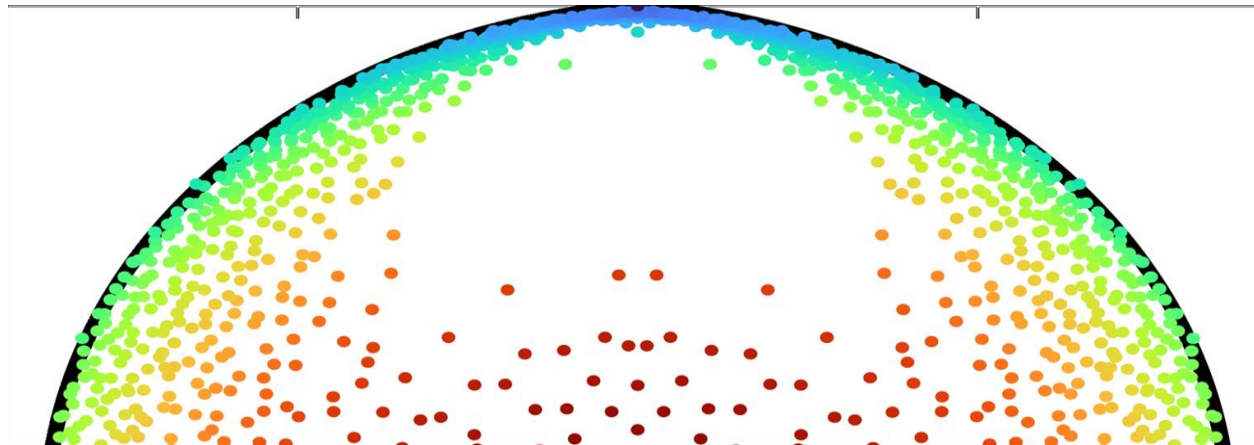


# Residual Dynamic Mode Decomposition

Rigorous data-driven computation of spectral properties  
of Koopman operators for dynamical systems

**Matthew Colbrook** (m.colbrook@damtp.cam.ac.uk)  
University of Cambridge

Joint work with **Alex Townsend** (Cornell University)

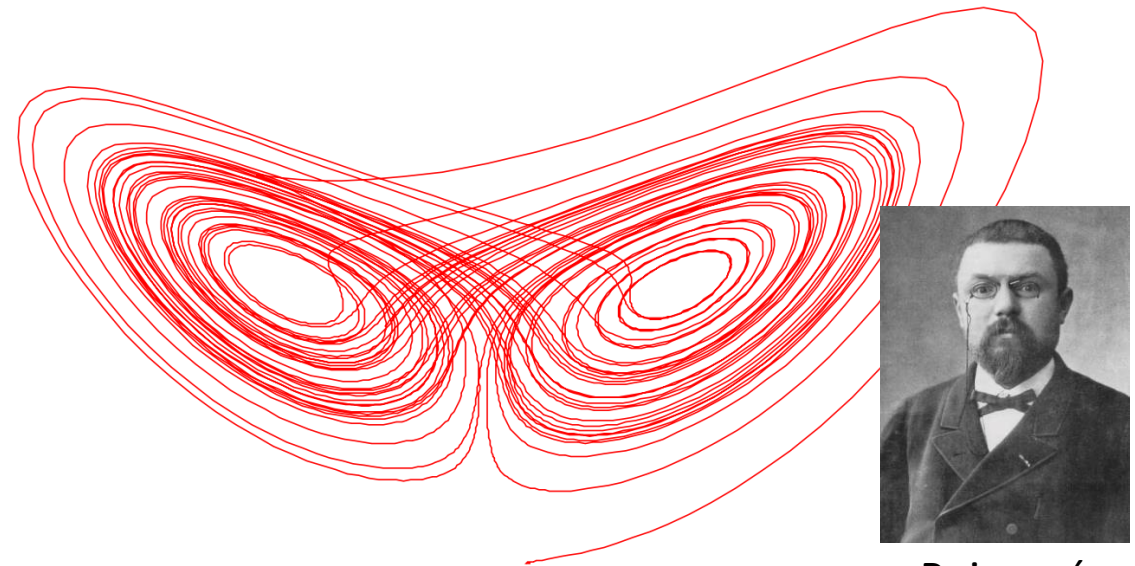


# Data-driven dynamical systems

- State  $x \in \Omega \subseteq \mathbb{R}^d$ , **unknown** function  $F: \Omega \rightarrow \Omega$  governs dynamics

$$x_{n+1} = F(x_n)$$

- **Goal:** Learn about system from data  $\{x^{(m)}, y^{(m)} = F(x^{(m)})\}_{m=1}^M$ 
  - **Data:** experimental measurements or numerical simulations
  - E.g., **used for** forecasting, control, design, understanding
- **Applications:** chemistry, climatology, electronics, epidemiology, finance, fluids, molecular dynamics, neuroscience, plasmas, robotics, video processing, etc.



Poincaré

# Operator viewpoint

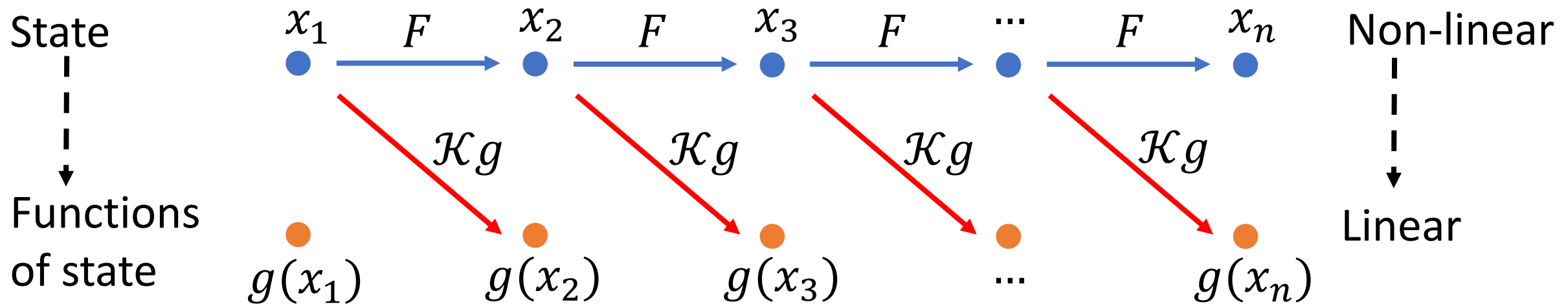
Koopman

von Neumann



- **Koopman operator**  $\mathcal{K}$  acts on functions  $g: \Omega \rightarrow \mathbb{C}$   

$$[\mathcal{K}g](x_n) = g(F(x_n)) = g(x_{n+1})$$
- $\mathcal{K}$  is **linear** but acts on an **infinite-dimensional** space.



- Work in  $L^2(\Omega, \omega)$  for positive measure  $\omega$ , with inner product  $\langle \cdot, \cdot \rangle$ .

• 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.

# Why is linear (much) easier?

$$x_{n+1} = F(x_n)$$

- Suppose  $F(x) = Ax, A \in \mathbb{R}^{d \times d}, A = V\Lambda V^{-1}$ .
- Set  $\xi = V^{-1}x$ ,

$$\xi_n = V^{-1}x_n = V^{-1}A^n x_0 = \Lambda^n V^{-1}x_0 = \Lambda^n \xi_0$$

- Let  $w^T A = \lambda w$ , set  $\varphi(x) = w^T x$ ,

$$[\mathcal{K}\varphi](x) = w^T Ax = \lambda \varphi(x)$$

Long-time dynamics  
become trivial!



**Eigenfunction**

Much more general (**non-linear** and even **chaotic**  $F$ ).



# Koopman mode decomposition

eigenfunction of  $\mathcal{K}$

generalised  
eigenfunction of  $\mathcal{K}$

$$g(x) = \sum_{\text{eigs } \lambda_j} c_{\lambda_j} \varphi_{\lambda_j}(x) + \int_{[-\pi, \pi]_{\text{per}}} \phi_{\theta, g}(x) d\theta$$

$$g(x_n) = [\mathcal{K}^n g](x_0) = \sum_{\text{eigs } \lambda_j} c_{\lambda_j} \lambda_j^n \varphi_{\lambda_j}(x_0) + \int_{[-\pi, \pi]_{\text{per}}} e^{in\theta} \phi_{\theta, g}(x_0) d\theta$$

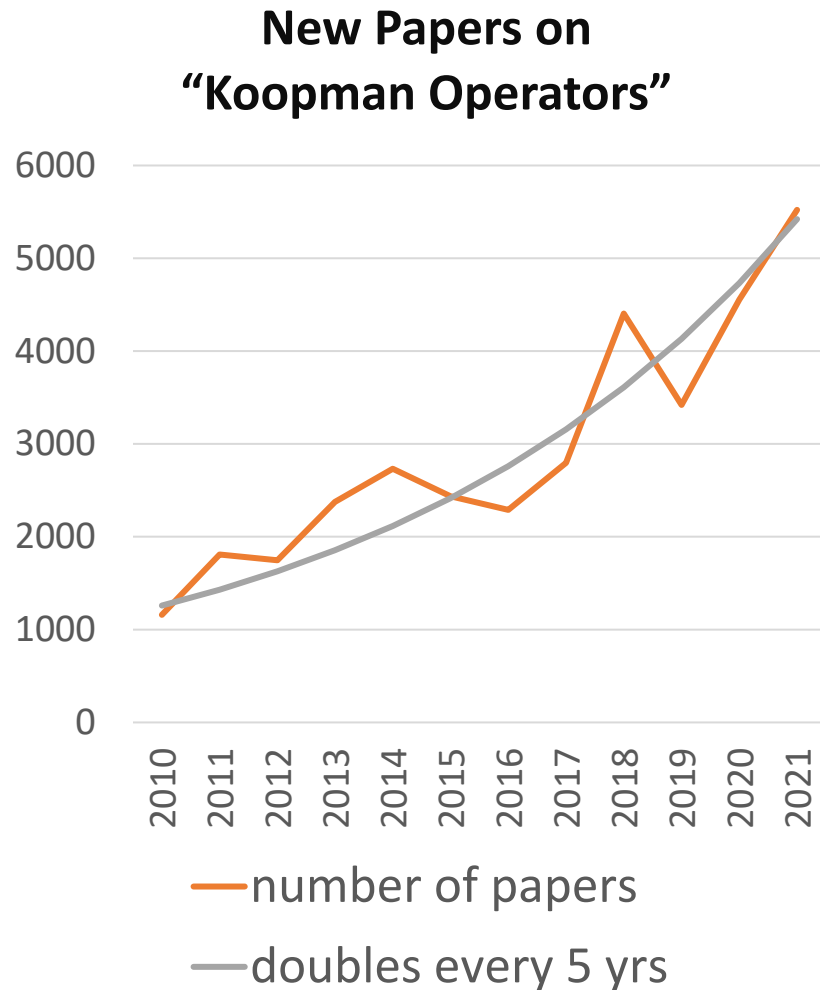
$$[\mathcal{K}g](x) = g(F(x))$$

**Encodes:** geometric features, invariant measures, transient behaviour, long-time behaviour, coherent structures, quasiperiodicity, etc.

**GOAL:** Data-driven approximation of  $\mathcal{K}$  and its spectral properties.

- Mezić, “Spectral properties of dynamical systems, model reduction and decompositions,” **Nonlinear Dynam.**, 2005.

# Koopmania\*: A revolution in the big data era?



≈35,000 papers over last decade!

