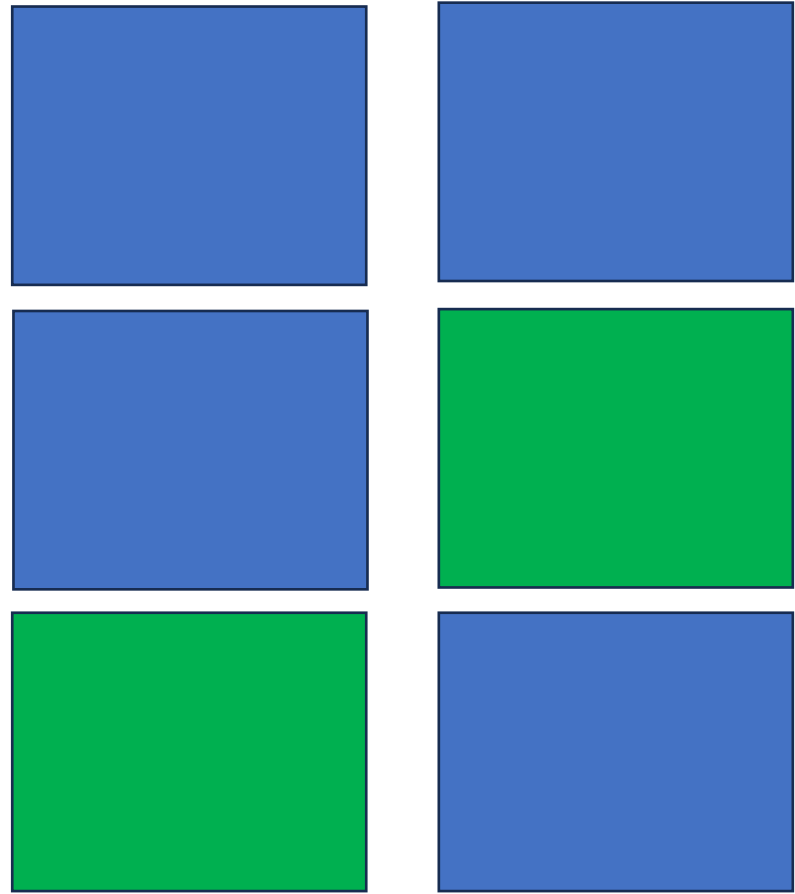
- C., Townsend, “Rigorous data-driven computation of spectral properties of Koopman operators for dynamical systems,” **Commun. Pure Appl. Math.**, 2023.
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adjoint



- C., Townsend, “Rigorous data-driven computation of spectral properties of Koopman operators for dynamical systems,” **Commun. Pure Appl. Math.**, 2023.
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- C., Townsend, “Rigorous data-driven computation of spectral properties of Koopman operators for dynamical systems,” **Commun. Pure Appl. Math.**, 2023.
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Residuals: $g = \sum_{j=1}^N \mathbf{g}_j \psi_j$, $\|\mathcal{K}g - \lambda g\|^2 = \langle \mathcal{K}g - \lambda g, \mathcal{K}g - \lambda g \rangle$

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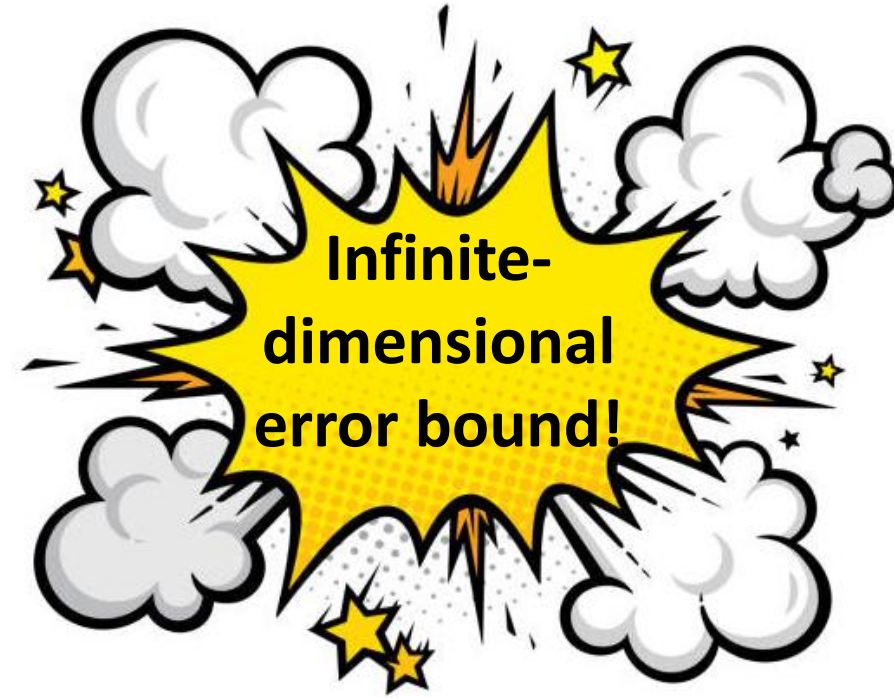
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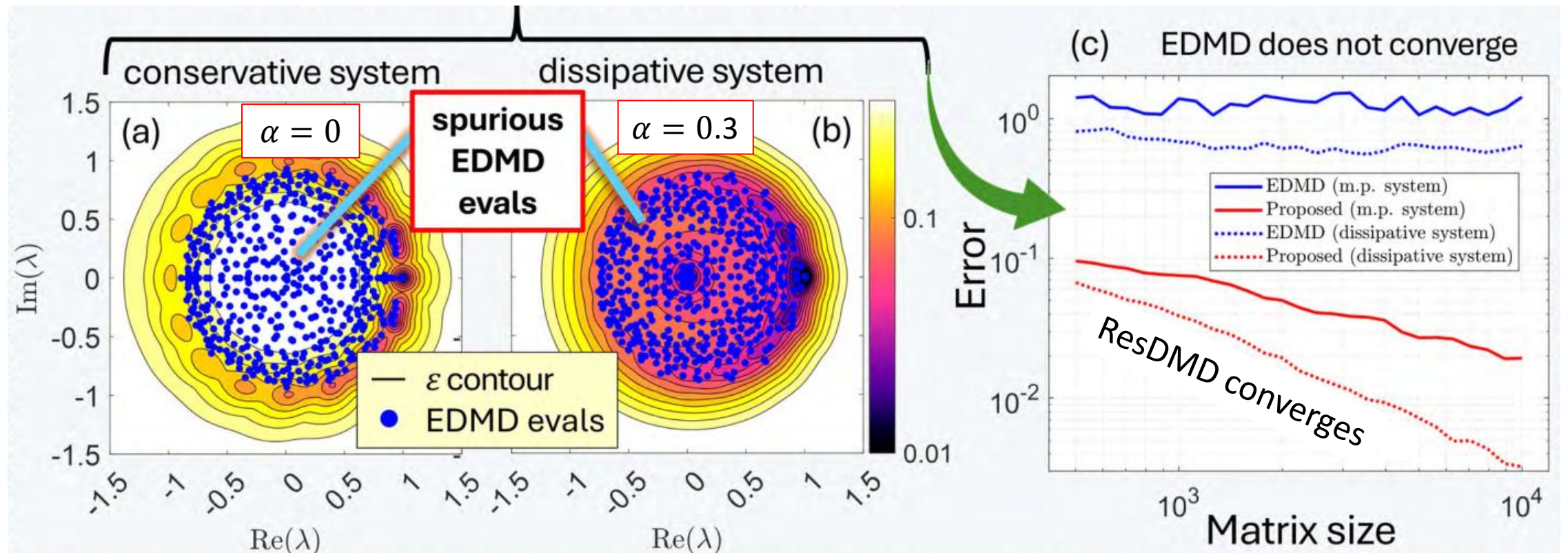
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ResDMD does converge!

- Duffing oscillator: $\dot{x} = y$, $\dot{y} = -\alpha y + x(1 - x^2)$, sampled $\Delta t = 0.3$.
- Gaussian radial basis functions, Monte Carlo integration ($M = 50000$)

Compute $\text{Sp}_\varepsilon(\mathcal{K})$, local adaptive control on $\varepsilon \downarrow 0$



What can we do?

Consider space of observables with finite energy: $L^2(\mathcal{X}, \omega)$

Theorem: There **exists** algorithms $\Gamma_{N,M}$ using snapshots such that

$$\lim_{N \rightarrow \infty} \lim_{M \rightarrow \infty} \Gamma_{N,M}(F) = \text{Sp}_\varepsilon(\mathcal{K}_F)$$

for all systems.



