

Spectral Learning for Dynamical Systems

Matthew Colbrook 11th Sep 2025



"To <u>classify</u> is to bring order into chaos." - **George Pólya**

What is a Koopman operator?

- $\mathcal{X} \subset \mathbb{R}^d$ the state space
- $X \ni x$ the state

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

Henri Poincaré (Sorbonne)



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cts
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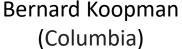
- Function space $L^2 = L^2(\mathcal{X}, \omega)$ (elements g called "observables")
- Koopman operator \mathcal{K}_F : $[\mathcal{K}_F g](x) = g(F(x))$

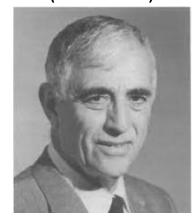




NB: Pointwise definition of \mathcal{K}_F needs $F\#\omega \ll \omega$ – this will hold throughout.

NB: \mathcal{K}_F bounded equivalent to $\mathrm{d}F\#\omega/\mathrm{d}\omega\in L^\infty$ – this will hold throughout (can be dropped).





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?

- $\mathcal{X} \subset \mathbb{R}^d$ the state space
- $X \ni x$ the state
- <u>Unknown</u> cts $F: \mathcal{X} \to \mathcal{X}$ the dynamics: $x_{n+1} = F(x_n)$
- Function space $L^2 = L^2(\mathcal{X}, \omega)$ (elements g called "observables")
- Koopman operator \mathcal{K}_F : $[\mathcal{K}_F g](x) = g(F(x))$ LINEAR!
- <u>Available</u> snapshot data: $\{(x^{(m)}, y^{(m)} = F(x^{(m)})) : m = 1, ..., M\}$

Can we compute spectral properties from trajectory data?

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

Why?

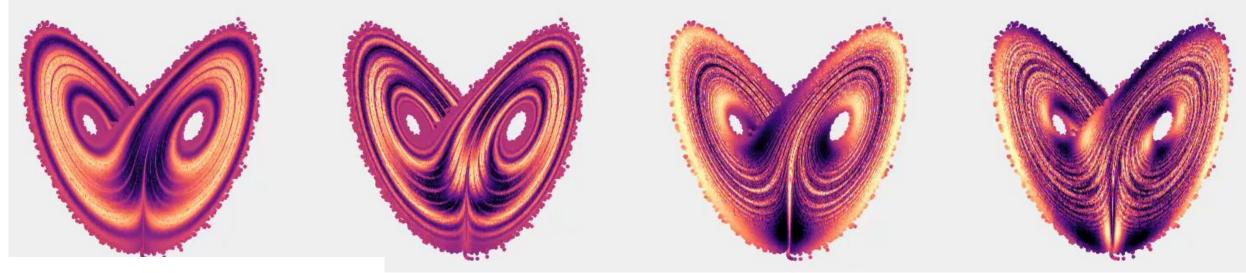
If
$$\|\mathcal{K}g - \lambda g\| \le \varepsilon$$
, then $g(x_n) = [\mathcal{K}^n g](x_0) = \lambda^n g(x_0) + \mathcal{O}(n\varepsilon)$

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

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Coherent features!

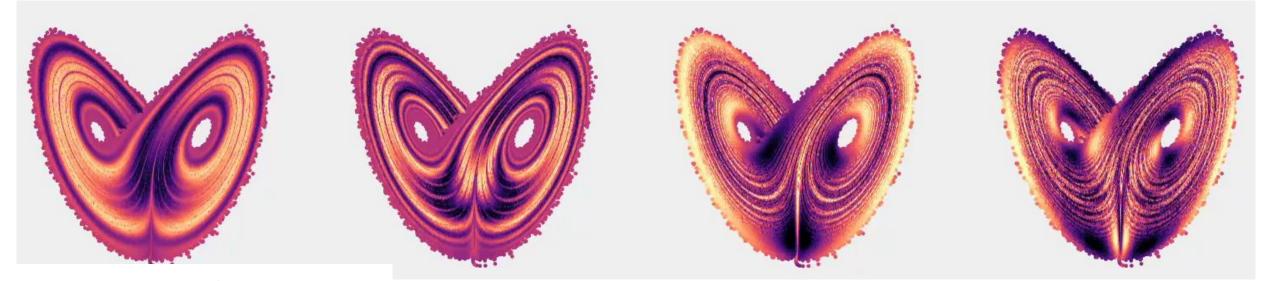
Lorenz attractor

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Coherent features!

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

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

Koopman Mode Decomposition

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

Verified Eigenfunctions

Coefficients $\in \mathbb{R}^d$, called "Koopman modes"

$$x \approx \sum_{j} c_{j} g_{j}(x) \in \mathbb{R}^{d}$$

Forecasts:

$$x_n = \sum_j \lambda_j^n c_j g_j(x_0) + \mathcal{O}(n\varepsilon)$$

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

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

GOAL: Compute spectral properties and figure out how hard this is.

DATA + DISCRETIZE ${\cal K}$

DATA + DISCRETIZE X



FINITE-DIMENSIONAL NUMERICAL LINEAR ALGEBRA

DATA + DISCRETIZE X



FINITE-DIMENSIONAL NUMERICAL LINEAR ALGEBRA



EIGENVALUES etc.

DATA + DISCRETIZE \mathcal{K}



FINITE-DIMEN NUMERICAL LINE

Works great if you have a selfadjoint operator that is compact or has compact resolvent!

Eigenvalue Problems

I. Babuška*

Institute for Physical Science and Technology and Department of Mathematics University of Maryland College Park, MD 20742, USA