***BUT: Very little on verified methods!***

**Computing spectra in infinite dimensions is notoriously hard!**

\*Wikipedia: "its wild surge in popularity is sometimes jokingly called 'Koopmania'"

# Challenges of computing

$$\text{Spec}(\mathcal{K}) = \{\lambda \in \mathbb{C}: \mathcal{K} - \lambda I \text{ is not invertible}\}$$

**Truncate:**  $\mathcal{K} \longrightarrow \mathbb{K} \in \mathbb{C}^{N_K \times N_K}$

- 1) **“Too much”:** Approximate spurious modes  $\lambda \notin \text{Spec}(\mathcal{K})$
- 2) **“Too little”:** Miss parts of  $\text{Spec}(\mathcal{K})$
- 3) **Continuous spectra.**

**Verification:** Is it right?

# Build the matrix: Dynamic Mode Decomposition (DMD)

Given dictionary  $\{\psi_1, \dots, \psi_{N_K}\}$  of functions  $\psi_j: \Omega \rightarrow \mathbb{C}$ ,

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

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

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

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

**Recall open problems:** too much, too little, continuous spectra, verification

- 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.
- Kutz, Brunton, Brunton, Proctor, "Dynamic mode decomposition: data-driven modeling of complex systems," **SIAM**, 2016.
- Williams, Kevrekidis, Rowley "A data-driven approximation of the Koopman operator: Extending dynamic mode decomposition," **J. Nonlinear Sci.**, 2015.

# Residual DMD (ResDMD): Approx. $\mathcal{K}$ and $\mathcal{K}^*\mathcal{K}$

$$\begin{aligned}\langle \psi_k, \psi_j \rangle &\approx \sum_{m=1}^M w_m \overline{\psi_j(x^{(m)})} \psi_k(x^{(m)}) = \left[ \underbrace{\Psi_X^* W \Psi_X}_G \right]_{jk} \\ \langle \mathcal{K}\psi_k, \psi_j \rangle &\approx \sum_{m=1}^M w_m \overline{\psi_j(x^{(m)})} \underbrace{\psi_k(y^{(m)})}_{[\mathcal{K}\psi_k](x^{(m)})} = \left[ \underbrace{\Psi_X^* W \Psi_Y}_{K_1} \right]_{jk} \\ \langle \mathcal{K}\psi_k, \mathcal{K}\psi_j \rangle &\approx \sum_{m=1}^M w_m \overline{\psi_j(y^{(m)})} \psi_k(y^{(m)}) = \left[ \underbrace{\Psi_Y^* W \Psi_Y}_{K_2} \right]_{jk}\end{aligned}$$

**Residuals:**  $g = \sum_{j=1}^{N_K} \mathbf{g}_j \psi_j, \quad \|\mathcal{K}g - \lambda g\|^2 \approx \mathbf{g}^* [K_2 - \lambda K_1^* - \bar{\lambda} K_1 + |\lambda|^2 G] \mathbf{g}$

- 
- C., Townsend, “Rigorous data-driven computation of spectral properties of Koopman operators for dynamical systems,” preprint.
  - C., Ayton, Szóke, “Residual Dynamic Mode Decomposition,” **J. Fluid Mech.**, to appear.
  - Code: <https://github.com/MColbrook/Residual-Dynamic-Mode-Decomposition>

# ResDMD: avoiding “too much”

$$\text{res}(\lambda, \mathbf{g})^2 = \frac{\mathbf{g}^* [K_2 - \lambda K_1^* - \bar{\lambda} K_1 + |\lambda|^2 G] \mathbf{g}}{\mathbf{g}^* G \mathbf{g}}$$

eigenvectors

eigenvalues

## Algorithm 1:

1. Compute  $G, K_1, K_2 \in \mathbb{C}^{N_K \times N_K}$  and eigendecomposition  $K_1 V = G V \Lambda$ .
2. For each eigenpair  $(\lambda, \mathbf{v})$ , compute  $\text{res}(\lambda, \mathbf{v})$ .
3. **Output:** subset of e-vectors  $V_{(\varepsilon)}$  & e-vals  $\Lambda_{(\varepsilon)}$  with  $\text{res}(\lambda, \mathbf{v}) \leq \varepsilon$  ( $\varepsilon = \text{input tol}$ ).

**Theorem (no spectral pollution):** Suppose quad. rule converges. Then

$$\limsup_{M \rightarrow \infty} \max_{\lambda \in \Lambda^{(\varepsilon)}} \|(\mathcal{K} - \lambda)^{-1}\|^{-1} \leq \varepsilon$$

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**BUT:** Typically, does not capture all of spectrum! (“too little”)



# ResDMD: avoiding “too little”

$$\text{Spec}_\varepsilon(\mathcal{K}) = \bigcup_{\|\mathcal{B}\| \leq \varepsilon} \text{Spec}(\mathcal{K} + \mathcal{B}), \quad \lim_{\varepsilon \downarrow 0} \text{Spec}_\varepsilon(\mathcal{K}) = \text{Spec}(\mathcal{K})$$

Algorithm 2:

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

1. Compute  $G, K_1, K_2 \in \mathbb{C}^{N_K \times N_K}$ .
2. For  $z_k$  in comp. grid, compute  $\tau_k = \min_{g = \sum_{j=1}^{N_K} \mathbf{g}_j \psi_j} \text{res}(z_k, g)$ , corresponding  $g_k$  (gen. SVD).
3. **Output:**  $\{z_k: \tau_k < \varepsilon\}$  (approx. of  $\text{Spec}_\varepsilon(\mathcal{K})$ ),  $\{g_k: \tau_k < \varepsilon\}$  ( $\varepsilon$ -pseudo-eigenfunctions).

**Theorem (full convergence):** Suppose the quadrature rule converges.

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

# Quadrature with trajectory data

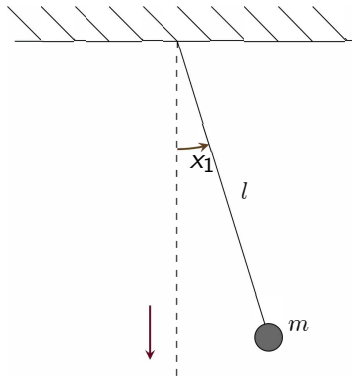
$$\text{E.g., } \langle \mathcal{K}\psi_k, \psi_j \rangle = \lim_{M \rightarrow \infty} \sum_{m=1}^M w_m \overline{\psi_j(x^{(m)})} \underbrace{\psi_k(y^{(m)})}_{[\mathcal{K}\psi_k](x^{(m)})}$$

Three examples:

- **High-order quadrature:**  $\{x^{(m)}, w_m\}_{m=1}^M$   $M$ -point quadrature rule.  
 Rapid convergence. Requires free choice of  $\{x^{(m)}\}_{m=1}^M$  and small  $d$ .
- **Random sampling:**  $\{x^{(m)}\}_{m=1}^M$  selected at random.  
 Large  $d$ . Slow Monte Carlo  $O(M^{-1/2})$  rate of convergence. ← Most common
- **Ergodic sampling:**  $x^{(m+1)} = F(x^{(m)})$ .  
 Single trajectory, large  $d$ . Requires ergodicity, convergence can be slow. ↘

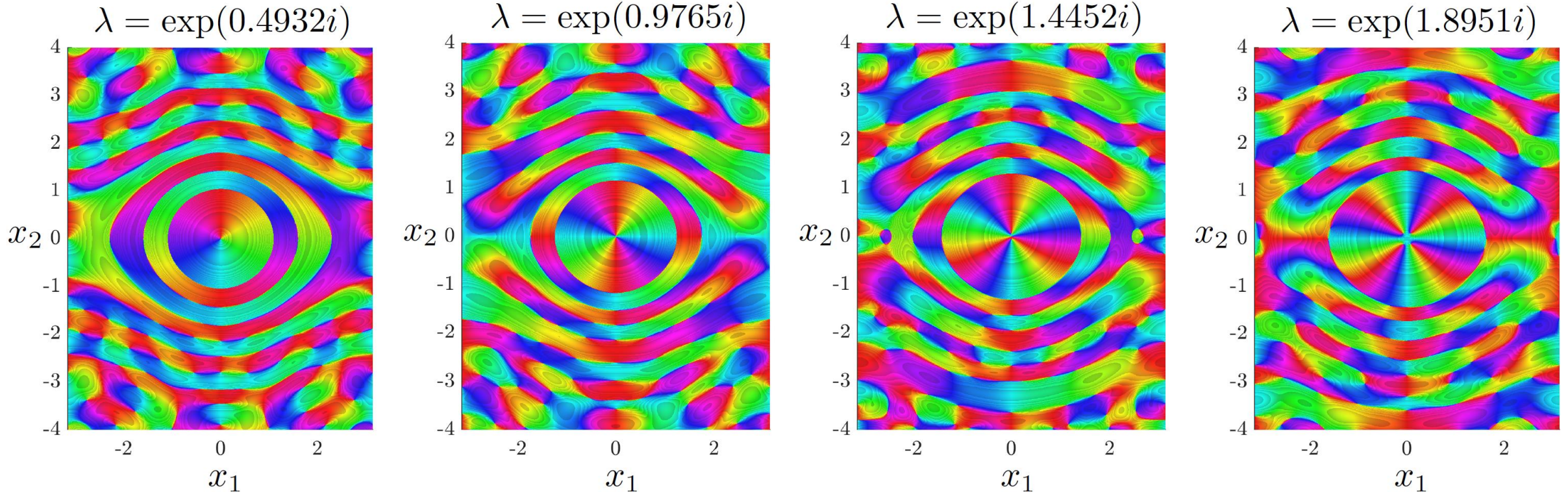
## Example: non-linear pendulum

$$\dot{x}_1 = x_2, \quad \dot{x}_2 = -\sin(x_1), \quad \Omega = [-\pi, \pi]_{\text{per}} \times \mathbb{R}$$



Computed pseudospectra ( $\varepsilon = 0.25$ ). Eigenvalues of  $\mathbb{K}$  shown as dots (spectral pollution).

# Approximate eigenfunctions



Colour represents complex argument, constant modulus shown as shadowed steps.  
All residuals smaller than  $\varepsilon = 0.05$  (made smaller by increasing  $N_K$ ).

# The Challenges

- ~~1) “Too much”: Approximate spurious modes  $\lambda \notin \text{Spec}(\mathcal{K})$~~  ✓
- ~~2) “Too little”: Miss parts of  $\text{Spec}(\mathcal{K})$~~  ✓
- 3) Continuous spectra.

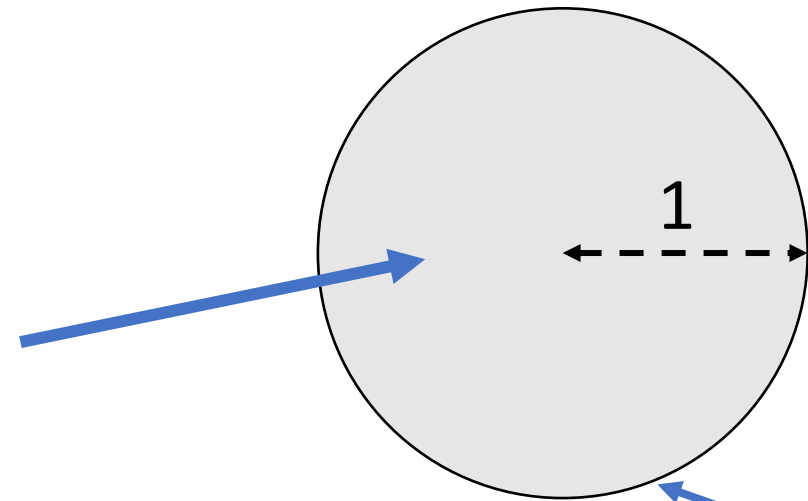
**Verification:** Is it right?