N = size of basis, M = amount of data (quadrature)

$$\text{Sp}_\varepsilon(\mathcal{K}) = \{z \in \mathbb{C} : \exists g, \|g\| = 1, \|\mathcal{K}g - zg\| \leq \varepsilon\}$$

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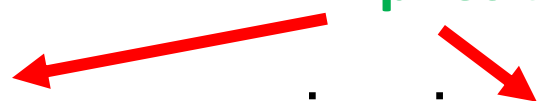
size of basis, M = amount of data (quadrature)

Double limit $\lim_{N \rightarrow \infty} \lim_{M \rightarrow \infty}$

Can we do better?

Adversaries: Double limit is necessary!

Implies $\mathcal{K}$ is unitary



Class of systems: $\Omega_{\mathbb{D}} = \{F: \bar{\mathbb{D}} \rightarrow \bar{\mathbb{D}} \mid F \text{ cts, measure preserving, invertible}\}.$

Data an algorithm can use: $\mathcal{T}_F = \{(x, y_m) \mid x \in \bar{\mathbb{D}}, \|F(x) - y_m\| \leq 2^{-m}\}.$

Theorem: There **does not exist** any sequence of deterministic algorithms $\{\Gamma_n\}$ using $\mathcal{T}_F$ such that $\lim_{n \rightarrow \infty} \Gamma_n(F) = \text{Sp}_{\varepsilon}(\mathcal{K}_F) \forall F \in \Omega_{\mathbb{D}}.$

NB:

- n can index anything.
- Universal - any type of algorithm or computational model.
- Similarly, no random algorithms converging with probability $> 1/2$.

Proof idea: Constructing an adversary

$$F_0: \text{rotation by } \pi, \operatorname{Sp}(\mathcal{K}_{F_0}) = \{\pm 1\}$$

Phase transition lemma: Let $X = \{x_1, \dots, x_N\}, Y = \{y_1, \dots, y_N\}$ be distinct points in annulus $\mathcal{A} = \{x \in \mathbb{D} \mid 0 < R < \|x\| < r < 1\}$ with $X \cap Y = \emptyset$. There exists a measure-preserving homeomorphism H such that H acts as the identity on $\mathbb{D} \setminus \mathcal{A}$ and $H(y_j) = F_0(H(x_j)), j = 1, \dots, N$.

Conjugacy of data ($x_j \rightarrow y_j$) with F_0

Idea: Use lemma to trick any algorithm into oscillating between spectra.

Proof idea: Constructing an adversary

Suppose (for contradiction) $\{\Gamma_n\}$ uses $\mathcal{T}_F$, $\lim_{n \rightarrow \infty} \Gamma_n(F) = \text{Sp}(\mathcal{K}_F) \quad \forall F \in \Omega_{\mathbb{D}}$.

Build an **adversarial** F ...

$$\mathcal{T}_F = \{(x, y_m) \mid \|F(x) - y_m\| \leq 2^{-m}\}$$

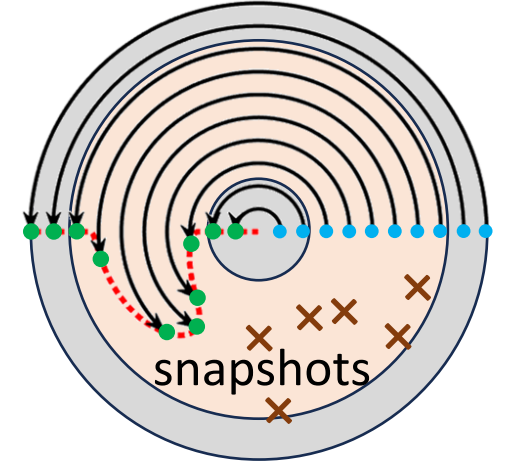
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Build an **adversarial** F ...

$$\widetilde{F}_1(r, \theta) = (r, \theta + \pi + \phi(r)), \text{ supp}(\phi) \subset [1/4, 3/4]$$

$$\text{Sp}(\mathcal{K}_{\widetilde{F}_1}) = \mathbb{T} \text{ (unit circle).}$$



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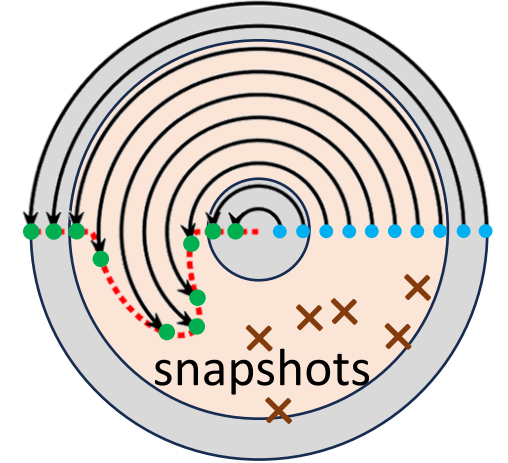
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$$\lim_{n \rightarrow \infty} \Gamma_n(\widetilde{F}_1) = \text{Sp}(\mathcal{K}_{\widetilde{F}_1}) \Rightarrow \exists n_1 \text{ s.t. } \text{dist}(i, \Gamma_{n_1}(\widetilde{F}_1)) \leq 1.$$

BUT Γ_{n_1} uses finite amount of info to output $\Gamma_{n_1}(\widetilde{F}_1)$.

Let X, Y correspond to these snapshots.



$$\mathcal{T}_F = \{(x, y_m) \mid \|F(x) - y_m\| \leq 2^{-m}\}$$

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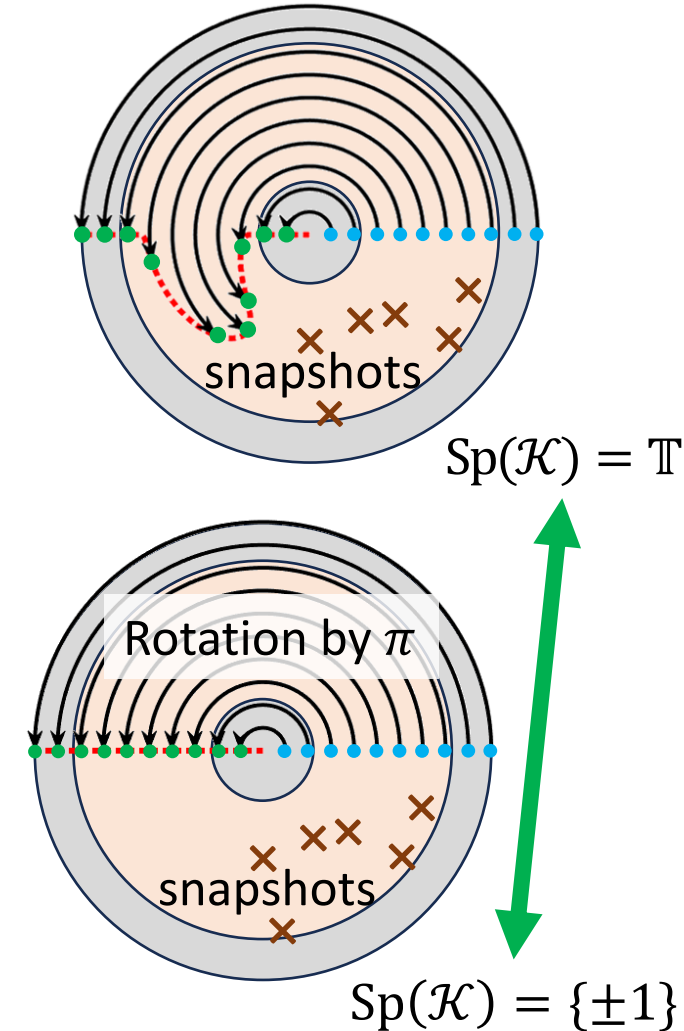
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Let X, Y correspond to these snapshots.

Lemma: $F_1 = H_1^{-1} \circ F_0 \circ H_1$ on annulus $\mathcal{A}_1$.