J. Osborn**

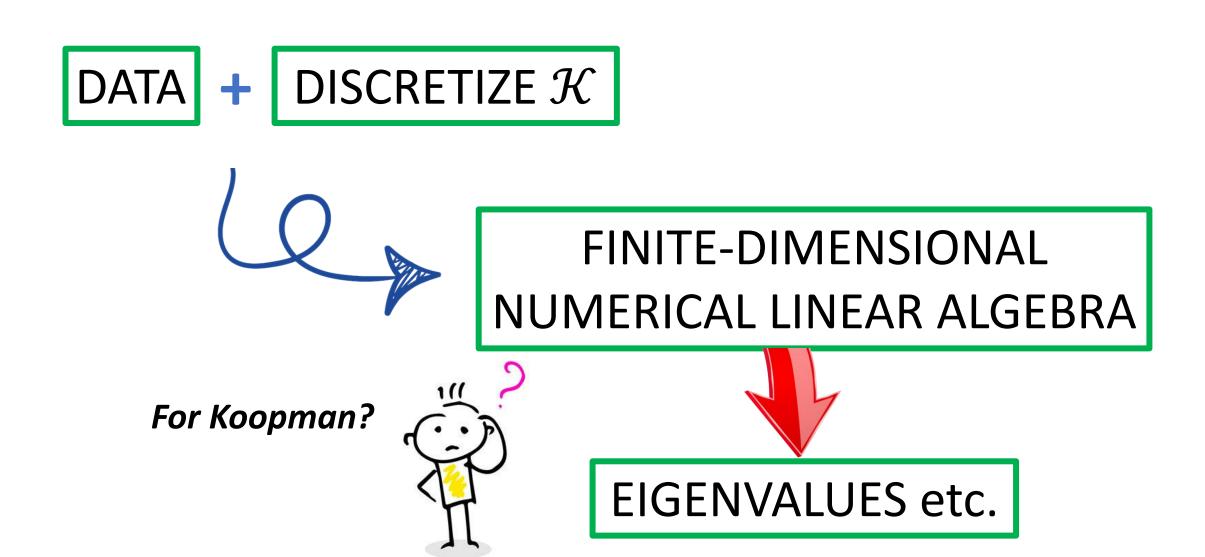
Department of Mathematics University of Maryland College Park, MD 20742, USA

*Partially supported by the Office of Naval Research under contract N00014-85-K-0169 and by the National Science Foundation under grant DMS-85-16191

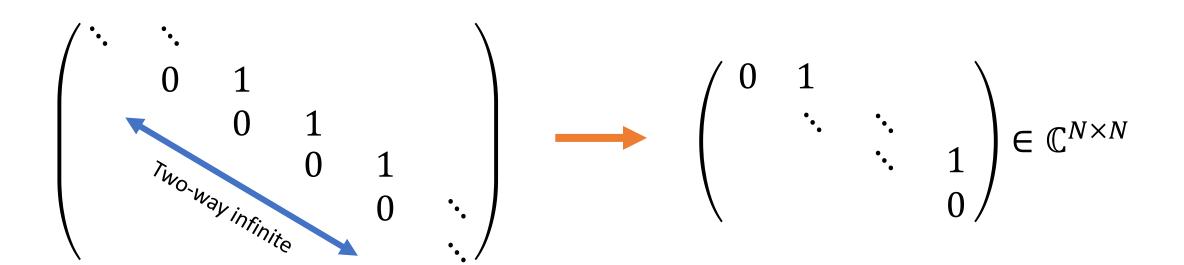
**Partially supported by the National Science Foundation under grant DMS-84-10324.

HANDBOOK OF NUMERICAL ANALYSIS, VOL. II Finite Blement Methods (Part 1) Edited by P.G. Carlet and J.L. Lions © 1991. Elsevier Science Publishers B.V. (North-Holland)

EIGENVALUES etc.



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



- 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!

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

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

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

quadrature points

$$\langle \psi_k, \psi_j \rangle \approx \sum_{m=1}^M w_m \overline{\psi_j(x^{(m)})} \psi_k(x^{(m)}) = \begin{bmatrix} \begin{pmatrix} \psi_1(x^{(1)}) & \cdots & \psi_N(x^{(1)}) \\ \vdots & \ddots & \vdots \\ \psi_1(x^{(M)}) & \cdots & \psi_N(x^{(M)}) \end{pmatrix}^* \underbrace{\begin{pmatrix} w_1 & & \\ & \ddots & \\ & & w_M \end{pmatrix}}_{\hat{W}} \underbrace{\begin{pmatrix} \psi_1(x^{(1)}) & \cdots & \psi_N(x^{(1)}) \\ \vdots & \ddots & \vdots \\ \psi_1(x^{(M)}) & \cdots & \psi_N(x^{(M)}) \end{pmatrix}}_{jk}$$
quadrature weights

$$\langle \mathcal{K}\psi_{k},\psi_{j}\rangle \approx \sum_{m=1}^{M} w_{m}\overline{\psi_{j}(x^{(m)})}\underbrace{\psi_{k}(y^{(m)})}_{[\mathcal{K}\psi_{k}](x^{(m)})} = \underbrace{\begin{bmatrix} \psi_{1}(x^{(1)}) & \cdots & \psi_{N}(x^{(1)}) \\ \vdots & \ddots & \vdots \\ \psi_{1}(x^{(M)}) & \cdots & \psi_{N}(x^{(M)}) \end{bmatrix}^{*}}_{\psi_{X}} \underbrace{\begin{pmatrix} w_{1} & & & \\ & \ddots & & \\ & & w_{M} \end{pmatrix}}_{W} \underbrace{\begin{pmatrix} \psi_{1}(y^{(1)}) & \cdots & \psi_{N}(y^{(1)}) \\ \vdots & \ddots & \vdots \\ \psi_{1}(y^{(M)}) & \cdots & \psi_{N}(y^{(M)}) \end{pmatrix}}_{ik}$$

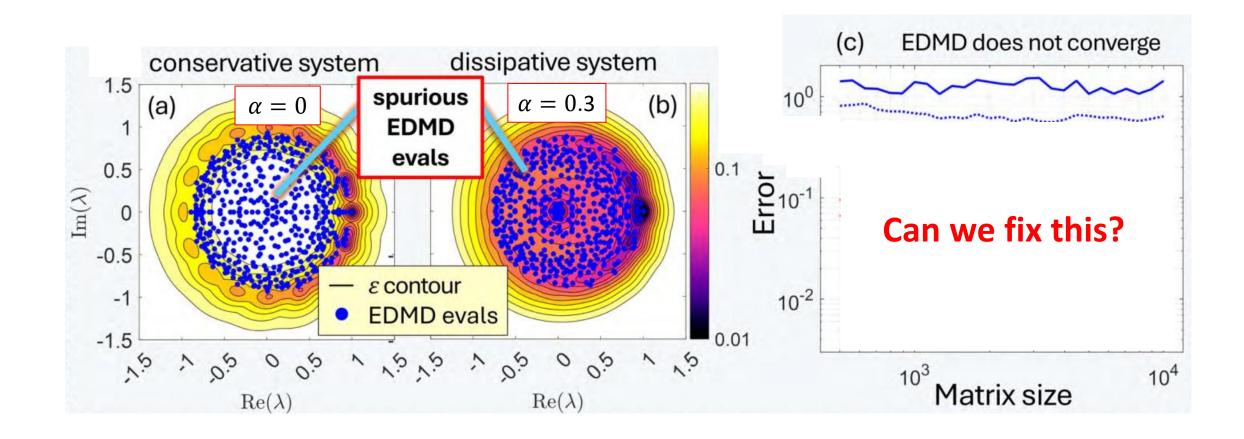
Galerkin Approximation

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

- 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)