# Setup for continuous spectra

Suppose system is measure preserving (e.g., Hamiltonian, ergodic, post-transient etc.)

$$\Leftrightarrow \mathcal{K}^* \mathcal{K} = I \text{ (isometry)}$$

$$\Rightarrow \text{Spec}(\mathcal{K}) \subseteq \{z: |z| \leq 1\}$$



(NB: we consider unitary extensions via Wold decomposition.)

spectral  
measure  
supp. on  
boundary

# Spectral decomposition of operators

$A \in \mathbb{C}^{n \times n}$  normal  $\Rightarrow$  O.N. basis of eigenvectors  $v_1, \dots, v_n$ :

$$v = \left( \sum_{k=1}^n \underset{\substack{\uparrow \\ \text{Projector onto Span}(v_k)}}{v_k v_k^*} \right) v, \quad Av = \left( \sum_{k=1}^n \underset{\substack{\uparrow \\ \text{eigenvalues}}}{\lambda_k} v_k v_k^* \right) v, \quad v \in \mathbb{C}^n$$



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Energy of “v” in each eigenvector:  $\mu_v(\lambda_j) = \langle v_j v_j^* v, v \rangle = |v_j^* v|^2$

This is called the spectral measure with respect to a vector  $v$ .

# Spectral decomposition of operators

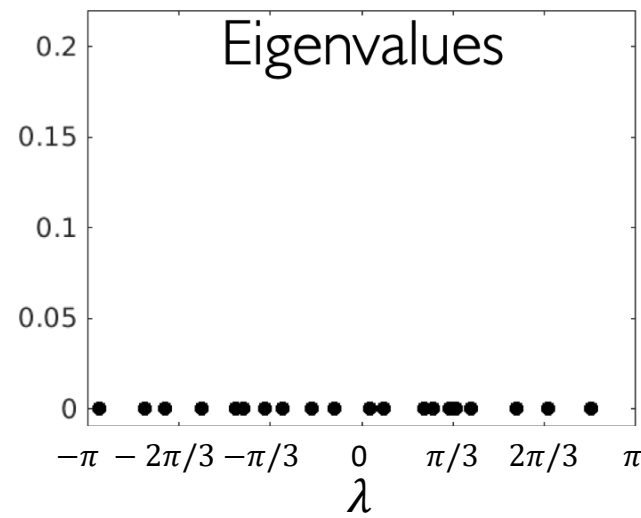
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↑  
Projector onto  $\text{Span}(v_k)$ 
↑  
eigenvalues

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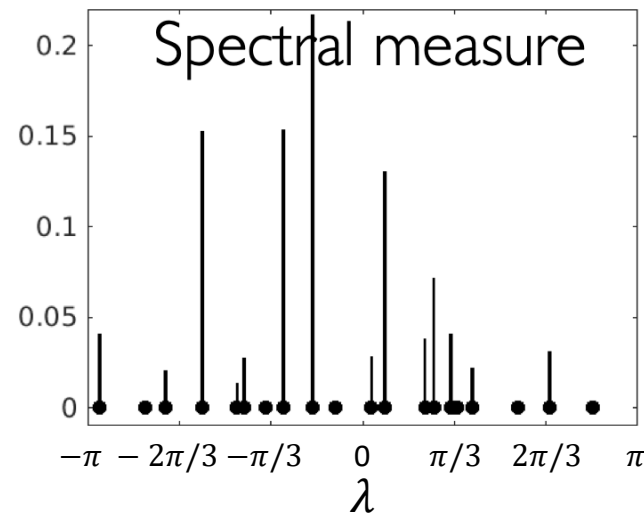
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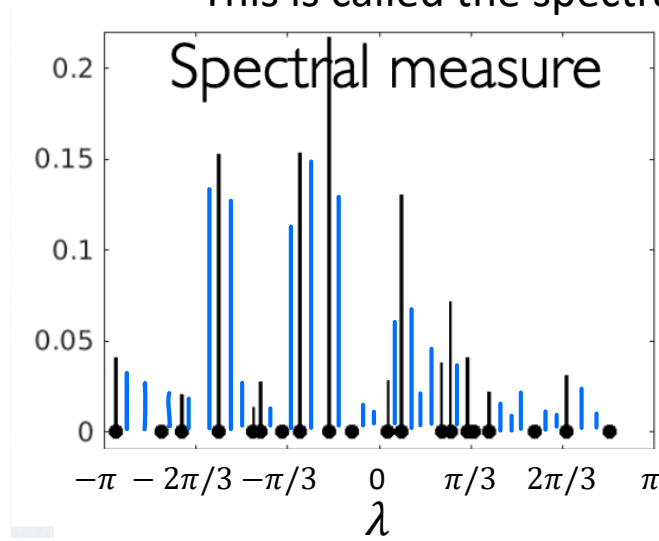
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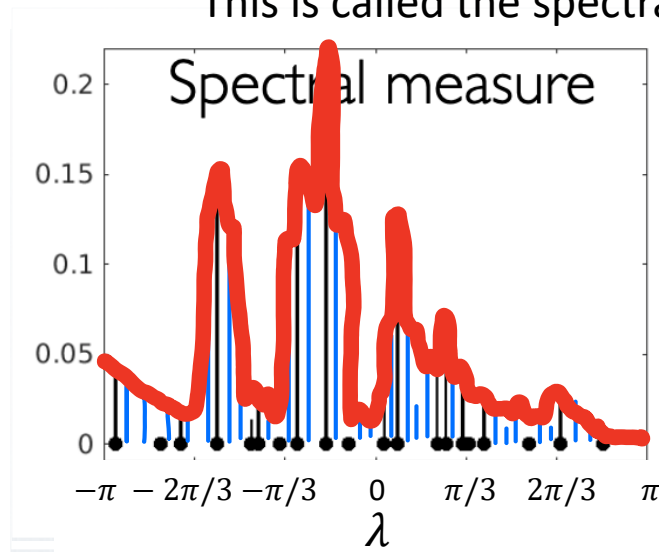
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$\mathcal{K}$  is unitary  $\Rightarrow$  projection-valued measure  $\xi$

$$g = \left( \int_{\mathbb{T}} d\xi(y) \right) g, \quad \mathcal{K}g = \left( \int_{\mathbb{T}} y d\xi(y) \right) g$$

Spectral measure  $\nu_g(B) = \langle \xi(B)g, g \rangle$

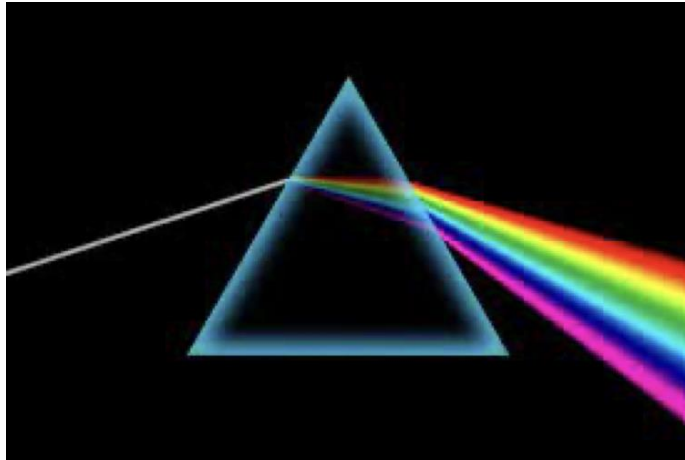
# Spectral decomposition of operators

$A \in \mathbb{C}^{n \times n}$  normal  $\Rightarrow$

$\Rightarrow$

O

White light contains a continuous spectra



$v,$

$A$

$\text{an}(v_k)$

eigenvector:

This is ca

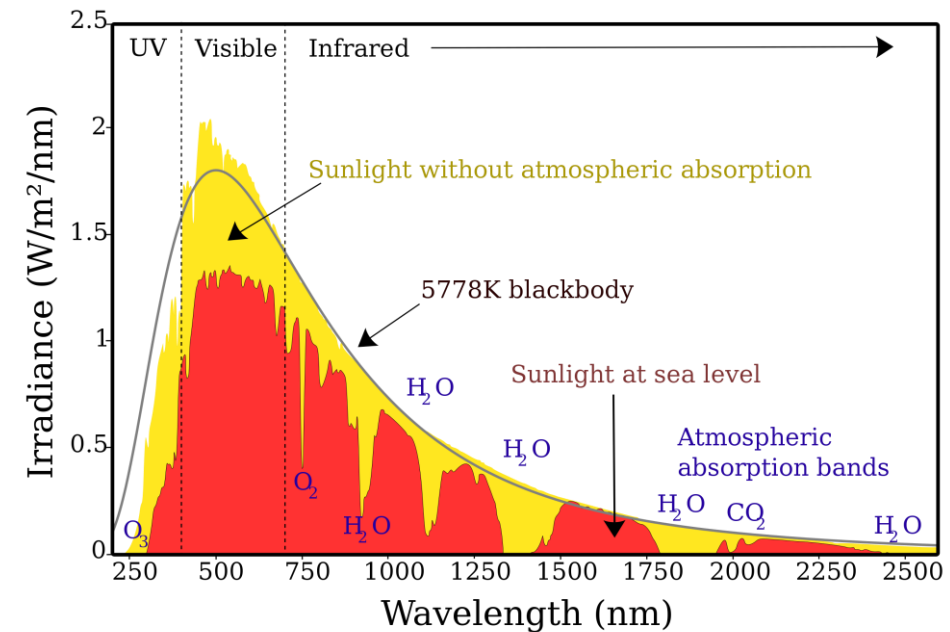
$\Rightarrow$

$$g = \left( \int_{\mathbb{T}} d\xi(y) \right) g,$$

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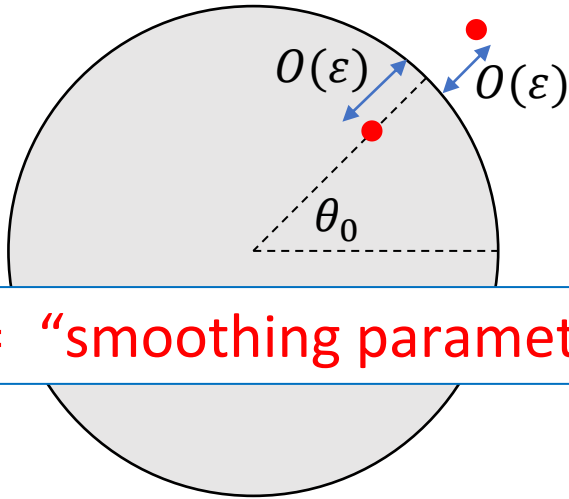
Often interesting to look at  
the intensity of each wavelength

Spectrum of Solar Radiation (Earth)





# Evaluating spectral measure



$\varepsilon =$  “smoothing parameter”

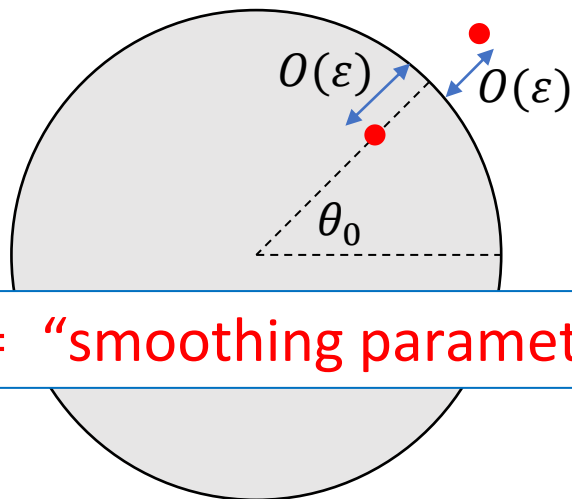
$$[P_\varepsilon * \nu_g](\theta_0) = \int_{[-\pi, \pi]_{\text{per}}} P_\varepsilon(\theta_0 - \theta) d\nu_g(\theta)$$

Smoothing convolution

Poisson kernel for  
unit disk

$$P_\varepsilon(\theta_0) = \frac{1}{2\pi} \frac{(1 + \varepsilon)^2 - 1}{1 + (1 + \varepsilon)^2 - 2(1 + \varepsilon)\cos(\theta_0)}$$

# Evaluating spectral measur



$\epsilon =$  "smoothing parameter"