Consistent data $\Rightarrow \Gamma_{n_1}(F_1) = \Gamma_{n_1}(\widetilde{F}_1)$, $\text{dist}(i, \Gamma_{n_1}(F_1)) \leq 1$

BUT $\text{Sp}(\mathcal{K}_{F_1}) = \text{Sp}(\mathcal{K}_{F_0}) = \{\pm 1\}$



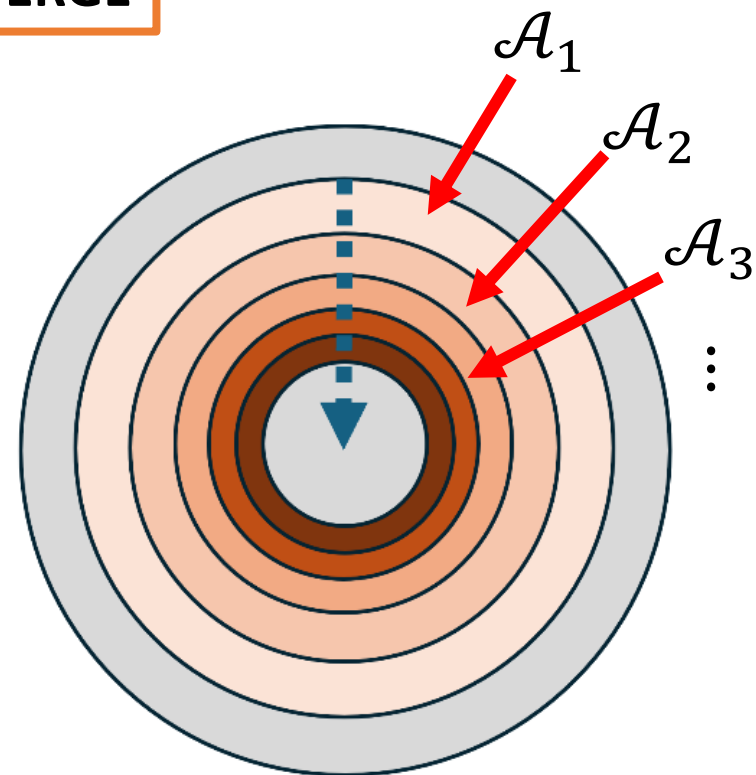
Proof idea: Constructing an adversary

Inductive step: Repeat on annuli, $F_k = H_k^{-1} \circ F_0 \circ H_k$ on $\mathcal{A}_k$. $F = \lim_{k \rightarrow \infty} F_k$

Consistent data $\Rightarrow \Gamma_{n_k}(F) = \Gamma_{n_k}(\widetilde{F_k})$, $\text{dist}(i, \Gamma_{n_k}(F)) \leq 1$, $n_k \rightarrow \infty$

BUT $\text{Sp}(\mathcal{K}_F) = \text{Sp}(\mathcal{K}_{F_0}) = \{\pm 1\}$

CANNOT CONVERGE



Cascade of disks

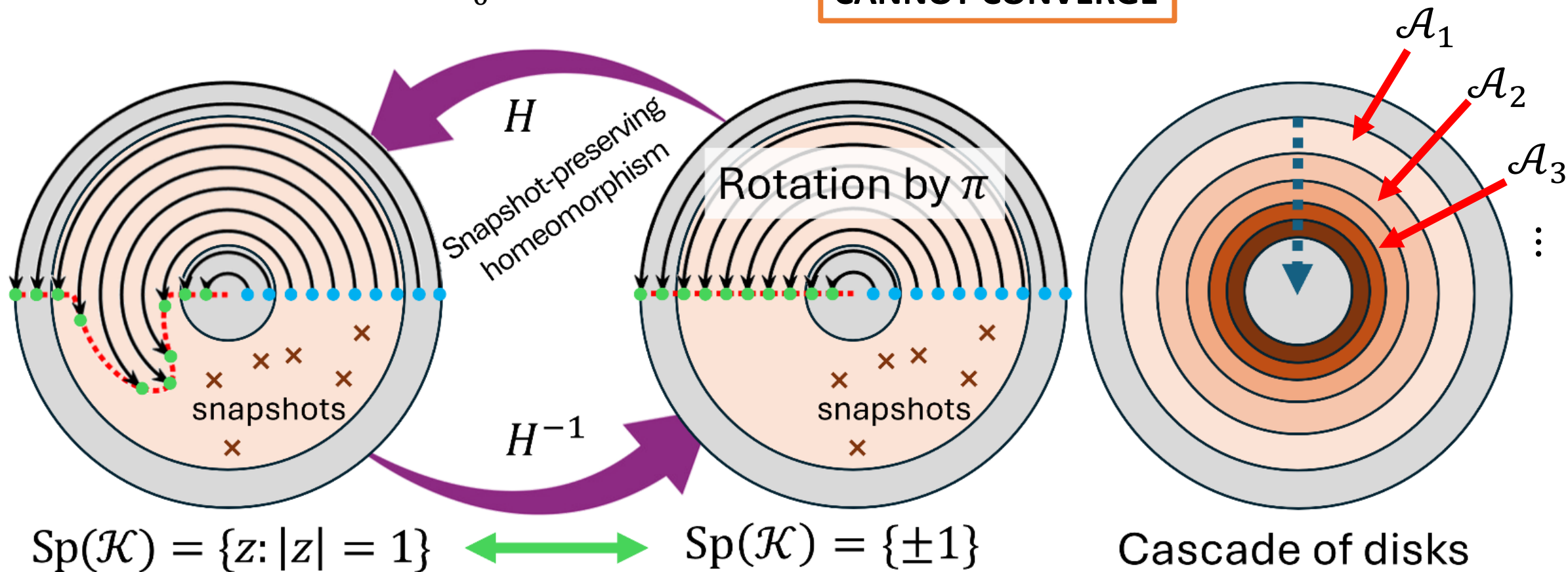
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BUT $\text{Sp}(\mathcal{K}_F) = \text{Sp}(\mathcal{K}_{F_0}) = \{\pm 1\}$

CANNOT CONVERGE



Johann Wolfgang von Goethe:

“Mathematicians are like Frenchmen: whatever you say to them they translate into their own language and forthwith it is something entirely different.”

Let's change the space...

Reproducing kernel Hilbert space (RKHS)

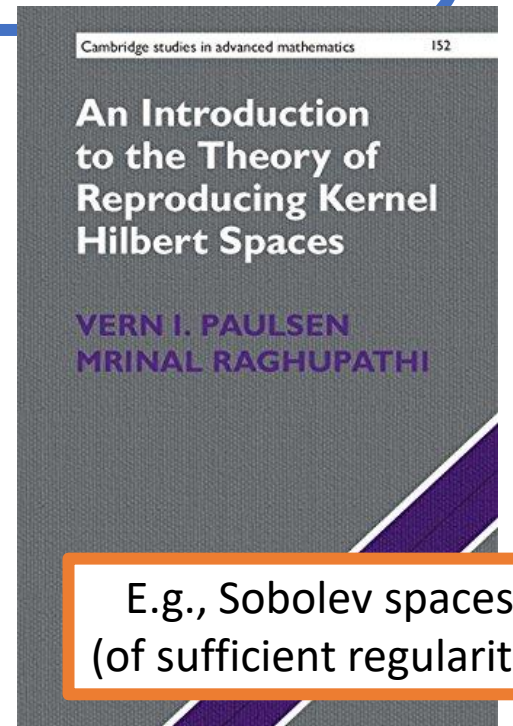
Hilbert space of functions on $\mathcal{X}$ s.t. $g \mapsto g(x)$ bounded $\forall x \in \mathcal{X}$.

Generated by a kernel $\mathfrak{K}: \mathcal{X} \times \mathcal{X} \rightarrow \mathbb{C}$

$$g(x) = \langle g, \mathfrak{K}_x \rangle, \quad \mathfrak{K}(x, y) = \langle \mathfrak{K}_x, \mathfrak{K}_y \rangle = \mathfrak{K}_x(y)$$

Advantages over $L^2(\mathcal{X}, \omega)$:

- Forecasts: space bounds $\Rightarrow$ pointwise bounds.
- High-dimensional systems practical through kernel trick.
- Fast methods for evaluating $\mathfrak{K}$.
- Different $\mathfrak{K} \Rightarrow$ different $\mathcal{K}$! Can be tailored to application.
(This is where the community is currently heading.)
- Couples M and N ...