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

$$\langle \mathcal{K} \psi_k, \psi_j \rangle \approx \sum_{m=1}^{M} w_m \overline{\psi_j(x^{(m)})} \, \underbrace{\psi_k(y^{(m)})}_{[\mathcal{K} \psi_k](x^{(m)})} = \left[\underbrace{\Psi_X^* W \Psi_Y}_{K_1} \right]_{jk}$$

- C., Townsend, "Rigorous data-driven computation of spectral properties of Koopman operators for dynamical systems," Commun. Pure Appl. Math., 2023.
- C., Ayton, Szőke, "Residual Dynamic Mode Decomposition," J. Fluid Mech., 2023.
- Code: https://github.com/MColbrook/Residual-Dynamic-Mode-Decomposition

$$\langle \psi_{k}, \psi_{j} \rangle \approx \sum_{m=1}^{M} w_{m} \overline{\psi_{j}(x^{(m)})} \psi_{k}(x^{(m)}) = \left[\underline{\Psi_{x}^{*}W\Psi_{x}} \right]_{jk}$$

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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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, $\|\mathcal{K}g - \lambda g\|^2 = \sum_{k,j=1}^{N} \mathbf{g}_k \overline{\mathbf{g}_j} \langle \mathcal{K}\psi_k - \lambda \psi_k, \mathcal{K}\psi_j - \lambda \psi_j \rangle$

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$$\langle \psi_k, \psi_j \rangle \approx \sum_{m=1}^M w_m \overline{\psi_j(x^{(m)})} \, \psi_k(x^{(m)}) = \left[\underbrace{\Psi_X^* W \Psi_X}_{\widehat{G}} \right]_{jk}$$
 Infinite-dimensional error bound!
$$\langle \mathcal{K} \psi_k, \psi_j \rangle \approx \sum_{m=1}^M w_m \overline{\psi_j(x^{(m)})} \, \underbrace{\psi_k(y^{(m)})}_{[\mathcal{K} \psi_k](x^{(m)})} = \left[\underbrace{\Psi_X^* W \Psi_Y}_{\widehat{K}_1} \right]_{jk}$$
 error bound!
$$\langle \mathcal{K} \psi_k, \mathcal{K} \psi_j \rangle \approx \sum_{m=1}^M w_m \overline{\psi_j(y^{(m)})} \, \psi_k(y^{(m)}) = \left[\underbrace{\Psi_Y^* W \Psi_Y}_{\widehat{K}_2} \right]_{jk}$$

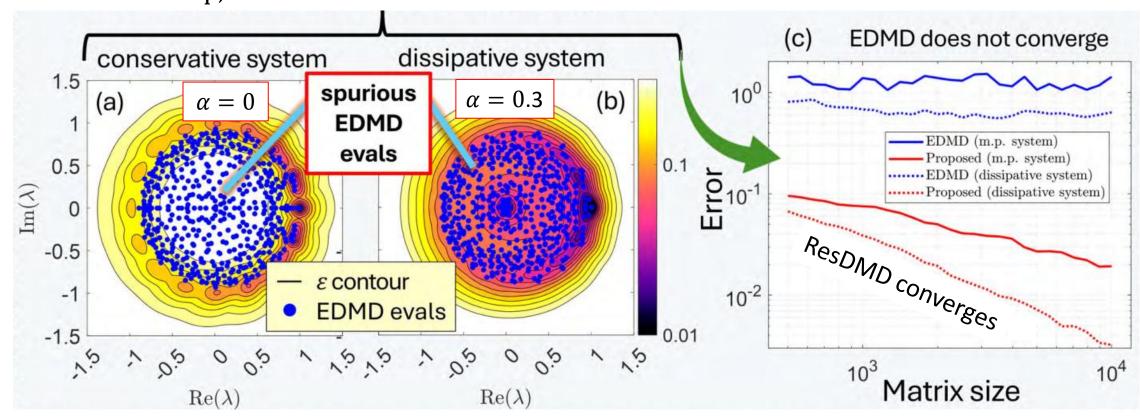
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, $\|\mathcal{K}g - \lambda g\|^{2} = \lim_{M \to \infty} \mathbf{g}^{*} [K_{2} - \lambda K_{1}^{*} - \bar{\lambda} K_{1} + |\lambda|^{2} G] \mathbf{g}$

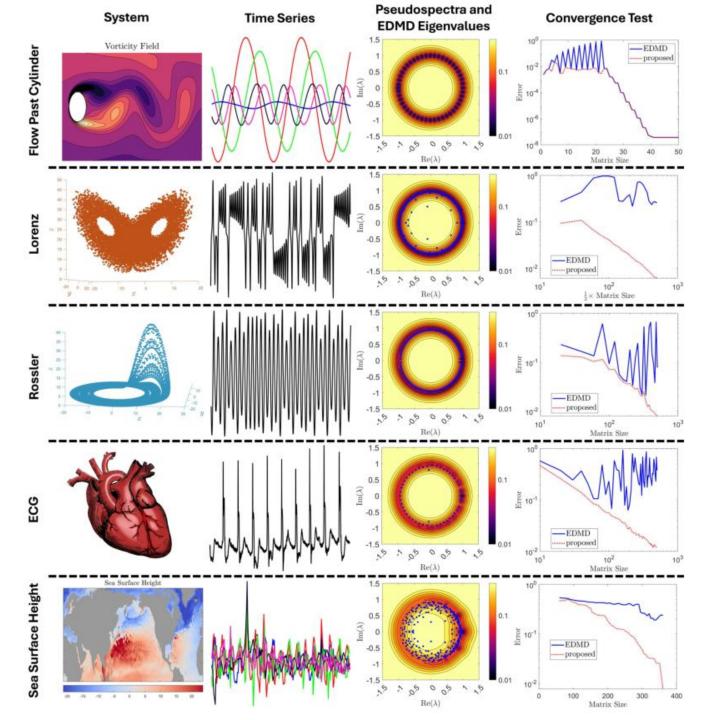
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ResDMD does converge!

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





Can maths help guide the way?

If (X, d) is a compact measure space and ω a Borel measure...

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

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

for all systems.



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

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

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

Double limit $\lim_{N\to\infty} \lim_{N\to\infty}$

Can we do better?

Adversaries: **Double** limit is necessary!

Implies ${\mathcal K}$ is unitary

Class of systems: $\Omega_{\mathbb{D}} = \{F : \overline{\mathbb{D}} \to \overline{\mathbb{D}} | F \text{ cts, measure preserving, invertible} \}$.