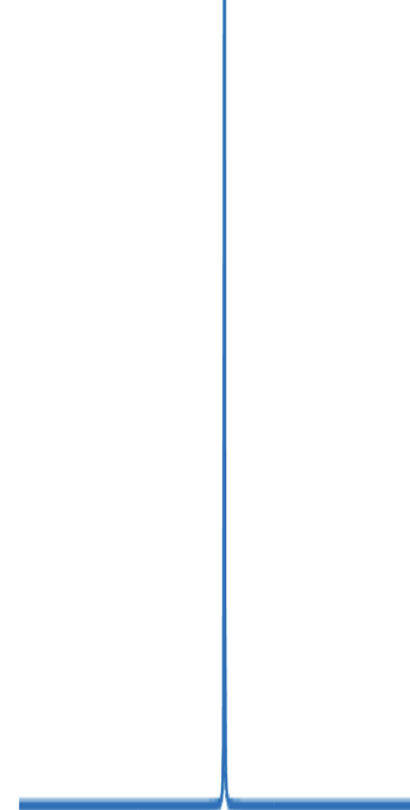
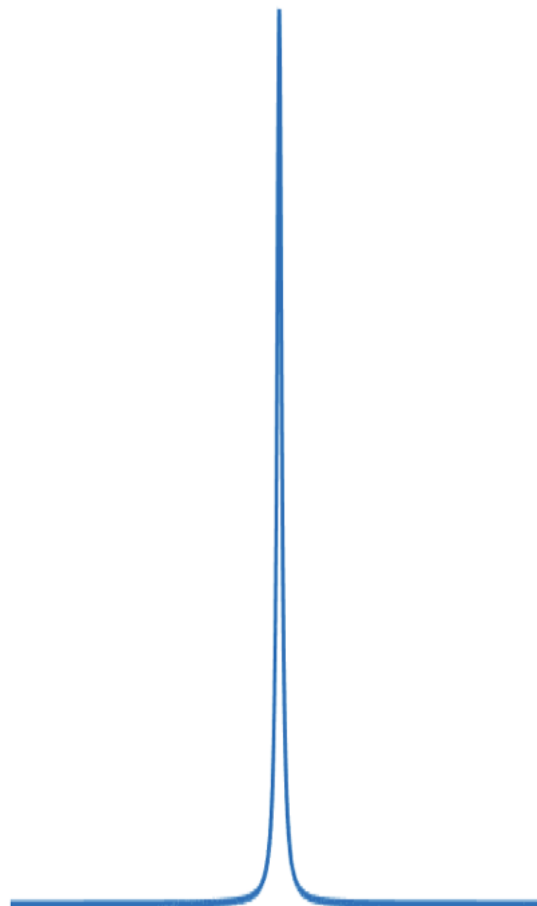
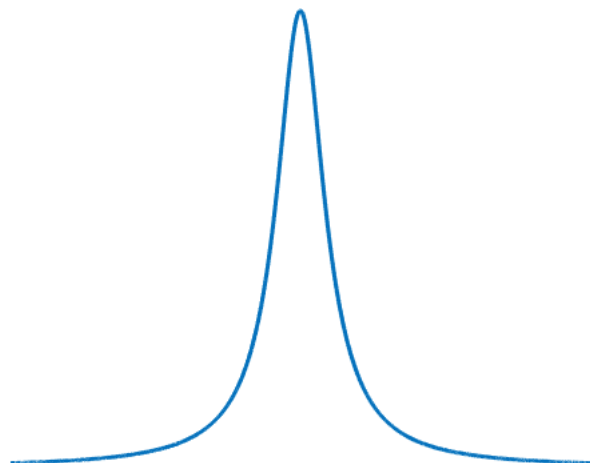
Poisso  
unit di

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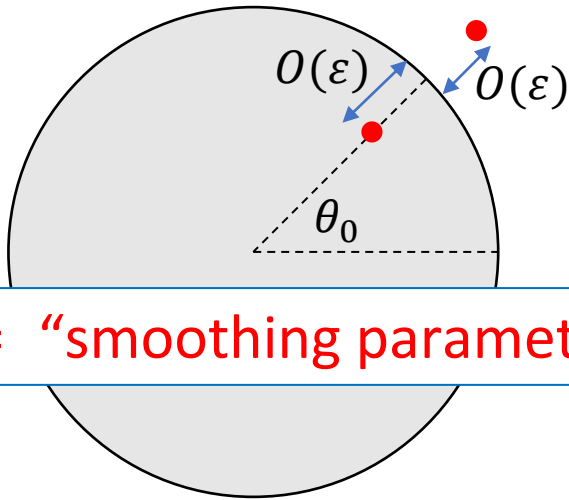
Smoothing co

$$\frac{1}{1 + \epsilon}$$

$$\overline{0}$$



# Evaluating spectral measure



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Smoothing convolution

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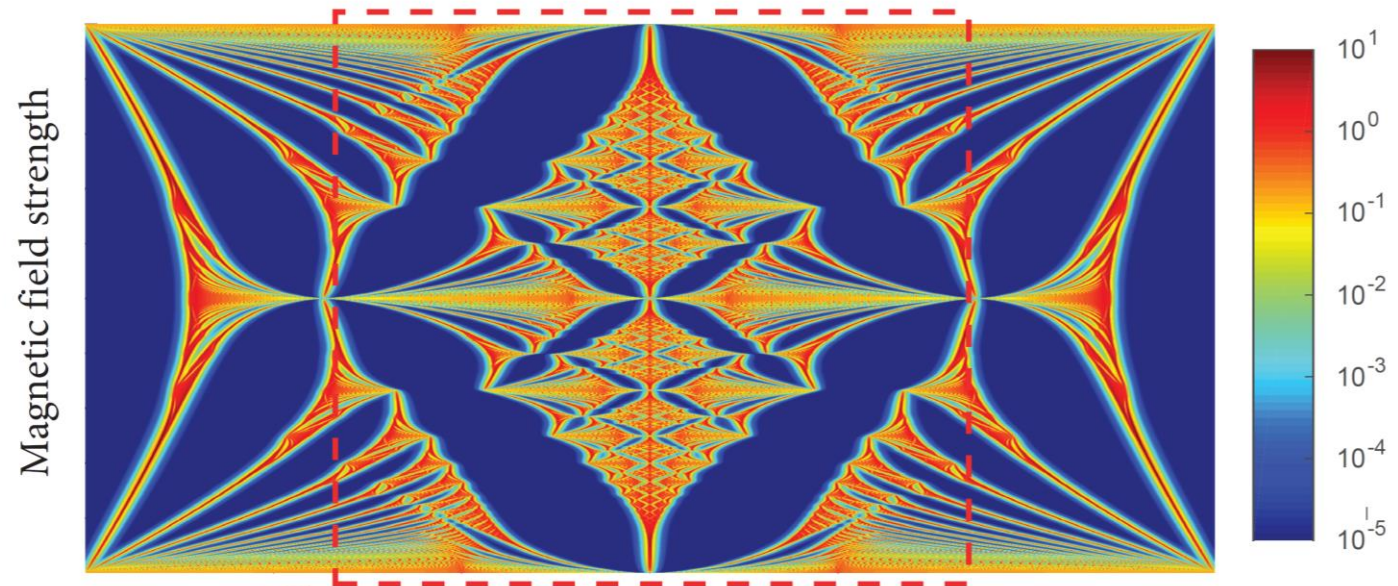
$$P_\varepsilon(\theta_0) = \frac{1}{2\pi} \frac{(1 + \varepsilon)^2 - 1}{1 + (1 + \varepsilon)^2 - 2(1 + \varepsilon)\cos(\theta_0)}$$

$$[P_\varepsilon * \nu_g](\theta_0) = \mathcal{C}_g(e^{i\theta_0}(1 + \varepsilon)^{-1}) - \mathcal{C}_g(e^{i\theta_0}(1 + \varepsilon))$$

$$\mathcal{C}_g(z) = \int_{[-\pi, \pi]_{\text{per}}} \frac{e^{i\theta} d\nu_g(\theta)}{e^{i\theta} - z} = \begin{cases} \langle (\mathcal{K} - zI)^{-1}g, \mathcal{K}^*g \rangle, & \text{if } |z| > 1 \\ -z^{-1} \langle g, (\mathcal{K} - \bar{z}^{-1}I)^{-1}g \rangle, & \text{if } 0 < |z| < 1 \end{cases}$$

ResDMD computes  
with error control

# Spectral measures of self-adjoint operators



Horizontal slice = spectral measure at constant magnetic field strength.

## Software package

SpecSolve available at <https://github.com/SpecSolve>  
 Capabilities: ODEs, PDEs, integral operators, discrete operators.

# Example

$$\mathcal{K} = \begin{pmatrix} \overline{\alpha_0} & \overline{\alpha_1}\rho_0 & \rho_0\rho_1 & & & \\ \rho_0 & -\overline{\alpha_1}\alpha_0 & -\alpha_0\rho_1 & & & \\ & \overline{\alpha_2}\rho_1 & -\overline{\alpha_2}\alpha_1 & \overline{\alpha_3}\rho_2 & \rho_3\rho_2 & \\ & \rho_2\rho_1 & -\alpha_1\rho_2 & -\overline{\alpha_3}\alpha_2 & -\rho_3\alpha_2 & \ddots \\ & & & \overline{\alpha_4}\rho_3 & -\overline{\alpha_4}\alpha_3 & \ddots \\ & & & \ddots & \ddots & \ddots \end{pmatrix}$$

$$\alpha_j = (-1)^j 0.95^{(j+1)/2}, \quad \rho_j = \sqrt{1 - |\alpha_j|^2}$$

Generalised shift, typical building block of many dynamical systems.

Fix  $N_K$ , vary  $\varepsilon$ : unstable!

Fix  $\varepsilon$ , vary  $N_K$ : too smooth!