SpecRKHS: Avoiding large data limit $M \rightarrow \infty$

Look at “Left eigenpairs” through $\mathcal{K}^*$:

$$\mathcal{K}^* \mathfrak{K}_x = \mathfrak{K}_{F(x)}$$

Evolution of functionals.

$$g(x) = \langle g, \mathfrak{K}_x \rangle_{\mathcal{H}}$$

No quadrature needed:

$$G_{jk} = \langle \mathfrak{K}_{x^{(k)}}, \mathfrak{K}_{x^{(j)}} \rangle = \mathfrak{K}(x^{(k)}, x^{(j)})$$

$$A_{jk} = \langle \mathcal{K}^* \mathfrak{K}_{x^{(k)}}, \mathfrak{K}_{x^{(j)}} \rangle = \langle \mathfrak{K}_{y^{(k)}}, \mathfrak{K}_{x^{(j)}} \rangle = \mathfrak{K}(y^{(k)}, x^{(j)})$$

$$L_{jk} = \langle \mathcal{K}^* \mathfrak{K}_{x^{(k)}}, \mathcal{K}^* \mathfrak{K}_{x^{(j)}} \rangle = \langle \mathfrak{K}_{y^{(k)}}, \mathfrak{K}_{y^{(j)}} \rangle = \mathfrak{K}(y^{(k)}, y^{(j)})$$

$$g = \sum_{m=1}^M \mathbf{g}_m \mathfrak{K}_{x^{(m)}}, \quad \|\mathcal{K}^* g - \lambda g\|_{\mathcal{H}}^2 = \mathbf{g}^* (L - \lambda A^* - \bar{\lambda} A + G) \mathbf{g}$$

SpecRKHS: Example algorithm

$$\text{res}^*(\lambda, \mathbf{g})^2 = \frac{\|\mathcal{K}^* g - \lambda g\|_{\mathcal{H}}^2}{\|g\|_{\mathcal{H}}^2} = \frac{\mathbf{g}^* [L - \lambda A^* - \bar{\lambda} A + |\lambda|^2 G] \mathbf{g}}{\mathbf{g}^* G \mathbf{g}}$$

1. Compute $G, A, L \in \mathbb{C}^{N \times N}$ ($N = M$)
2. For z_k in grid, compute $\tau_k = \min_{g = \sum_{m=1}^N \mathbf{g}_m \mathfrak{K}_{x(m)}}$ $\text{res}^*(z_k, \mathbf{g})$, corresponding g_k (gen. SVD).
3. **Output:** $\{z_k: \tau_k < \varepsilon\}, \{g_k: \tau_k < \varepsilon\}$ (ε -pseudoeigenfunctions).

Theorem:

- **Error control:** $\{z_k: \tau_k < \varepsilon\} \subseteq \text{Sp}_\varepsilon(\mathcal{K}^*)$
- **Convergence:** Converges locally uniformly to $\text{Sp}_\varepsilon(\mathcal{K}^*)$ (as $N \rightarrow \infty$)

$$\text{Sp}_\varepsilon(\mathcal{K}^*) = \{z \in \mathbb{C}: \exists g, \|g\|_{\mathcal{H}} = 1, \|\mathcal{K}^* g - z g\|_{\mathcal{H}} \leq \varepsilon\}$$

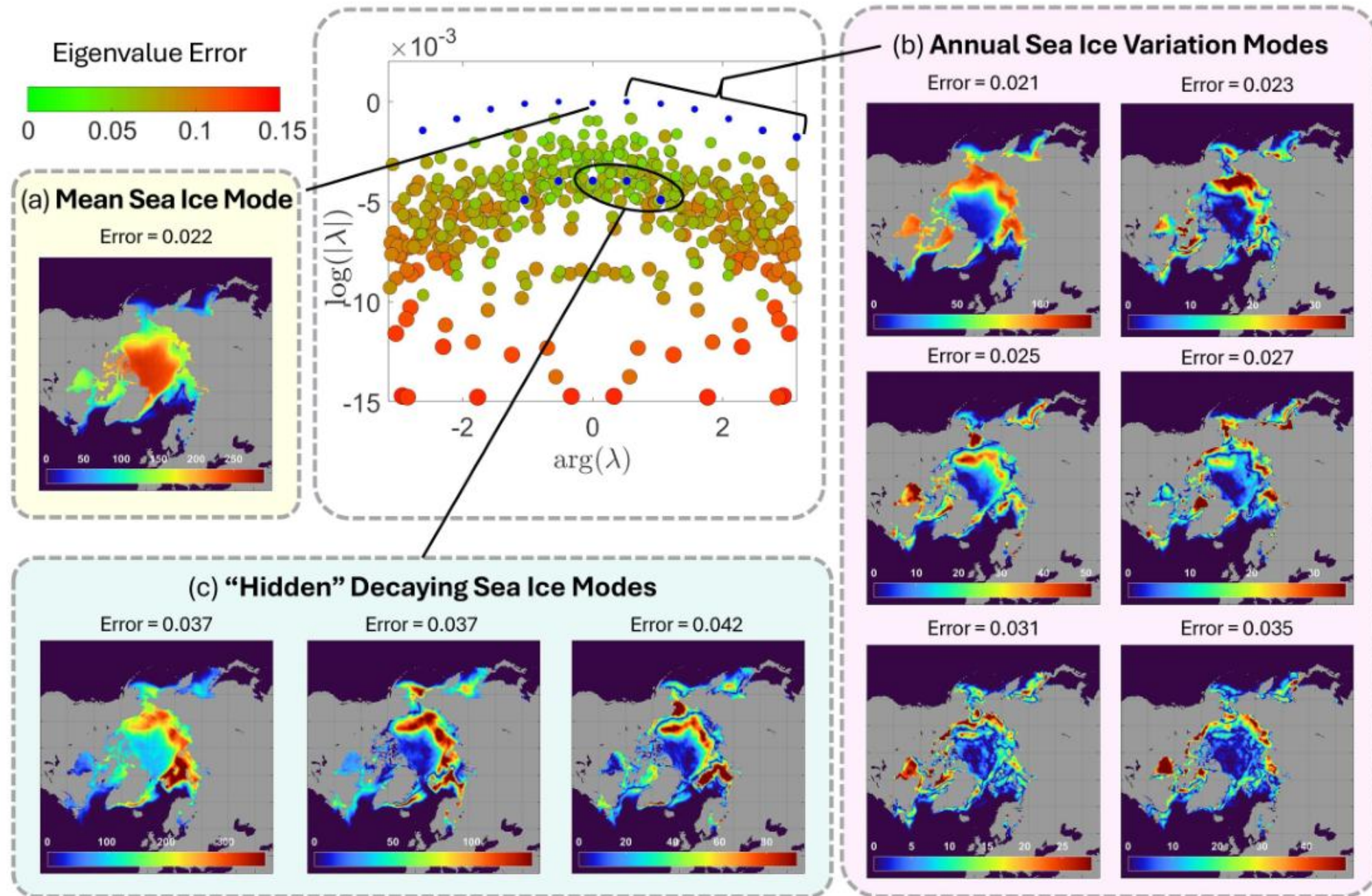
Practical gains: Arctic sea ice forecasting



Satellite data

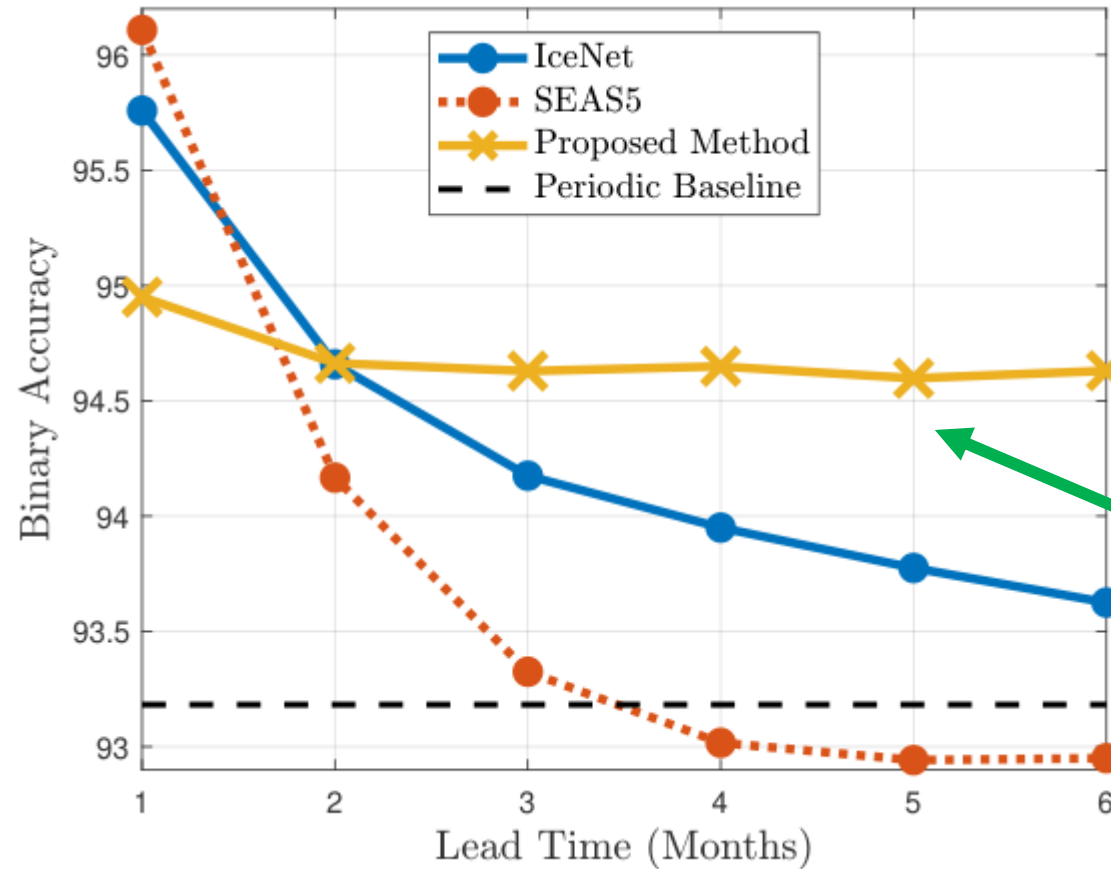


- Problems:**
1. Very hard to locate geographical significant regions.
 2. Very hard to predict more than two months in advance.