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

Theorem: There does not exist any sequence of deterministic algorithms $\{\Gamma_n\}$ using \mathcal{T}_F such that $\lim_{n\to\infty}\Gamma_n(F)=\operatorname{Sp}_{\mathrm{ap},\epsilon}(\mathcal{K}_F)\ \forall F\in\Omega_{\mathbb{D}}.$

NB:

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

• C., Mezić, Stepanenko, "Adversarial Dynamical Systems Reveal Limits and Rules for Trustworthy Data-Driven Learning," preprint, 2025.

$$F_0$$
: rotation by π , $\mathrm{Sp}(\mathcal{K}_{F_0})=\{\pm 1\}$

Phase transition lemma: Let $X = \{x_1, ..., x_N\}, Y = \{y_1, ..., y_N\}$ be distinct points in annulus $\mathcal{A} = \{x \in \mathbb{D} | 0 < R < \|x\| < r < 1\}$ 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 $H(y_j) = F_0(H(x_j)), j = 1, ..., N$.

Conjugacy of data $(x_i \rightarrow y_i)$ with F_0

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

• Brown and Halperin. "On certain area-preserving maps." Annals of Mathematics, 1935.

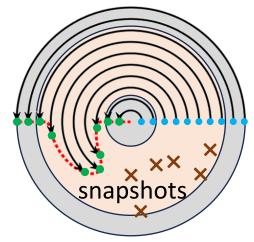
Suppose (for contradiction) $\{\Gamma_n\}$ uses \mathcal{T}_F , $\lim_{n\to\infty}\Gamma_n(F)=\operatorname{Sp}(\mathcal{K}_F)\ \forall F\in\Omega_{\mathbb{D}}$. Build an adversarial F...

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Build an adversarial F...

$$\widetilde{F_1}(r,\theta) = (r,\theta + \pi + \phi(r)), \operatorname{supp}(\phi) \subset [1/4, 3/4]$$

 $\operatorname{Sp}(\mathcal{K}_{\widetilde{F_1}}) = \mathbb{T}$ (unit circle).



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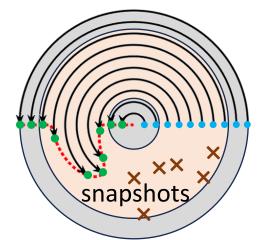
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 $\operatorname{Sp}(\mathcal{K}_{\widetilde{F_1}}) = \mathbb{T}$ (unit circle).

 $\lim_{n\to\infty}\Gamma_n\big(\widetilde{F_1}\big)=\operatorname{Sp}(\mathcal{K}_{\widetilde{F_1}})\Rightarrow \exists n_1 \text{ s.t. } \operatorname{dist}(i,\Gamma_{n_1}\big(\widetilde{F_1}\big))\leq 1.$

BUT Γ_{n_1} uses finite amount of info to output $\Gamma_{n_1}(\widetilde{F_1})$. Let X, Y correspond to these snapshots.



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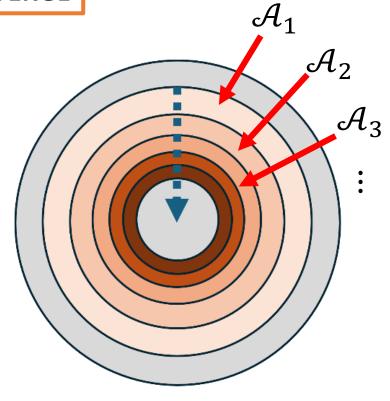
Lemma: $F_1 = H_1^{-1} \circ F_0 \circ H_1$ on annulus \mathcal{A}_1 . Consistent data $\Rightarrow \Gamma_{n_1}(F_1) = \Gamma_{n_1}(\widetilde{F_1})$, dist $(i, \Gamma_{n_1}(F_1)) \leq 1$ BUT $\operatorname{Sp}(\mathcal{K}_{F_1}) = \operatorname{Sp}(\mathcal{K}_{F_0}) = \{\pm 1\}$

snapsho^{*} $\operatorname{Sp}(\mathcal{K}) = \mathbb{T}$ Rotation by π $Sp(\mathcal{K}) = \{\pm 1\}$

Inductive step: Repeat on annuli, $F_k = H_k^{-1} \circ F_0 \circ H_k$ on \mathcal{A}_k . $F = \lim_{k \to \infty} F_k$ Consistent data $\Rightarrow \Gamma_{n_k}(F) = \Gamma_{n_k}(\widetilde{F_k})$, $\operatorname{dist}(i, \Gamma_{n_k}(F)) \leq 1$, $n_k \to \infty$

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

CANNOT CONVERGE

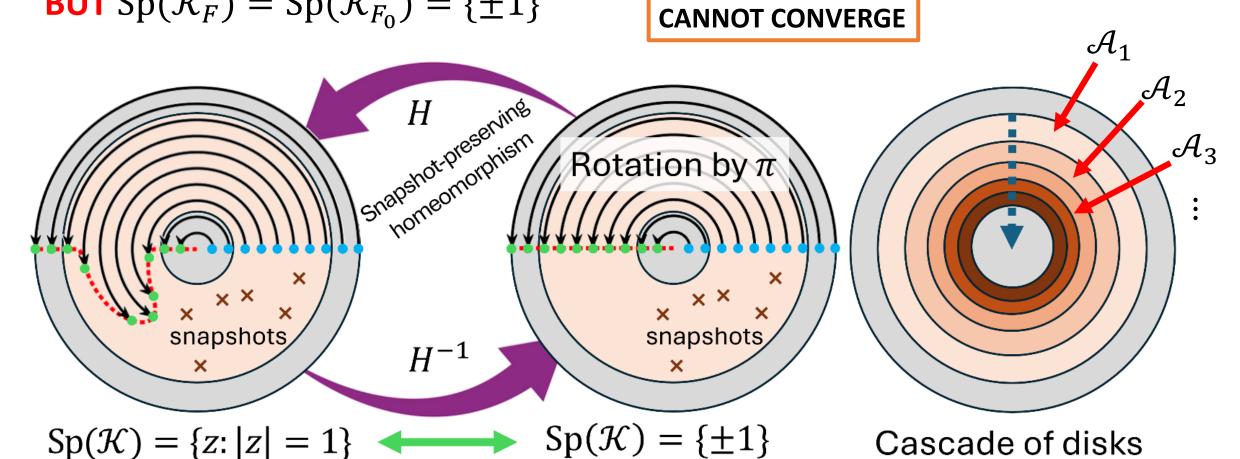


Cascade of disks

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

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

BUT Sp(\mathcal{K}_F) = Sp(\mathcal{K}_{F_0}) = {±1}



Classifications: Solvability Complexity Index (SCI)

SCI: Fewest number of limits needed to solve a computational problem.