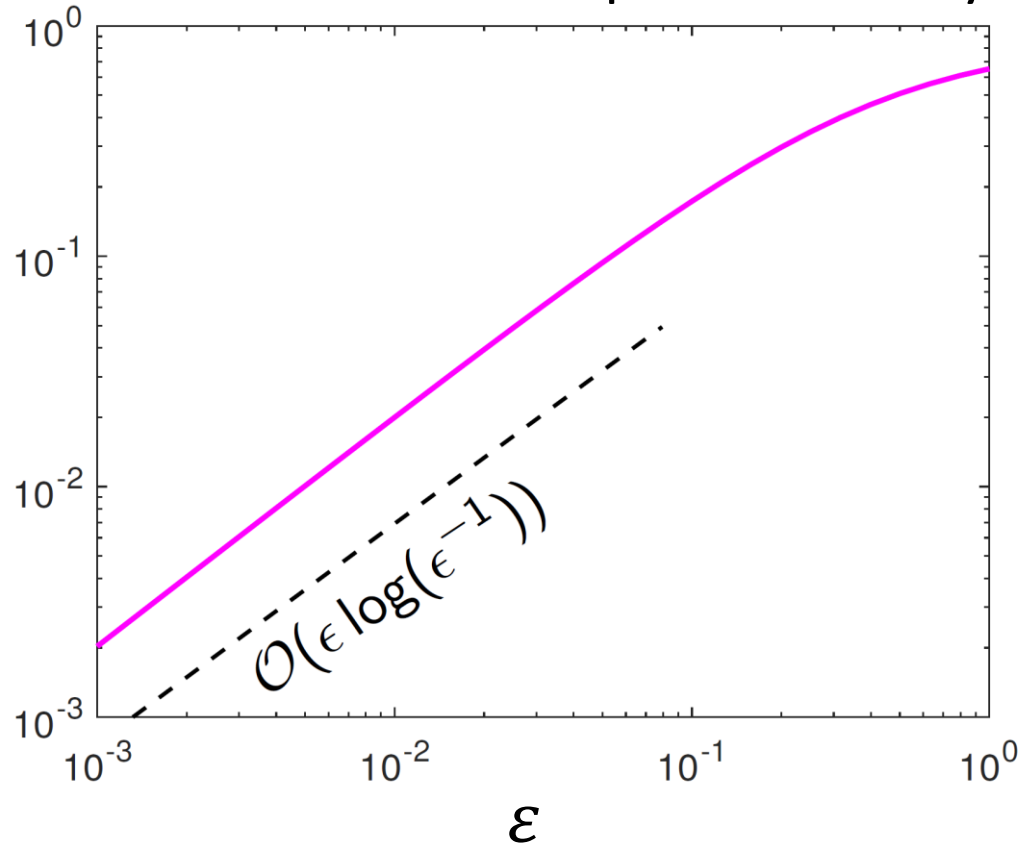


Adaptive: new matrix to compute residuals crucial

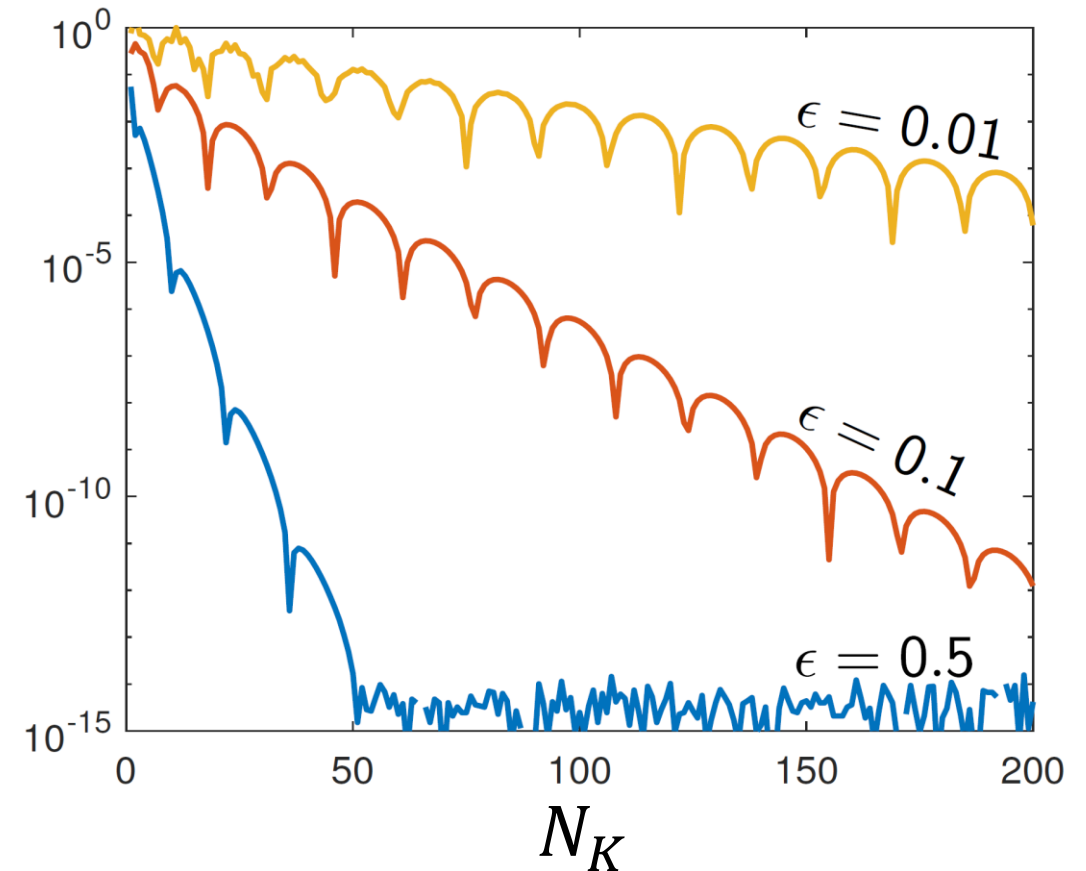
# But ... slow convergence

**Problem:** As  $\varepsilon \downarrow 0$ , error is  $O(\varepsilon \log(1/\varepsilon))$  and  $N_K(\varepsilon) \rightarrow \infty$ .

Pointwise error for spectral density



Error due to discretization

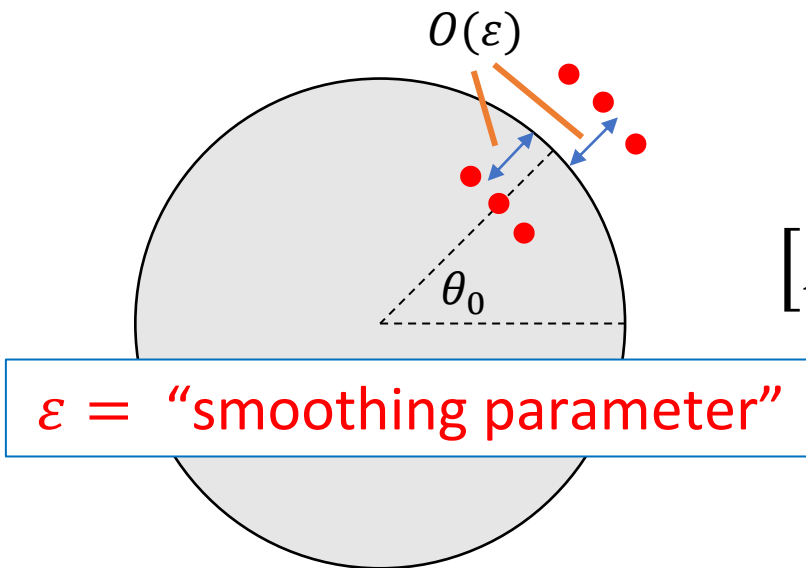


Small  $N_K$  critical in data-driven computations. Can we improve convergence rate?

# High-order rational kernels

$m$ th order rational kernels:

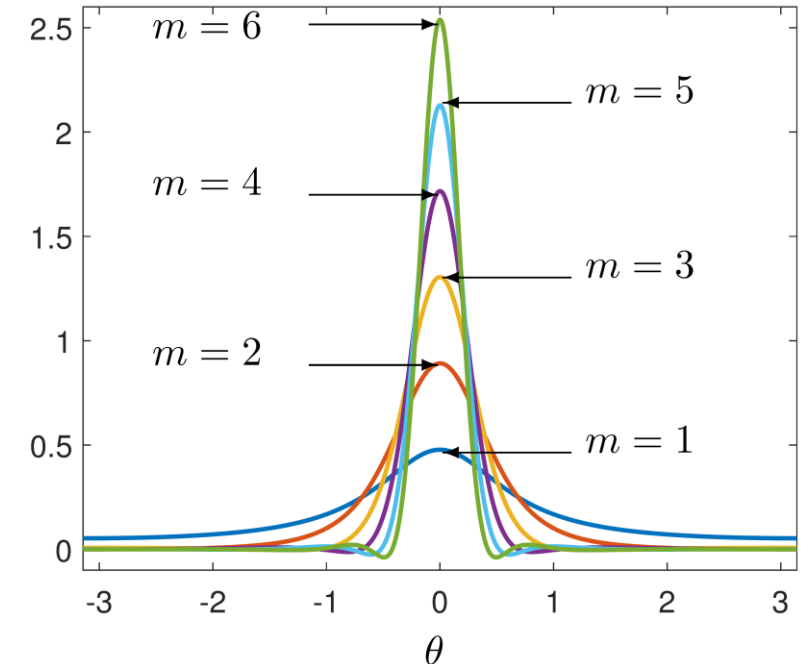
$$K_\varepsilon(\theta) = \frac{e^{-i\theta}}{2\pi} \sum_{j=1}^m \left[ \frac{c_j}{e^{-i\theta} - (1 + \varepsilon \bar{z}_j)^{-1}} - \frac{d_j}{e^{-i\theta} - (1 + \varepsilon z_j)} \right]$$



ResDMD computes  
with error control

$$[K_\varepsilon * \nu_g](\theta_0) = \sum_{j=1}^m \left[ c_j \mathcal{C}_g(e^{i\theta_0}(1 + \varepsilon \bar{z}_j)^{-1}) - d_j \mathcal{C}_g(e^{i\theta_0}(1 + \varepsilon z_j)) \right]$$

Kernels



Smaller  $N_K$  (larger  $\varepsilon$ )

# Convergence

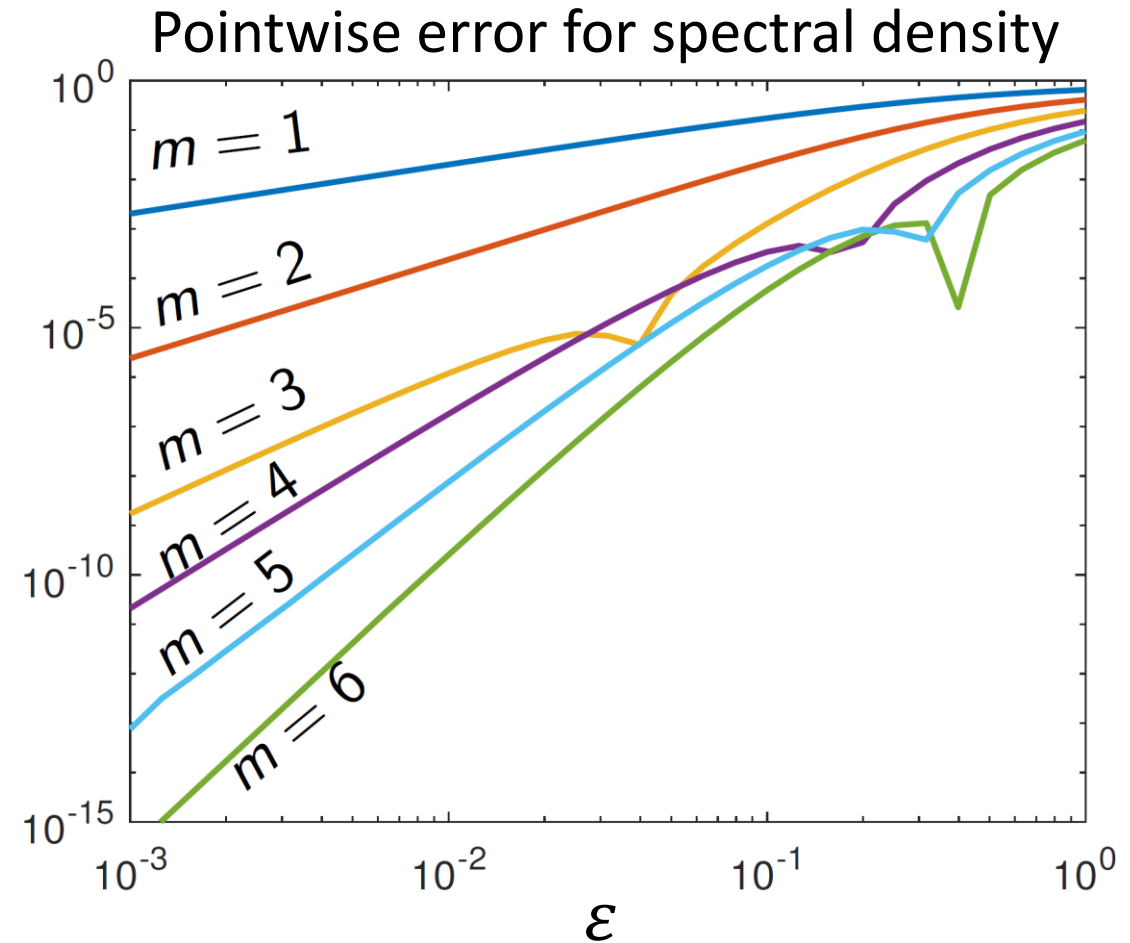
**Theorem:** Automatic selection of  $N_K(\varepsilon)$  with  $O(\varepsilon^m \log(1/\varepsilon))$  convergence:

- Density of continuous spectrum  $\rho_g$ .  
(pointwise and  $L^p$ )
- Integration against test functions.  
(weak convergence)

$$\int_{[-\pi, \pi]_{\text{per}}} h(\theta) [K_\varepsilon * \nu_g](\theta) \, d\theta$$

$$= \int_{[-\pi, \pi]_{\text{per}}} h(\theta) \, d\nu_g(\theta) + O(\varepsilon^m \log(1/\varepsilon))$$

Also recover discrete spectrum.