- C., Mezić, Stepanenko, “Adversarial dynamical systems characterize when data-driven learning succeeds or fails,” **Nature Com.**, 2026.

Avoid spurious evals $\Rightarrow$ State-of-the-art forecasts



$$g(x_n) = [\mathcal{K}^n g](x_0)$$

$$\|\mathcal{K}g - \lambda g\| \leq \varepsilon$$

$$\Rightarrow g(x_n) = \lambda^n g(x_0) + \mathcal{O}(n\varepsilon)$$

Use ε to filter evals!

Figure: Mean binary accuracy over test years 2012-2020.

(IceNet: Andersson et al, “Seasonal Arctic sea ice forecasting with probabilistic deep learning.” *Nature Communications*, 2021.)

What about forecast bounds and dictionary learning?

Principle Angle Decomposition (PAD)

Algorithm 3.1 Principal angles and observables between $\mathcal{V}$ and $\mathcal{KV}$.

Input: Matrices $\mathbf{G}$, $\mathbf{A}$ and $\mathbf{L}$ in (2.3) and (2.9) for dictionary $\{\psi_j\}_{j=1}^N$, subspace selection matrix $\mathbf{B} \in \mathbb{C}^{N \times n}$, cut-off tolerance $\epsilon_c \geq 0$. For RKHS, use instead the matrices $\mathbf{G}$, $\mathbf{A}$ and $\mathbf{R}$ (instead of $\mathbf{L}$) from (2.6) and (2.10).

- 1: Compress the matrices to form $\mathbf{G}_\mathcal{V} = \mathbf{B}^* \mathbf{G} \mathbf{B}$, $\mathbf{A}_\mathcal{V} = \mathbf{B}^* \mathbf{A} \mathbf{B}$, and $\mathbf{L}_\mathcal{V} = \mathbf{B}^* \mathbf{L} \mathbf{B}$.
- 2: Compute the matrix

$$\mathbf{J}_\mathcal{V} = \begin{pmatrix} \mathbf{G}_\mathcal{V} & \mathbf{A}_\mathcal{V} \\ \mathbf{A}_\mathcal{V}^* & \mathbf{L}_\mathcal{V} \end{pmatrix} \in \mathbb{C}^{2n \times 2n},$$

and its eigendecomposition $\mathbf{J}_\mathcal{V} = \mathbf{U} \mathbf{D} \mathbf{U}^*$.

- 3: Set $\mathcal{I} = \{j : \mathbf{D}_{jj} > \epsilon_c, j = 1, \dots, 2n\}$, and compute

$$\mathbf{C}_1 = \sqrt{\mathbf{D}(\mathcal{I}, \mathcal{I})} (\mathbf{U}(1:n, \mathcal{I}))^*, \quad \mathbf{C}_2 = \sqrt{\mathbf{D}(\mathcal{I}, \mathcal{I})} (\mathbf{U}(n+1:2n, \mathcal{I}))^*.$$

- 4: Compute the principal angles and observables

$$[\{\theta_j\}_{j=1}^q, \mathbf{U}_1, \mathbf{U}_2] = \text{subspacea}(\mathbf{C}_1, \mathbf{C}_2), \quad q \leq n,$$

and convert back to the original dictionary via (3.1) and (3.2).

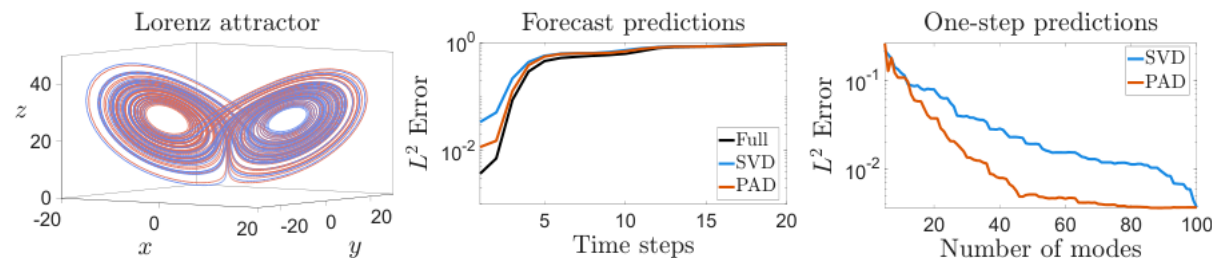
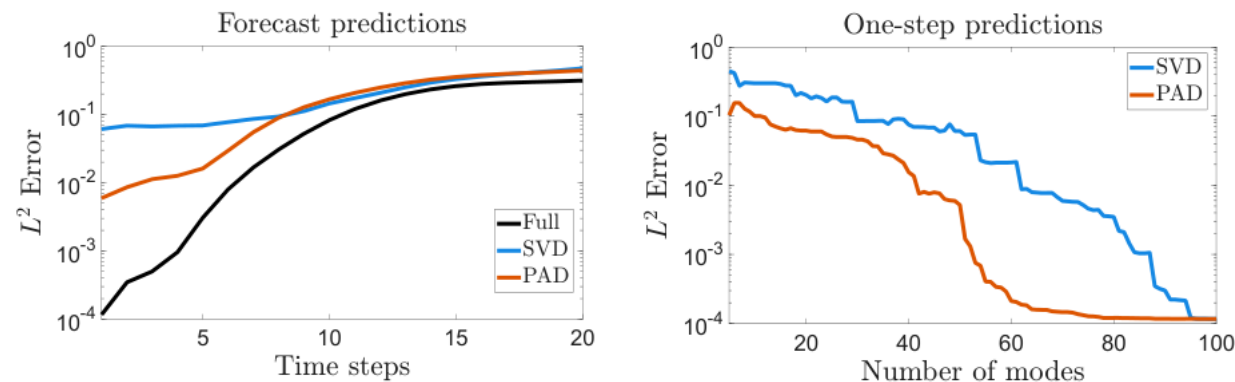
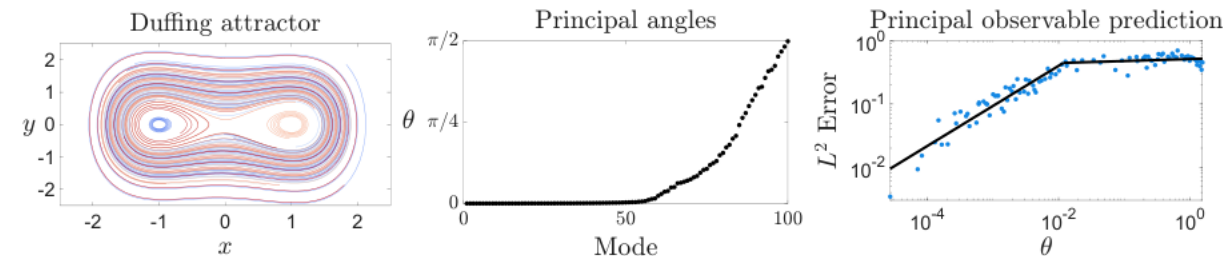
Output: Principal angles $\{\theta_j\}_{j=1}^q$ and principal observables $\{(u_j, v_j)\}_{j=1}^q$.

Algorithm 3.2 Principal Angle Decomposition (PAD).

Input: Matrices $\mathbf{G}$, $\mathbf{A}$, $\mathbf{K}$ and $\mathbf{L}$ in (2.3), (2.4), and (2.9) for dictionary $\{\psi_j\}_{j=1}^N$, cut-off tolerance $\epsilon_c \geq 0$, number of post-compression modes $r > 0$.