- Δ_1 : One limit, full error control. E.g., $d(\Gamma_n(F), \operatorname{Sp}(\mathcal{K}_F)) \leq 2^{-n}$.
- Δ_{m+1} : $SCI \leq m$.
- Σ_m : SCI $\leq m$, final limit from below.

E.g.,
$$\Sigma_1$$
: $\sup_{z \in \Gamma_n(F)} \operatorname{dist}(z, \operatorname{Sp}(\mathcal{K}_F)) \leq 2^{-n}$.

• Π_m : SCI $\leq m$, final limit from above.

E.g.,
$$\Pi_1$$
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- Hansen, "On the solvability complexity index, the n-pseudospectrum and approximations of spectra of operators." J. Am. Math. Soc., 2011.
- C., "The foundations of infinite-dimensional spectral computations," PhD diss., University of Cambridge, 2020.
- C., Hansen, "The foundations of spectral computations via the solvability complexity index hierarchy," J. Eur. Math. Soc., 2022.
- C., Antun, Hansen, "The difficulty of computing stable and accurate neural networks," Proc. Natl. Acad. Sci. USA, 2022.
- Ben-Artzi, C., Hansen, Nevanlinna, Seidel, "On the solvability complexity index hierarchy and towers of algorithms," arXiv, 2020.

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SCI: Fewest number of limits needed to solve a computational problem.

- Δ_1 : One limit, full error control. E.g., $d(\Gamma_n(F), \operatorname{Sp}(\mathcal{K}_F)) \leq 2^{-n}$.
- Δ_{m+1} : $SCI \leq m$.

trust output

• Σ_m : SCI $\leq m$, final limit from below.

E.g.,
$$\Sigma_1$$
: $\sup_{z \in \Gamma_n(F)} \operatorname{dist}(z, \operatorname{Sp}(\mathcal{K}_F)) \leq 2^{-n}$. al limit from above.

• Π_m : SCI $\leq m$, final limit from above.

E.g.,
$$\Pi_1$$
: $\sup_{z \in \operatorname{Sp}(\mathcal{K}_F)} \operatorname{dist}(z, \Gamma_n(F)) \leq 2^{-n}$.

covers spectrum

- Hansen, "On the solvability complexity index, the n-pseudospectrum and approximations of spectra of operators." J. Am. Math. Soc., 2011.
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- Ben-Artzi, C., Hansen, Nevanlinna, Seidel, "On the solvability complexity index hierarchy and towers of algorithms," arXiv, 2020.

Lots of SCI upper bounds lurking in Koopman literature!

SCI: Fewest number of limits needed to solve a computational problem.

Algorithm	Comments/Assumptions	Spectral Problem's Corresponding SCI Upper Bound			
Aigontiili		KMD	Spectrum	Spectral Measure (if m.p.)	Spectral Type (if m.p.)
Extended DMD [47]	general L^2 spaces	$SCI \le 2^*$	N/C	N/C	n/a
Residual DMD [44]	general L^2 spaces	$SCI \le 2^*$	$SCI \le 3^*$	$SCI \le 2^*$	varies, see [84] e.g., a.c. density: $SCI \le 2^*$
Measure-preserving EDMD [45]	m.p. systems	$SCI \le 1$	N/C	$SCI \le 2^*$ (general) $SCI \le 1$ (delay-embedding)	n/a
Hankel DMD [85]	m.p. ergodic systems	$SCI \le 2^*$	N/C	N/C	n/a
Periodic approximations [86]	m.p. $+\omega$ a.c.	$SCI \leq 2$	N/C	$SCI \le 2$ (see [87])	a.c. density: $SCI \leq 3$
Christoffel–Darboux kernel [40]	m.p. ergodic systems	$SCI \leq 3$	n/a	$SCI \leq 2$	e.g., a.c. density: $SCI \leq 2$
Generator EDMD [88]	ctstime, samples ∇F (otherwise additional limit)	$SCI \le 2$	N/C	$SCI \leq 2$ (see [89])	n/a
Compactification [42]	ctstime, m.p. ergodic systems	$SCI \le 4$	N/C	$ SCI \le 4$	n/a
Resolvent compactification [43]	ctstime, m.p. ergodic systems	$SCI \leq 5$	N/C	$SCI \leq 5$	n/a
Diffusion maps [90] (see also [10])		•	•••••	n/a	
D			•		Are these sharp?

Previous techniques prove upper bounds on SCI.

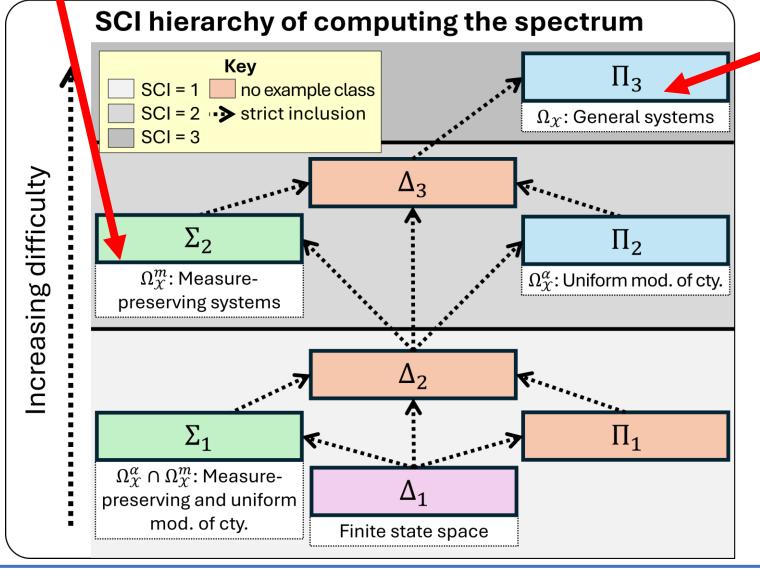
"N/C": method need not converge. "n/a": algorithm not applicable to problem.

Also in Ulam's method for Markov processes, SRB measure computation, control,...

Lower + upper bounds

Classification for Koopman

3 limits needed in general!



Different classes:

$$\Omega_{\mathcal{X}} = \{F: \mathcal{X} \to \mathcal{X} \mid F \text{ cts}\}$$

$$\Omega_{\mathcal{X}}^{m} = \{F: \mathcal{X} \to \mathcal{X} \mid F \text{ cts, m. p.}\}$$

$$\Omega_{\mathcal{X}}^{\alpha} = \{F: \mathcal{X} \to \mathcal{X} \mid F \text{ mod. cty. } \alpha\}$$

$$[d_{\mathcal{X}}(F(x), F(y)) \leq \alpha(d_{\mathcal{X}}(x, y))]$$

Optimal algorithms and classifications of dynamical systems.