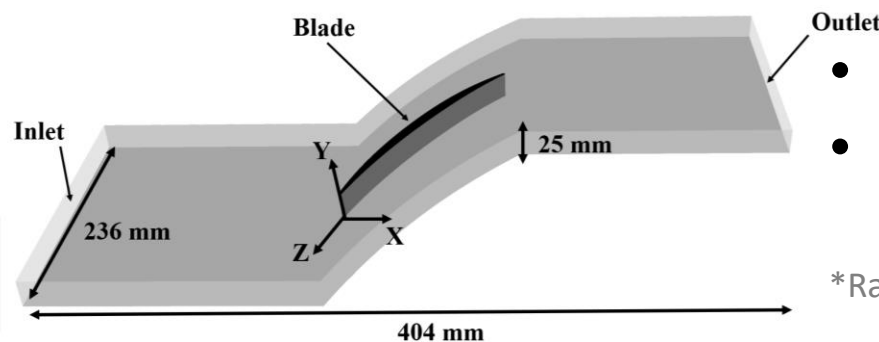
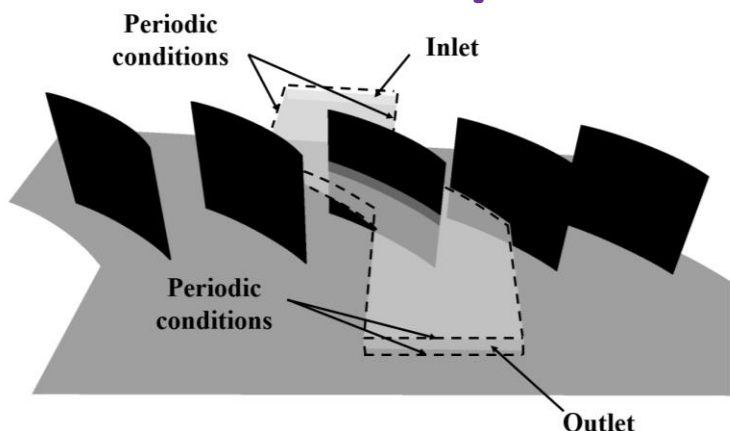


# The Challenges

- ~~1) “Too much”: Approximate spurious modes  $\lambda \notin \text{Spec}(\mathcal{K})$~~  ✓
- ~~2) “Too little”: Miss parts of  $\text{Spec}(\mathcal{K})$~~  ✓
- ~~3) Continuous spectra.~~ ✓

**Verification:** Is it right?

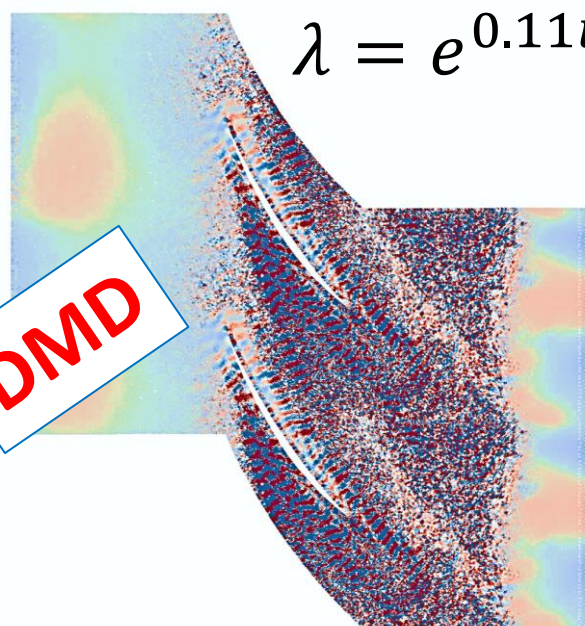
# Example: Trustworthy computation for large $d$



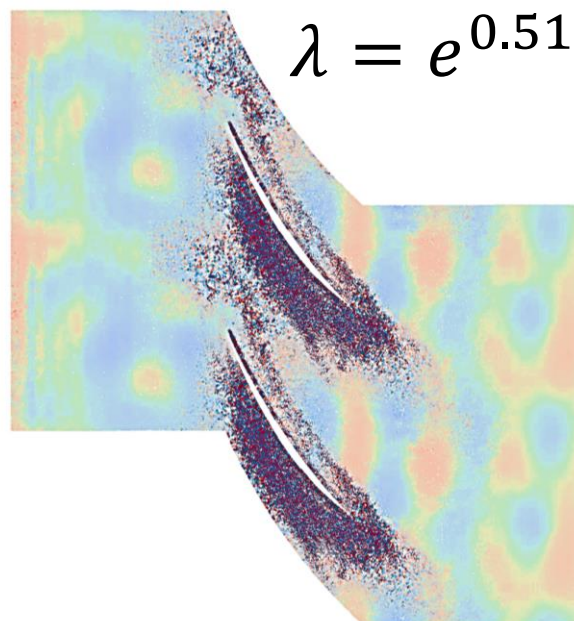
- Reynolds number  $\approx 3.9 \times 10^5$
- Ambient dimension ( $d$ )  $\approx 300,000$  (number of measurement points)

\*Raw measurements provided by Stephane Moreau (Sherbrooke)

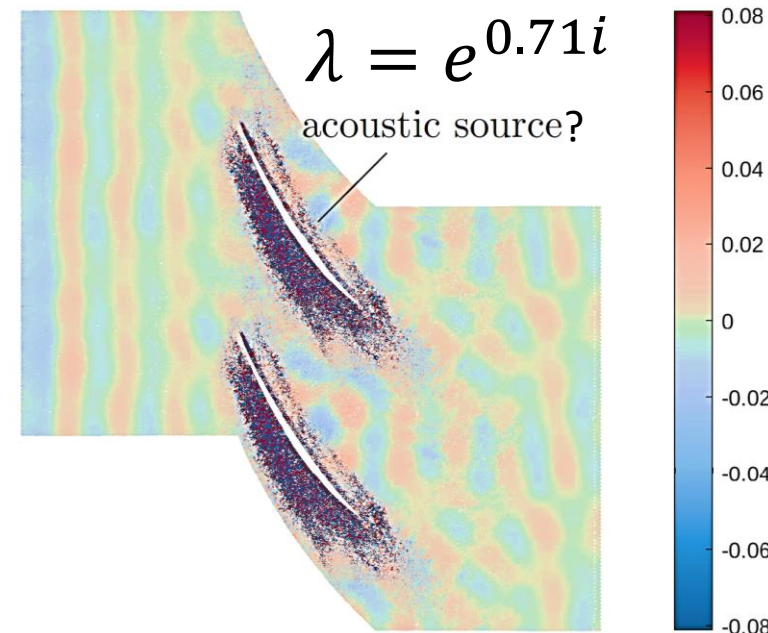
Rel. Error = ?  
 $\lambda = e^{0.11i}$



Rel. Error = ?  
 $\lambda = e^{0.51i}$

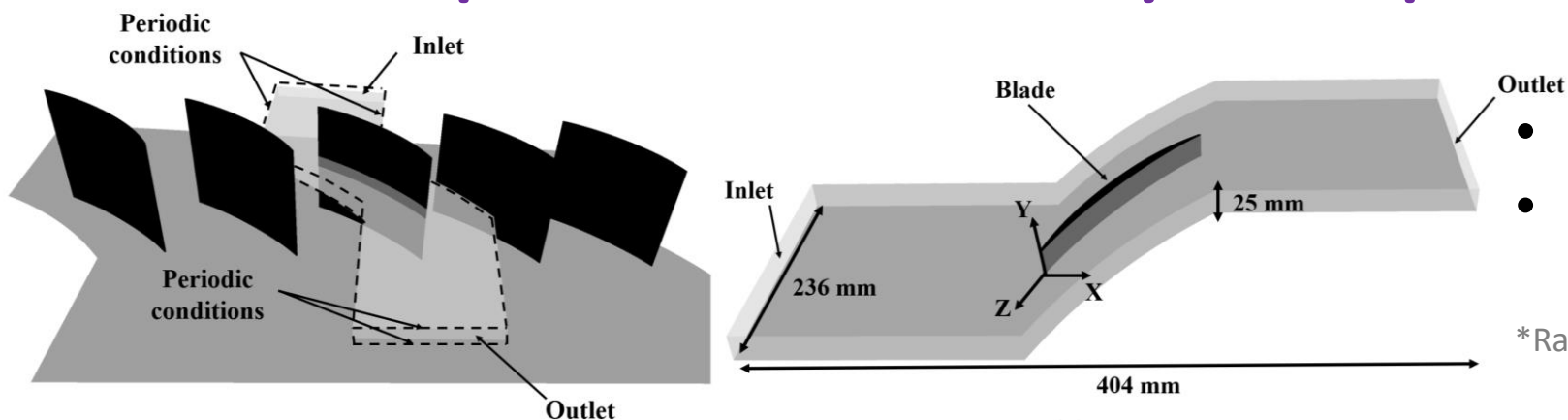


Rel. Error = ?  
 $\lambda = e^{0.71i}$   
 acoustic source?





# Example: Trustworthy computation for large $d$



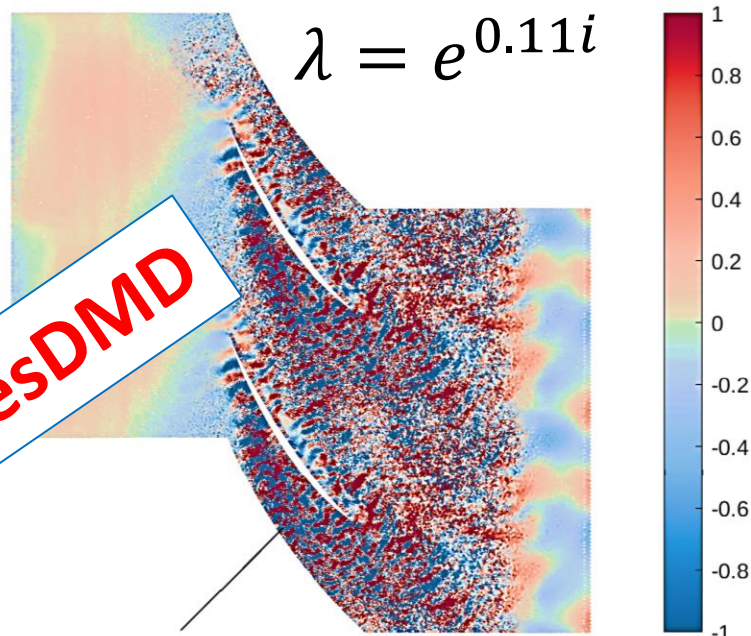
- Reynolds number  $\approx 3.9 \times 10^5$
- Ambient dimension ( $d$ )  $\approx 300,000$  (number of measurement points)

\*Raw measurements provided by Stephane Moreau (Sherbrooke)