- 1: Compute principal angles $\{\theta_j\}_{j=1}^q$ and observables $\mathbf{U}_1$ via Algorithm 3.1 ($\mathbf{B} = \mathbf{I}_N$).
- 2: Define $\mathbf{U}'_1 = \mathbf{U}_1(1:q, :) + \mathbf{K} \mathbf{U}_1(q+1:2q, :)$, and truncate to $\mathbf{U} = \mathbf{U}'_1(:, 1:r)$.
- 3: Construct the matrices $\mathbf{A}_{\text{pad}} = \mathbf{U}^* \mathbf{A} \mathbf{U} = \mathbf{K}_{\text{pad}}$ and $\mathbf{L}_{\text{pad}} = \mathbf{U}^* \mathbf{L} \mathbf{U}$.

Output: The compressed matrices $\mathbf{A}_{\text{pad}}$, $\mathbf{K}_{\text{pad}}$ and $\mathbf{L}_{\text{pad}}$.



Principle Angle Decomposition (PAD)

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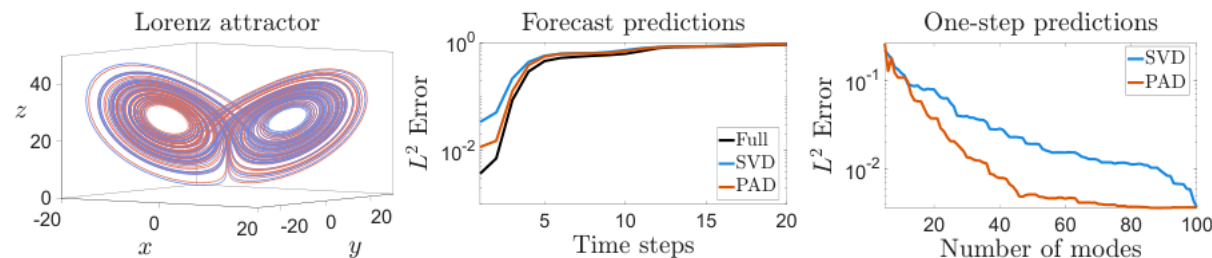
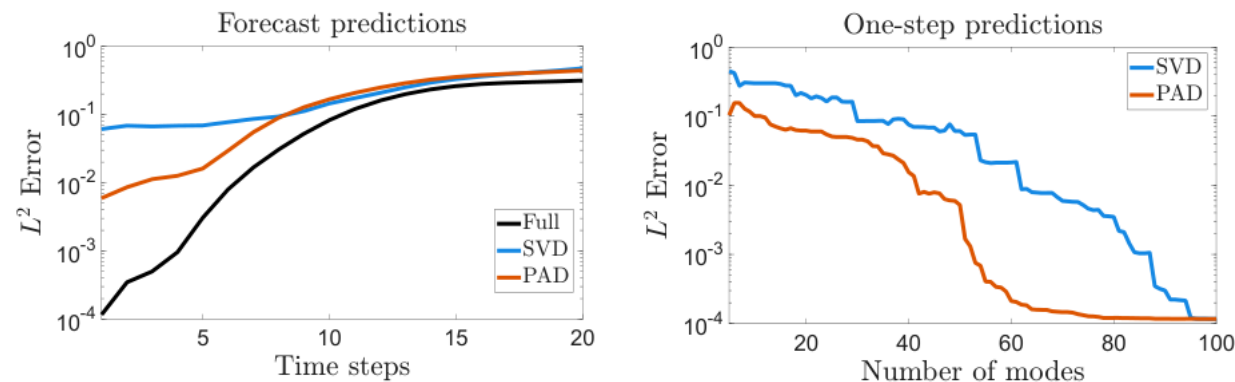
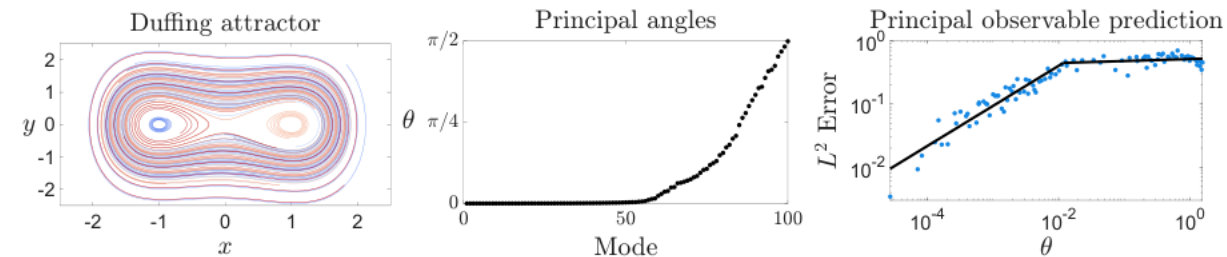
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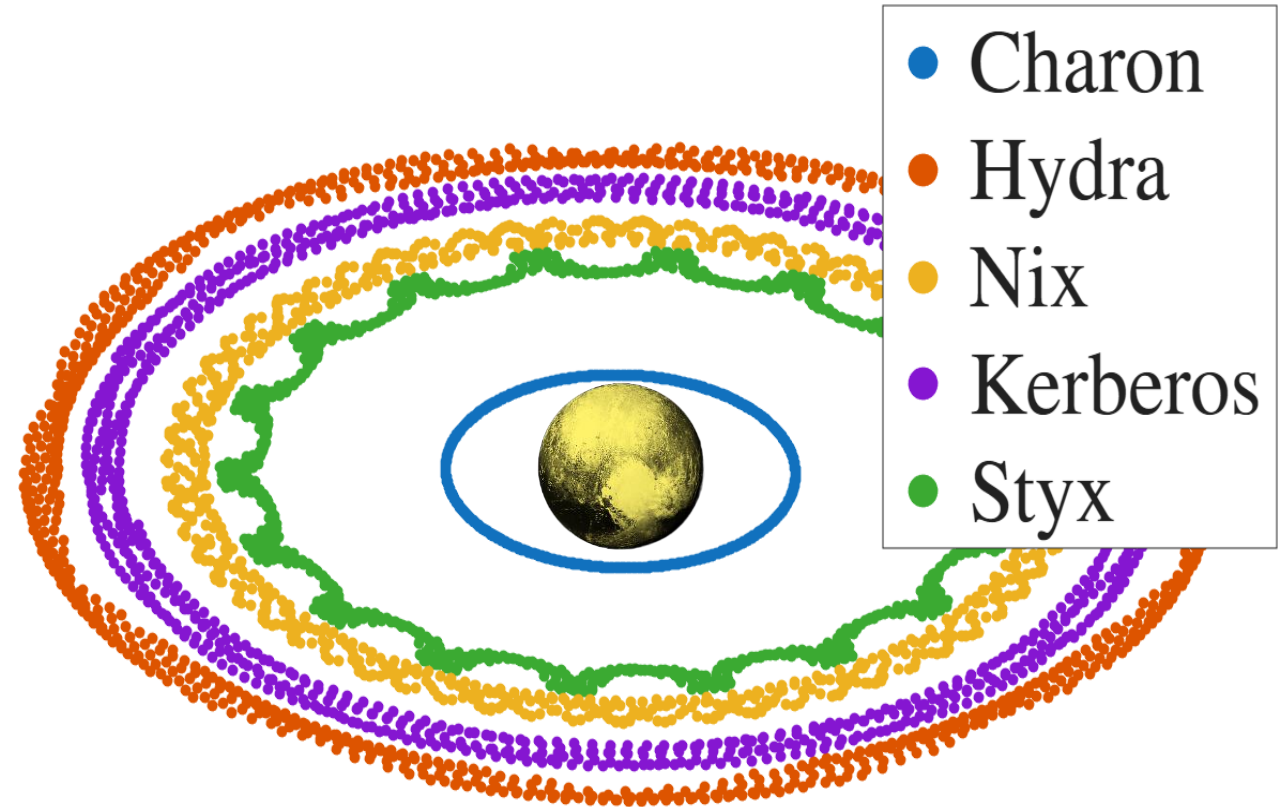


Dictionary learning and forecast multistep bounds

- Pluto and its 5 satellites
- Comparable mass of Pluto and Charon → **complex dynamics**
- Obtain error bounds for both L^2 and RKHS cases