C., Mezić, Stepanenko, "Adversarial Dynamical Systems Reveal Limits and Rules for Trustworthy Data-Driven Learning," preprint, 2025.

Peter Lax:

"The trick of the successful mathematician is to turn the question being asked into one he knows how to answer."

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 perform this trick by changing 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 $\Re: \mathcal{X} \times \mathcal{X} \to \mathbb{C}$

$$g(x) = \langle g, \mathfrak{K}_{\chi} \rangle, \qquad \mathfrak{K}(x, y) = \langle \mathfrak{K}_{\chi}, \mathfrak{K}_{\gamma} \rangle = \mathfrak{K}_{\chi}(y)$$

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

- Forecasts: space bounds ⇒ pointwise bounds.
- High-dimensional systems practical through kernel trick.
- Fast methods for evaluating \Re .
- Different $\Re \Rightarrow$ different $\Re!$ Can be tailored to application. (This is where the community is currently heading.)
- Leads to fundamental "possibility" gains...

An Introduction to the Theory of Reproducing Kernel Hilbert Spaces

VERN I. PAULSEN MRINAL RAGHUPATHI

E.g., Sobolev spaces (of sufficient regularity)

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An Introduction to the Theory of Reproducing Kernel Hilbert Spaces

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E.g., Sobolev spaces (of sufficient regularity)

SpecRKHS: Avoiding large was Look at "Left eigenpairs" through \mathcal{K}^* : $\mathcal{K}^*\mathfrak{K}_\chi = \mathfrak{K}_{F(\chi)}$

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

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

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

$$R_{jk} = \left\langle \mathcal{K}^* \mathfrak{K}_{\chi(k)}, \mathcal{K}^* \mathfrak{K}_{\chi(j)} \right\rangle = \left\langle \mathfrak{K}_{y(k)}, \mathfrak{K}_{y(j)} \right\rangle = \mathfrak{K}(y^{(k)}, y^{(j)})$$

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

Boullé, C., Conradie, "Convergent Methods for Koopman Operators on Reproducing Kernel Hilbert Spaces," preprint, 2025.

SpecRKHS: Example algorithm

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

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

Theorem:

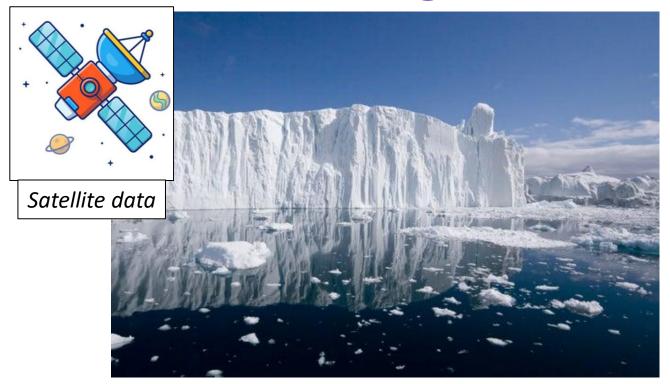
First convergent method for general ${\mathcal K}$

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

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

• Boullé, C., Conradie, "Convergent Methods for Koopman Operators on Reproducing Kernel Hilbert Spaces," preprint, 2025.

Practical gains: Sea ice forecasting

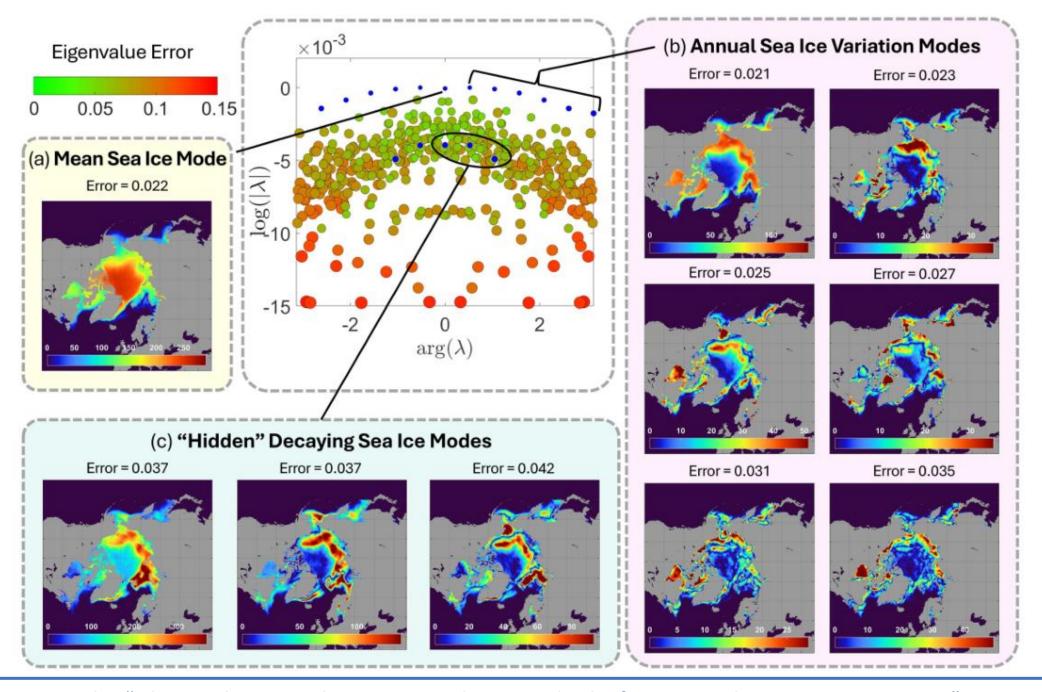




Motivation: Arctic amplification, polar bears, local communities, effect on extreme weather in Northern hemisphere,...