Rel. Error  $\leq 0.0054$

$$\lambda = e^{0.11i}$$

ResDMD

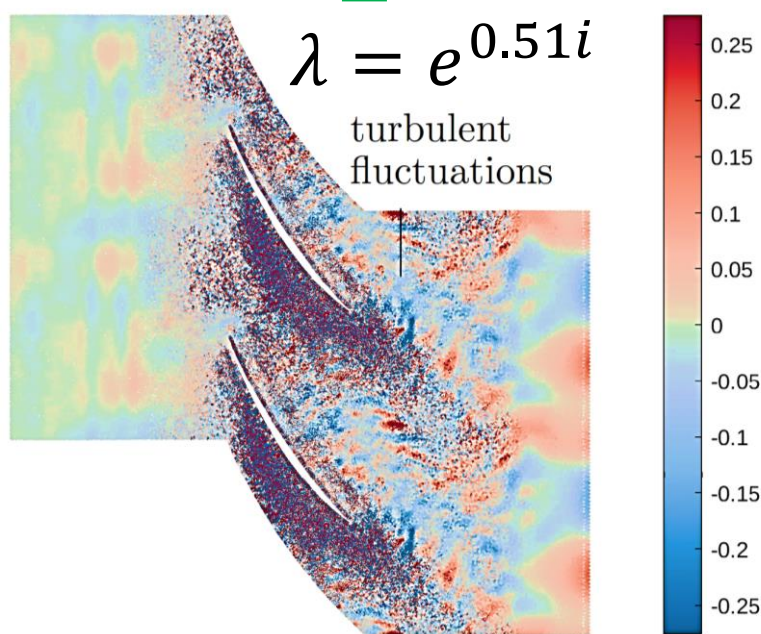


acoustic vibrations

Rel. Error  $\leq 0.0128$

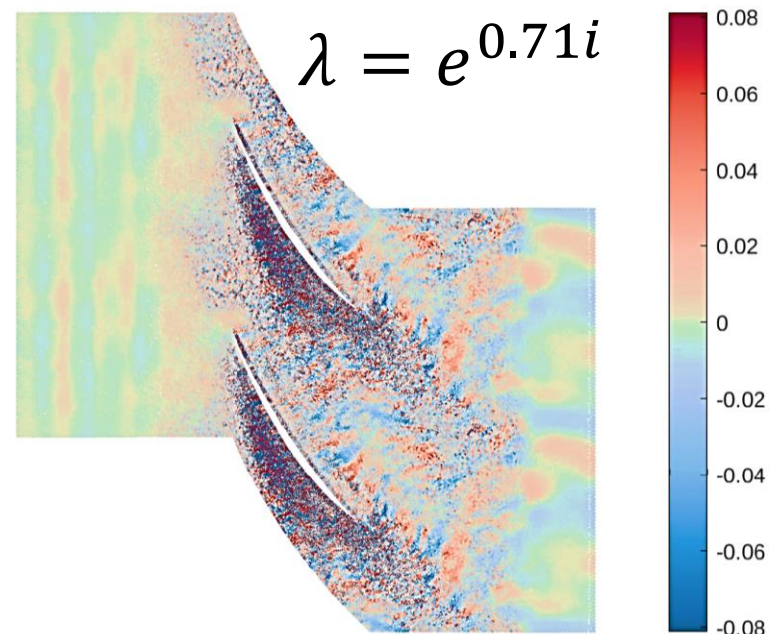
$$\lambda = e^{0.51i}$$

turbulent  
fluctuations



Rel. Error  $\leq 0.0196$

$$\lambda = e^{0.71i}$$





# Large $d$ ( $\Omega \subseteq \mathbb{R}^d$ ): robust and scalable

Popular to learn dictionary  $\{\psi_1, \dots, \psi_{N_K}\}$

E.g., DMD with truncated SVD (linear dictionary, most popular), kernel methods (this talk), neural networks, etc.

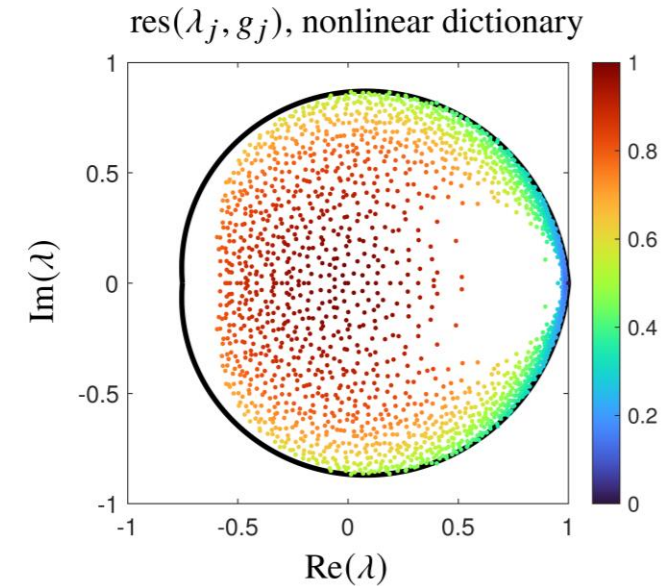
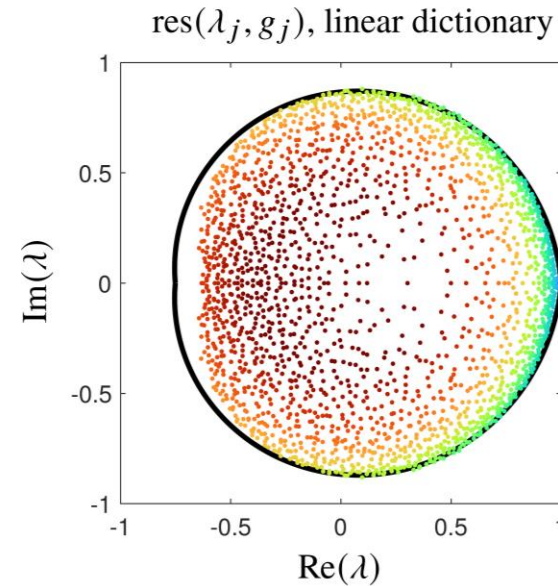
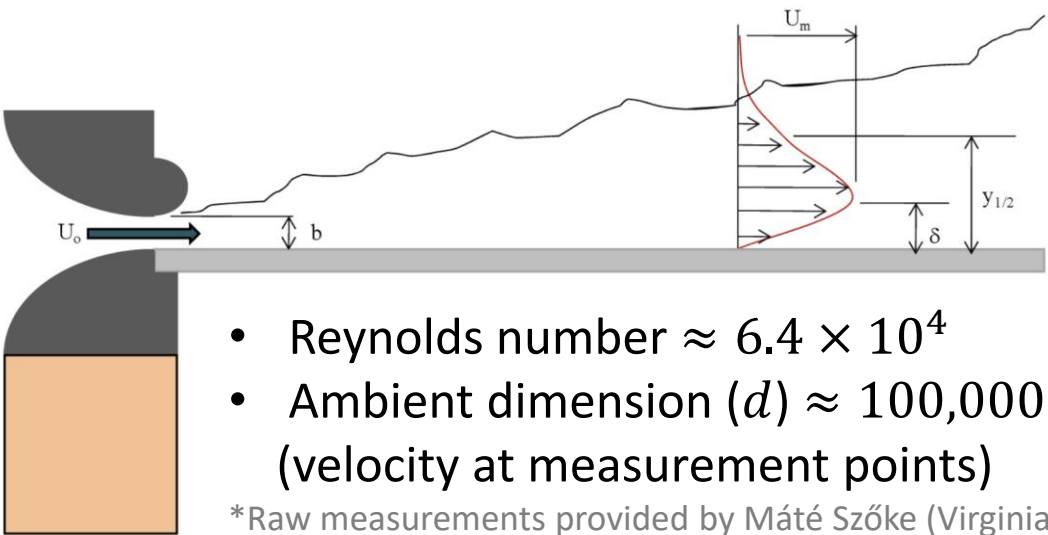
**Q: Is discretisation  $\text{span}\{\psi_1, \dots, \psi_{N_K}\}$  large/rich enough?**

## *Above algorithms:*

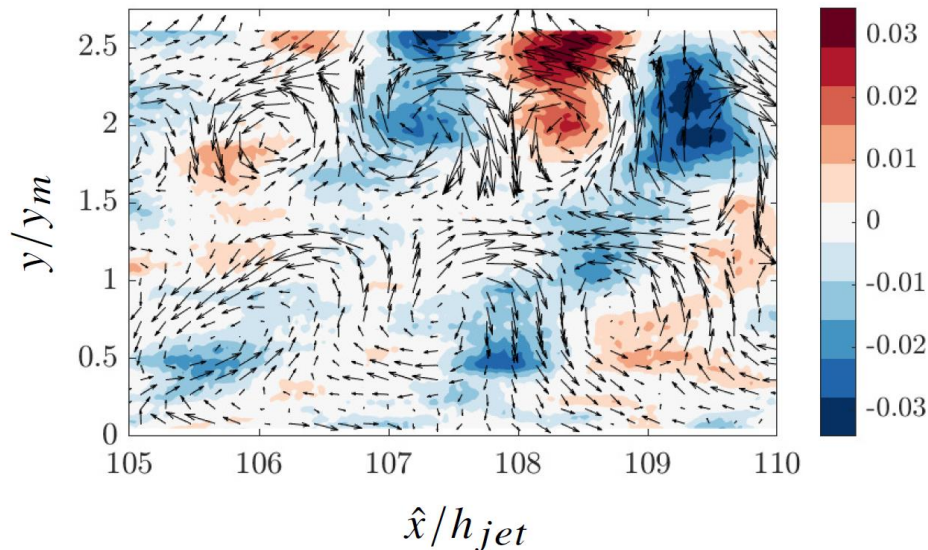
- Pseudospectra:  $\{z_k : \tau_k < \varepsilon\} \subseteq \text{Spec}_\varepsilon(\mathcal{K})$  **error control**
- Spectral measures:  $\mathcal{C}_g(z)$  and smoothed measures **adaptive check**

$\Rightarrow$  Rigorously **verify** learnt dictionary  $\{\psi_1, \dots, \psi_{N_K}\}$

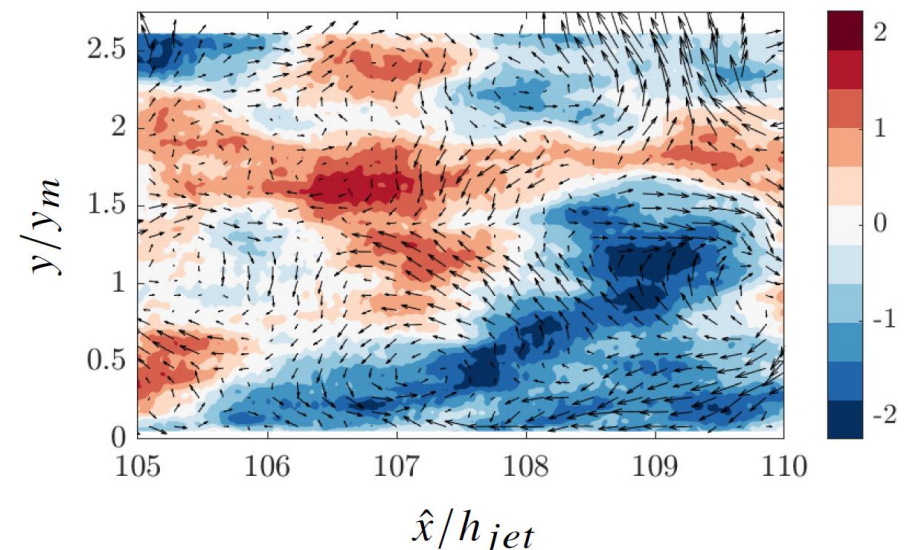
# Example: Verify the dictionary



$$\lambda = 0.9439 + 0.2458i, \text{ error} \leq 0.0765$$

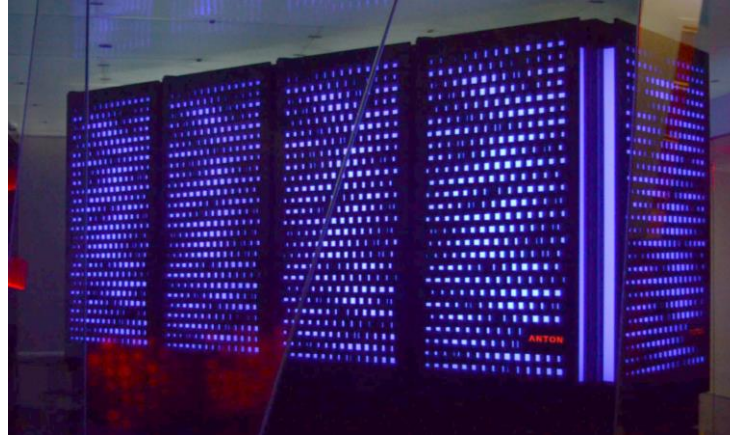
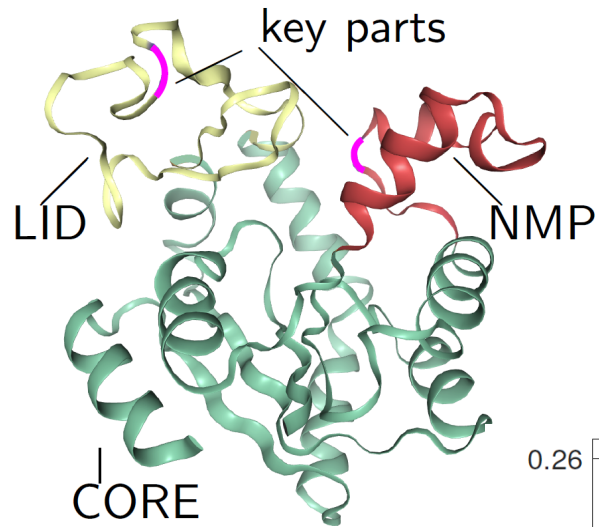