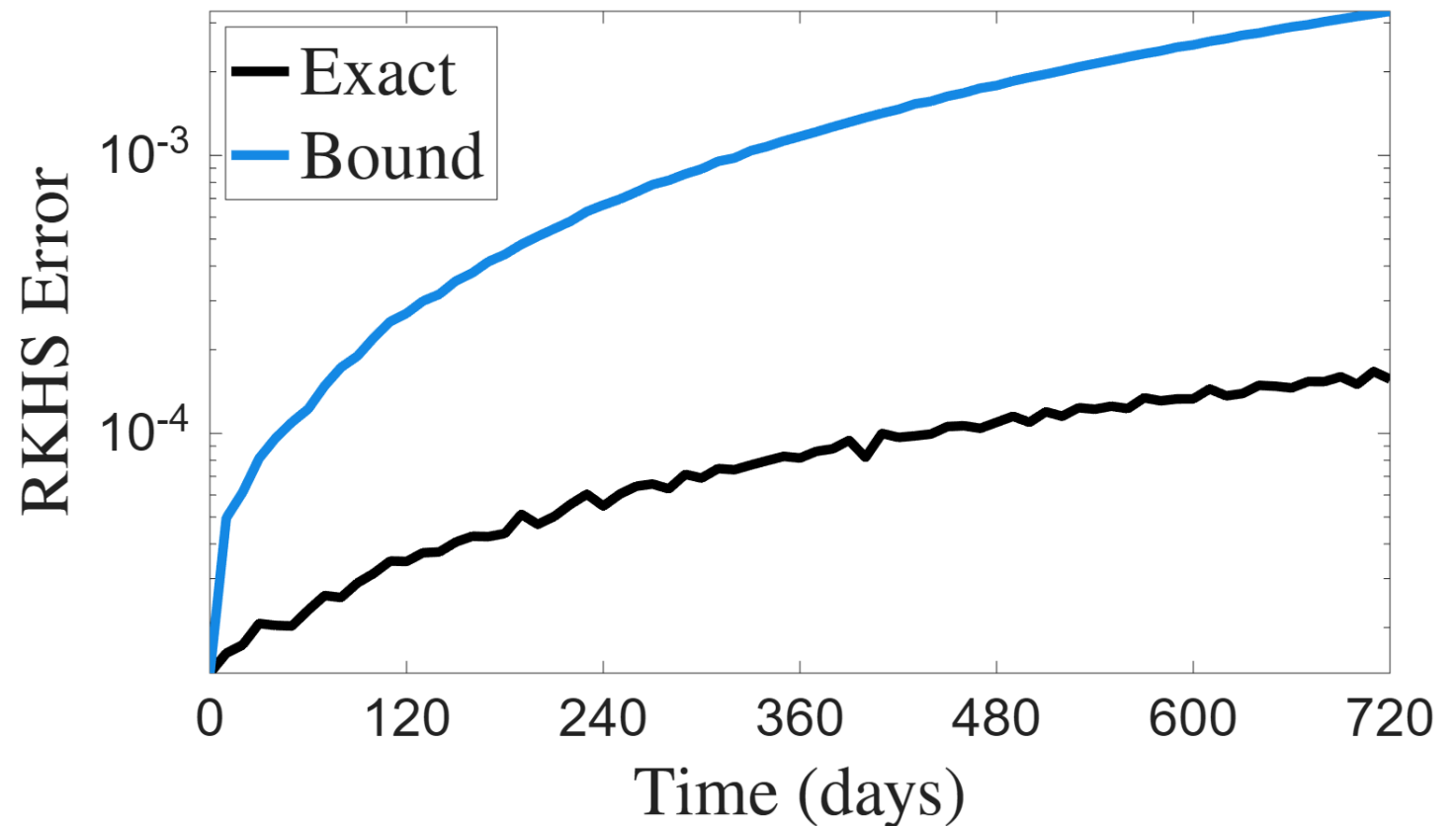
apply to dictionary refinement

obtain pointwise error bounds, expected error surrogates



Dictionary learning and forecast bounds

- Apply algorithms to RKHS setting using Matérn kernel

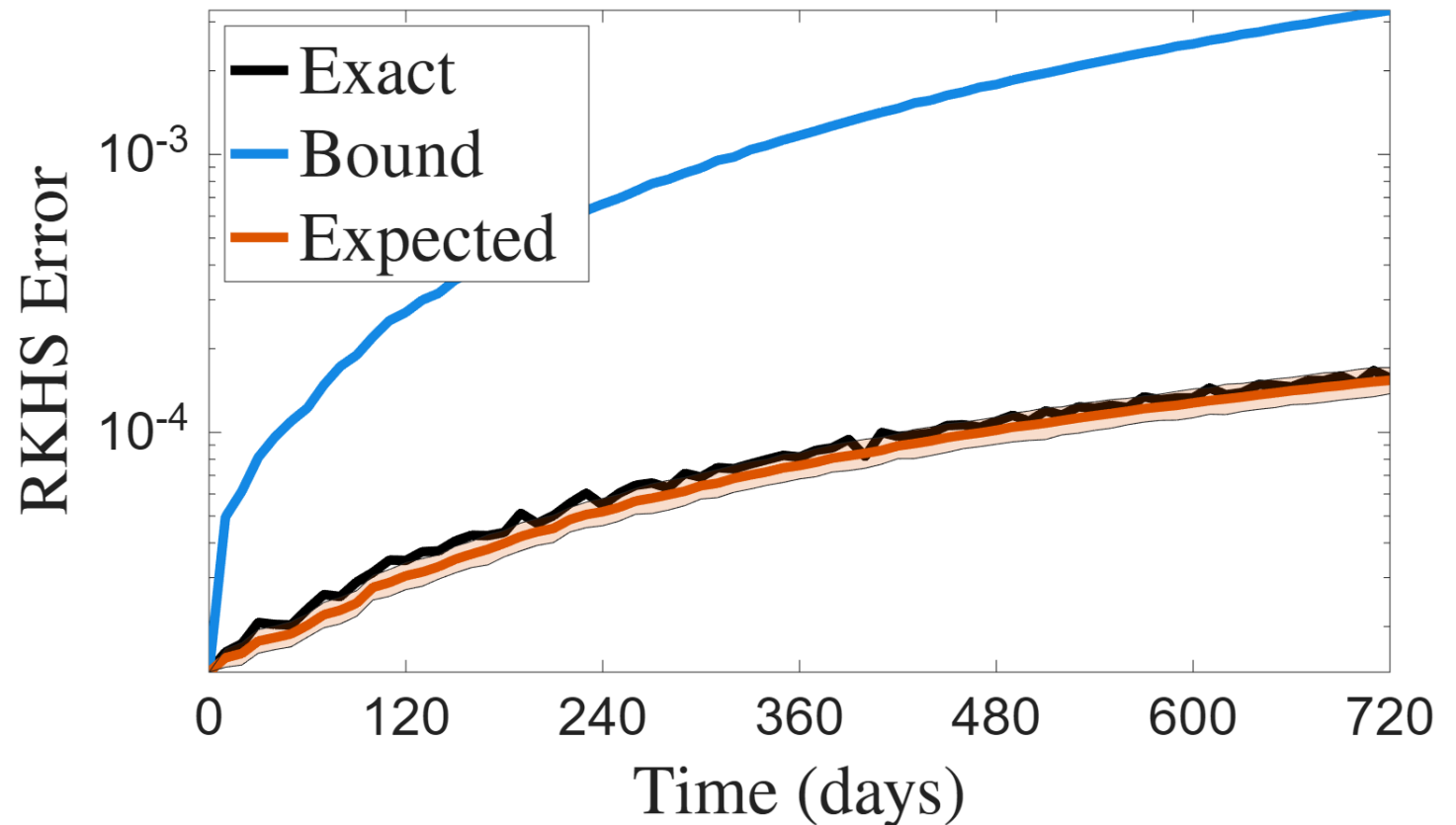


Dictionary learning and forecast bounds

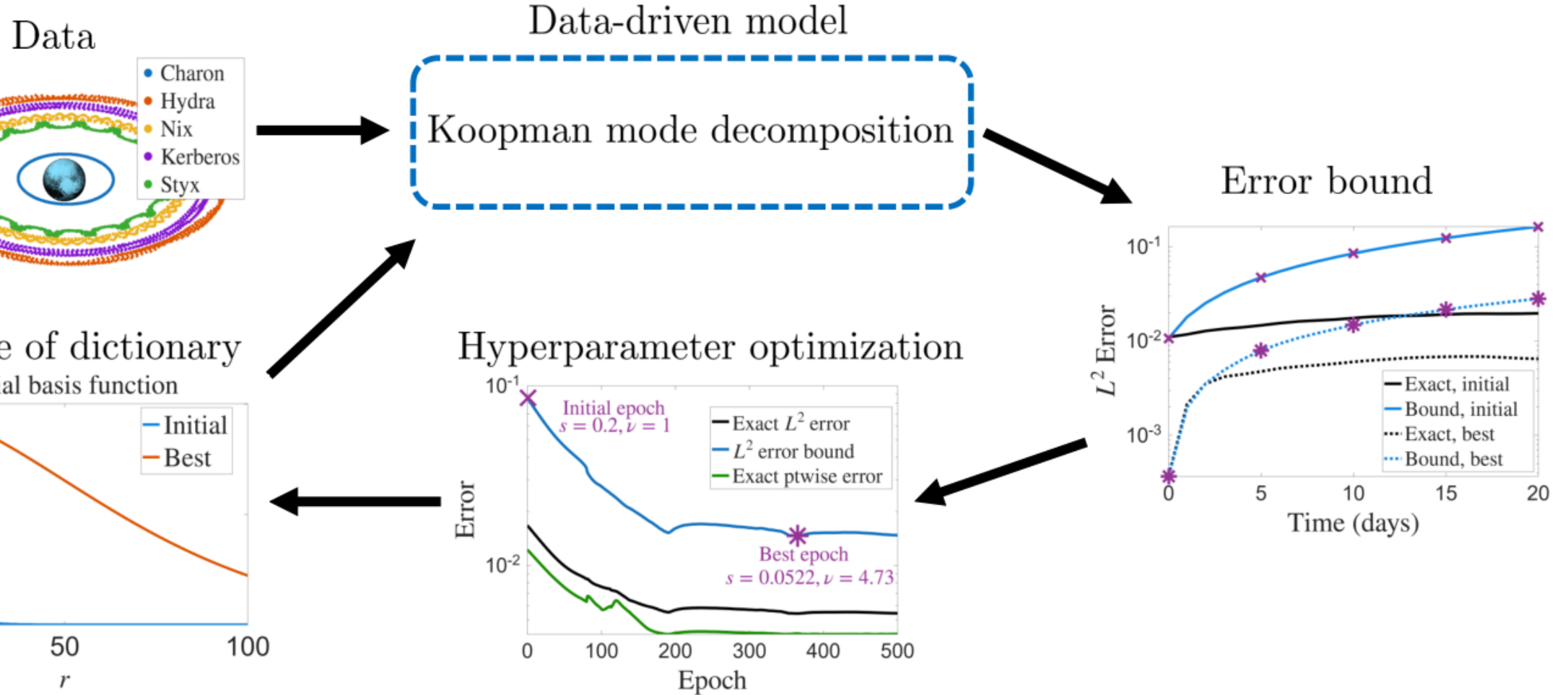
- Apply algorithms to RKHS setting using Matérn kernel
- Expected errors using **Gaussian processes** avoid overestimation

- E.g., replace $\|Ax\| \leq \|A\|\|x\|$ with $\|Ax\| \approx \mathbb{E}_{\mathcal{C}}[A]\|x\|$ where

$$\mathbb{E}_{\mathcal{C}}[A] = \mathbb{E} \left[\frac{\|Ax\|}{\|x\|} \right], \quad x \sim \mathcal{GP}(0, \mathcal{C})$$



Dictionary learning and forecast bounds

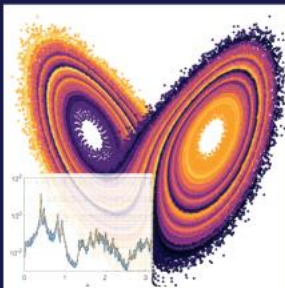
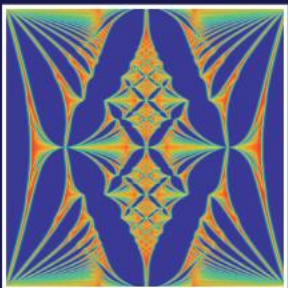
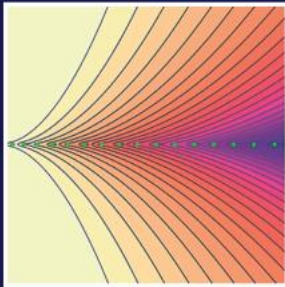
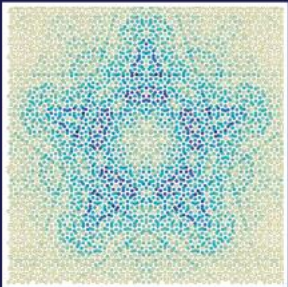


Shameless plug 1: CUP book out August 2026...

MATTHEW J. COLBROOK

Infinite-Dimensional Spectral Computations

Foundations, Algorithms, and Modern Applications



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100s of:

- Classifications
- Algorithms
- Examples (full code)
- Exercises (full solutions)

If something of interest – speak to me!

This talk

Shameless plug 2: Issac Newton Institute Programme

43

Operator Methods for Dynamical Systems

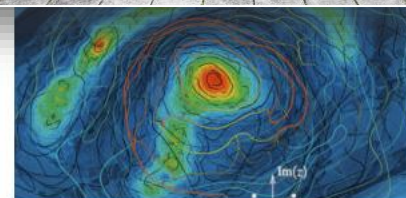
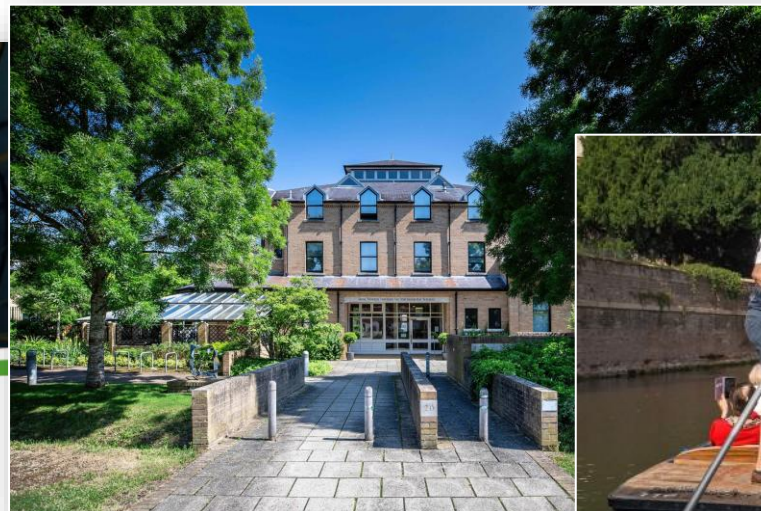
OMD

3 August 2026 to 28 August 2026

Programme theme

Dynamical systems lie at the heart of our understanding of complex phenomena, whether in modelling weather and ocean currents, molecular dynamics, population growth or stock market fluctuations. However, the nonlinear nature of systems poses a challenge to describing their behaviour. Operator methods offer a powerful way to tackle this challenge. By representing a finite-dimensional nonlinear system's evolution as a linear operator acting on an infinite-dimensional space of functions, tools from linear algebra and spectral analysis can be used to gain insights into the system's long-term behaviour. This operator-based perspective has roots in 20th-century mathematics. It has proven fruitful in classical settings like statistical mechanics, where it helped connect chaotic systems with well-understood linear techniques.