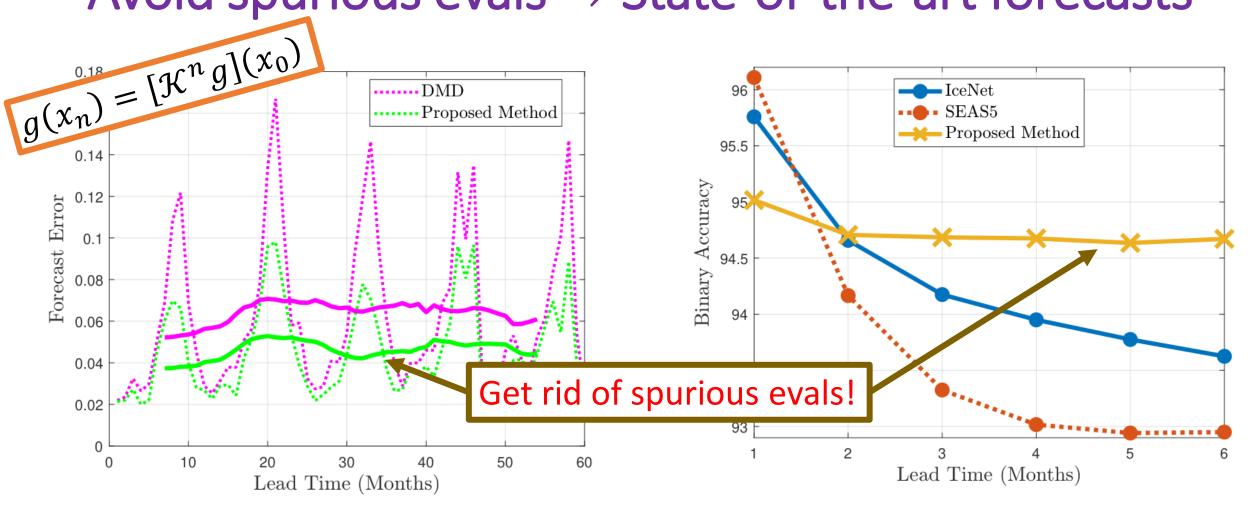
Problems: 1. Very hard to locate geographical significant regions.

2. Very hard to predict more than two months in advance.



• C., Mezić, Stepanenko, "Adversarial Dynamical Systems Reveal Limits and Rules for Trustworthy Data-Driven Learning," preprint, 2025.

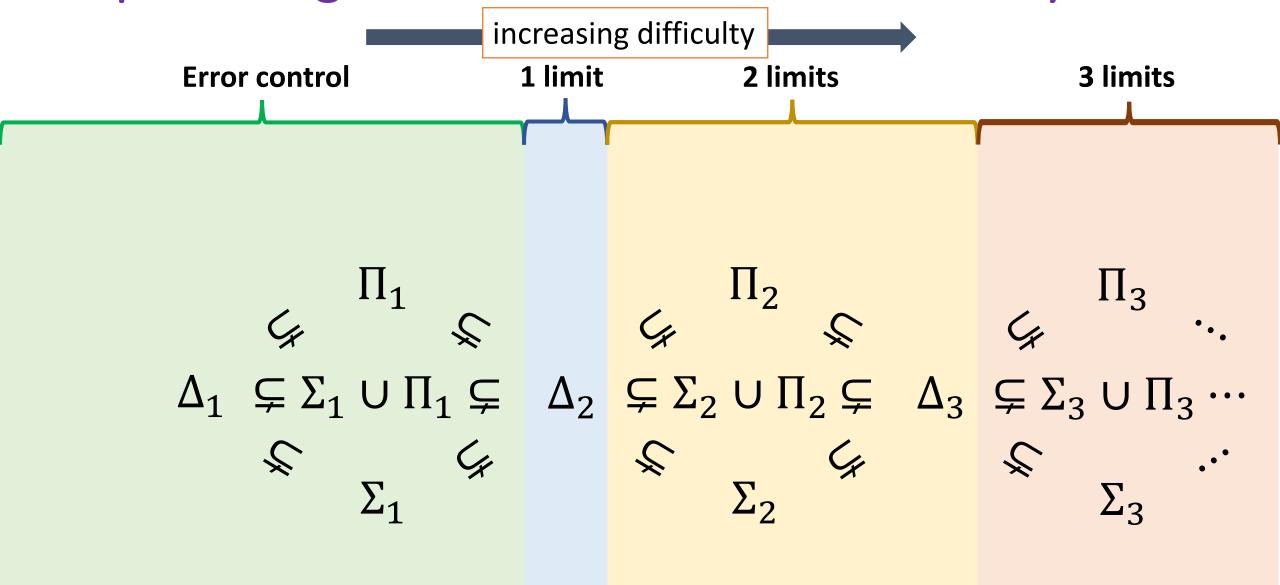
Avoid spurious evals ⇒ State-of-the-art forecasts