In recent years, interest in operator methods for dynamical systems has surged, driven by data-driven techniques to approximate and analyse these operators from real-world or simulated data. A wide range of frameworks and algorithms have emerged, creating an exciting opportunity and need to develop a unifying foundation for these approaches. This programme aims to bring together experts and young researchers from various communities who use operator-theoretic perspectives (for example, the Koopman and transfer operator frameworks) to study dynamics. By uniting participants from pure theory to practical applications, the programme will spark new collaborations and jointly tackle key open questions, building a more cohesive research community. In particular, a significant focus will be on exploring the spectral properties of these operators—essentially, understanding their eigenvalues, modes, and related features—as these provide crucial insights into a system's long-term behaviour and how such behaviour can be effectively analysed and computed.



Organisers

- Matthew Colbrook *University of Cambridge*
- Gary Froyland *University of New South Wales*
- Nathan Kutz *University of Washington*
- Julia Slipantschuk *University of Warwick; Universität Bayreuth*
- Caroline Wormell *University of Sydney*

Participants

- Wael Bahsoun *Loughborough University*
- Steve Brunton *University of Washington*
- Christopher Budd *University of Bath; Institute of Mathematics and its Applications*

Visit to find out more
about operator theory
and machine learning in
dynamical systems!

(and to enjoy Cambridge
in the summer!)

Pointers

1. Data-driven spectral problems for Koopman operators are hugely popular.
BUT: Standard truncation methods can fail – NEED TO GO INF-DIM!
2. **General methods with convergence for spectral properties**
 (spectra, pseudospectra, spectral measures, etc.) of K. operators!
E.g., Verification of approximate eigenfunctions leads to practical gains.
3. **SCI hierarchy** classifies computational problems:
Lower bounds through method of adversarial dynamics.
Upper bounds $\Rightarrow$ new “inf.-dim.” algorithms. Rigorous, optimal, practical.
 $\rightarrow$ We now have a near complete spectral picture for K. on $L^2(\mathcal{X}, \omega)$ and RKHS!
4. **Matrix L** leads to forecast bounds, ways to train dictionary, and PAD.

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