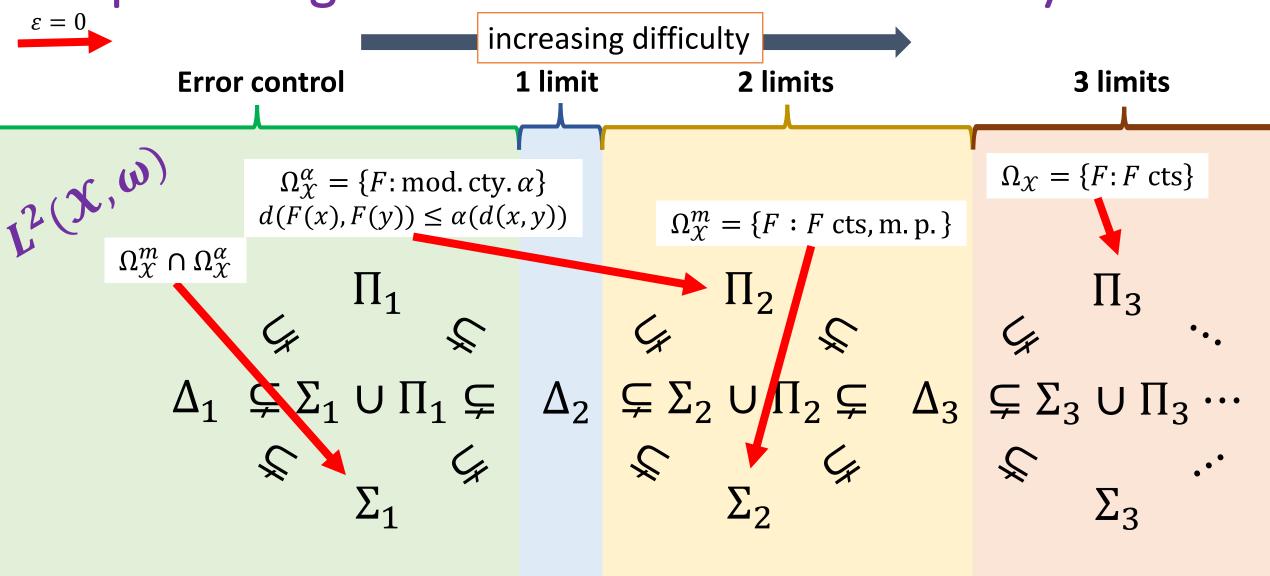


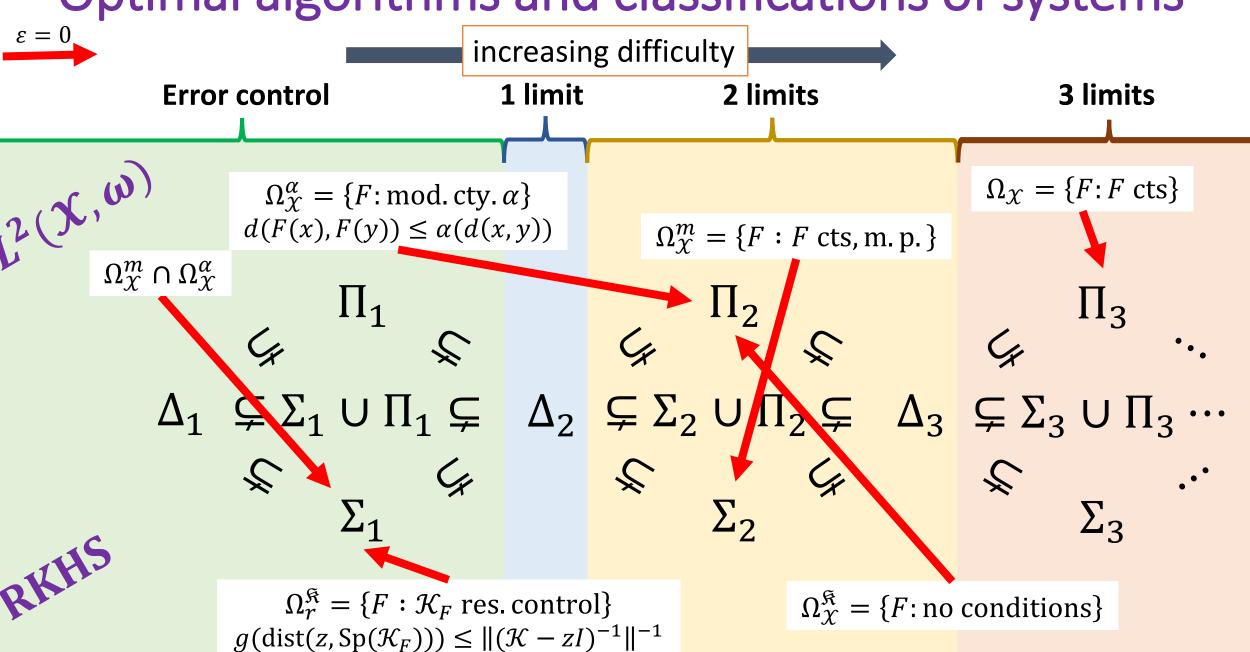
Relative mean squared error over 2016-2020. Model built from 2005-2015 data. (Solid lines moving 12-month mean.)

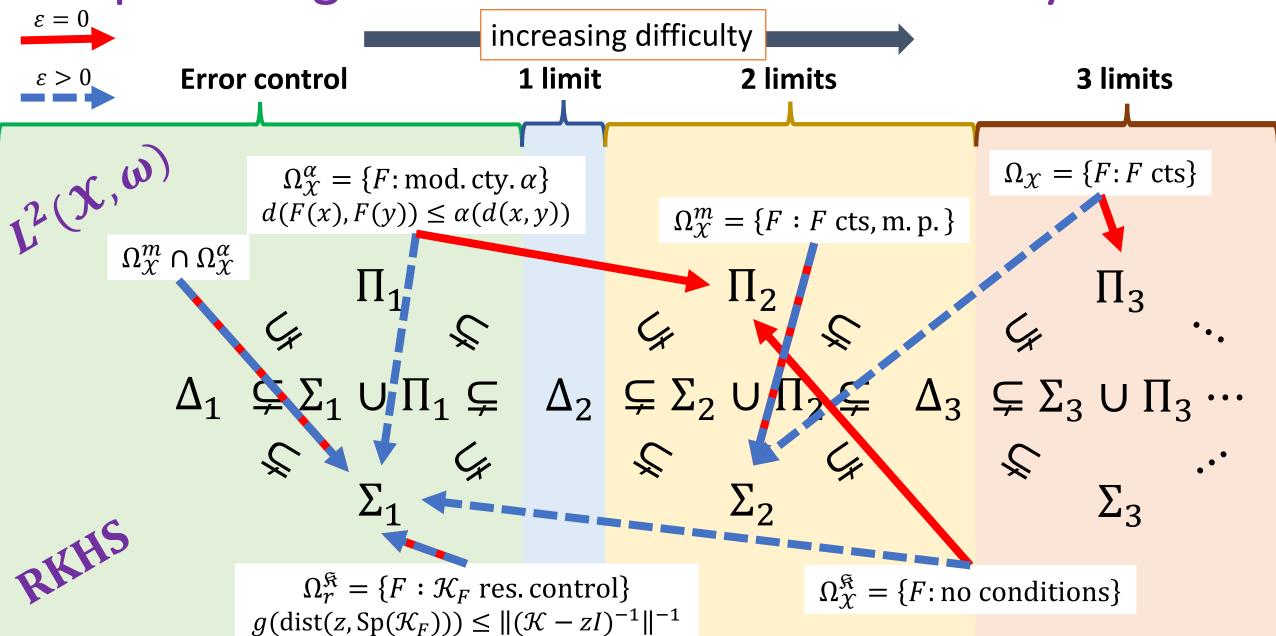
Mean binary accuracy over test years 2012-2020. (IceNet: Andersson et al, "Seasonal Arctic sea ice forecasting with probabilistic deep learning." Nature Communications, 2021.)

• C., Mezić, Stepanenko, "Adversarial Dynamical Systems Reveal Limits and Rules for Trustworthy Data-Driven Learning," preprint, 2025.









Pointers

- Data-driven spectral problems for Koopman operators are hugely popular.
 BUT: Standard truncation methods often fail.
- 2. General method 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 <u>adversarial dynamics</u>.
 Upper bounds ⇒ new "inf.-dim." algorithms. <u>Rigorous, optimal, practical.</u>
- \longrightarrow We now have a near complete picture for Koopman on $L^2(\mathcal{X},\omega)$ and RKHS!

NB: Similar picture has emerged for spectral measures, dealing with continuous spectra (versus eigenvalues) and spectral type (different flavors of dynamics).

Shameless plug...

Upcoming book with CUP:

INFINITE-DIMENSIONAL SPECTRAL COMPUTATIONS

Foundations, Algorithms, and Modern Applications

100s of: classifications, algorithms, examples (webpage: full code), figures, exercises (webpage: full solutions).

**Out by end of 2025 (hopefully!)... **

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