$$\lambda = 0.8948 + 0.1065i, \text{ error} \leq 0.1105$$



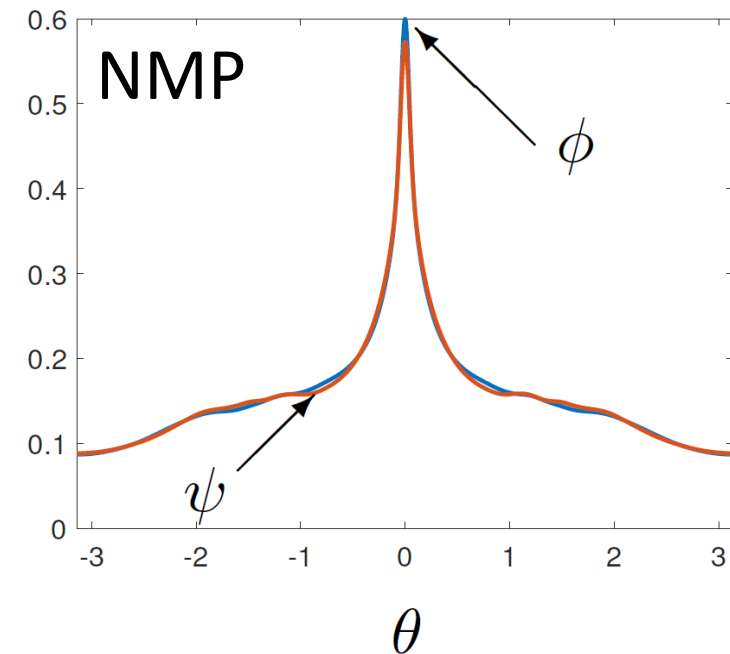
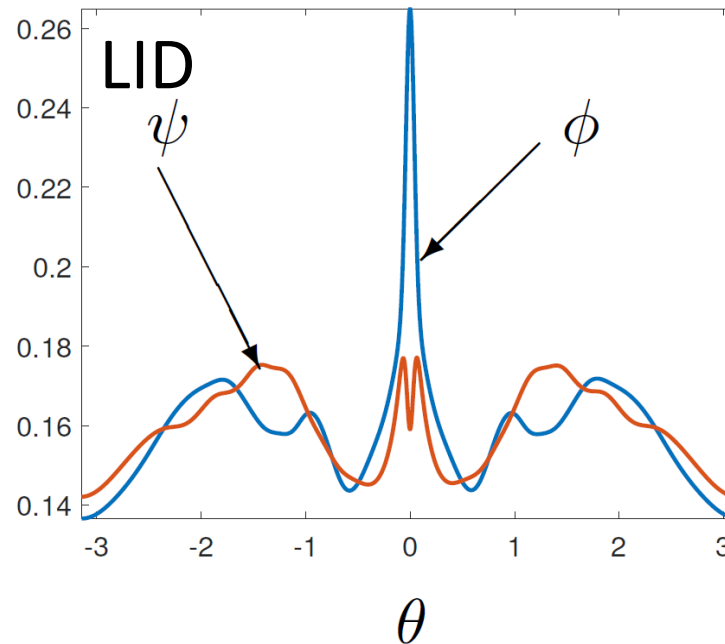
# Example: Spectral measures in large $d$

## Adenylate Kinase

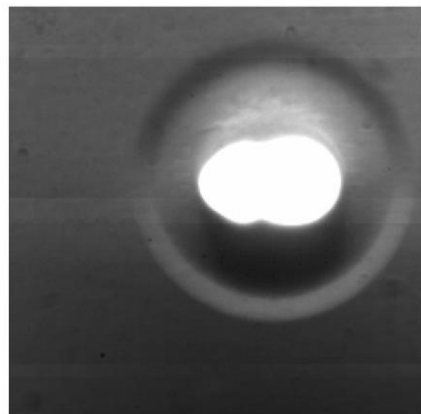
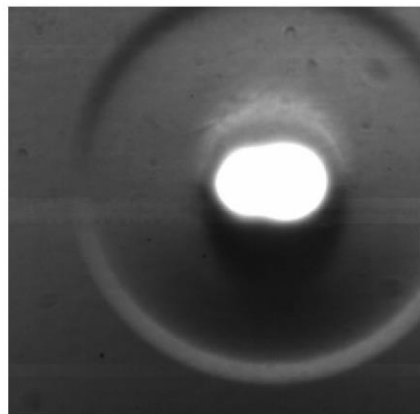
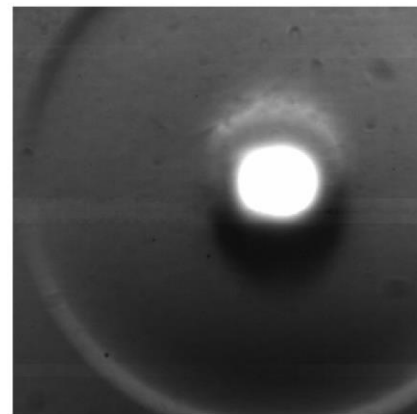
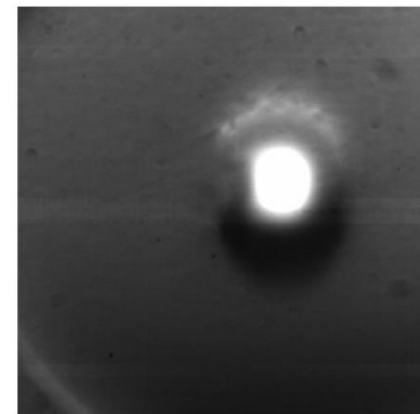
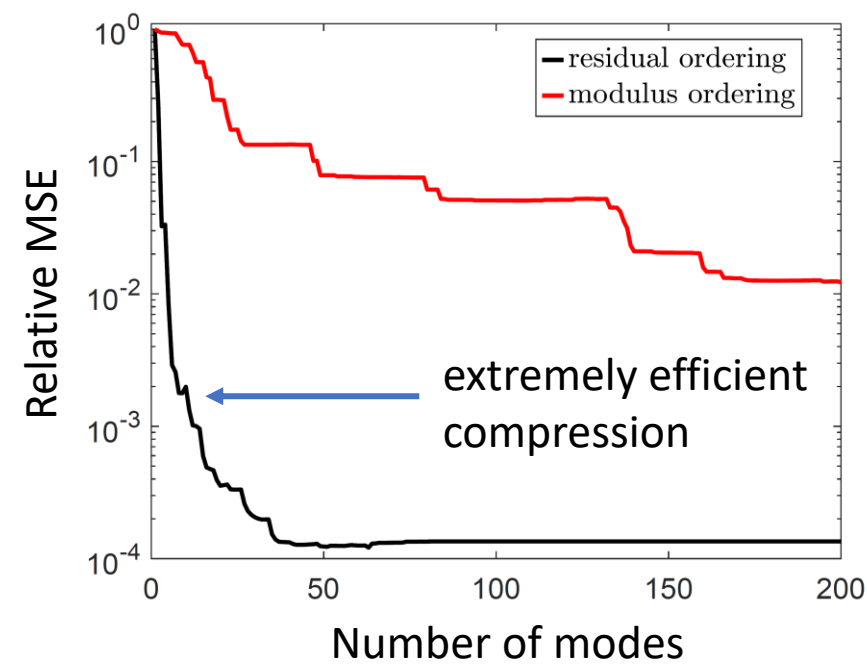
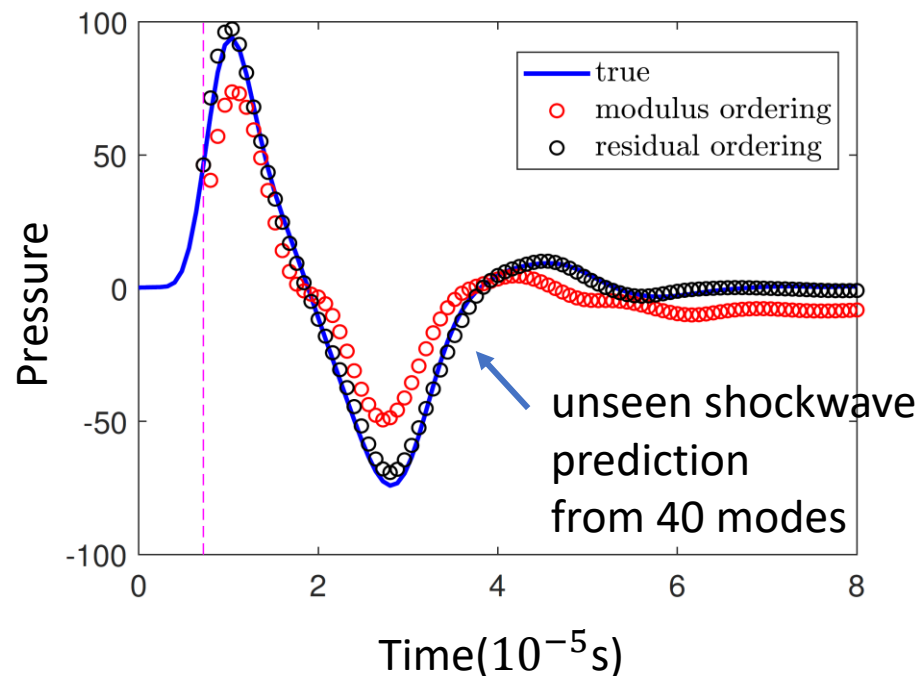


- Ambient dimension ( $d$ )  $\approx 20,000$  (positions and momenta of atoms)
- 6th order kernel (spec res  $10^{-6}$ )

\*Dataset: [www.mdanalysis.org/MDAnalysisData/adk\\_equilibrium.html](http://www.mdanalysis.org/MDAnalysisData/adk_equilibrium.html)



# Example: Trustworthy Koopman mode decomposition

a)  $t = 5 \mu\text{s}$ b)  $t = 10 \mu\text{s}$ c)  $t = 15 \mu\text{s}$ d)  $t = 20 \mu\text{s}$ 

- C., Ayton, Szőke, "Residual Dynamic Mode Decomposition," **J. Fluid Mech.**, to appear.



# Wider programme

- Inf.-dim. computational analysis  $\Rightarrow$  **Compute spectral properties rigorously.**
- Continuous linear algebra  $\Rightarrow$  **Avoid the woes of discretization**
- Solvability Complexity Index hierarchy  $\Rightarrow$  **Classify diff. of comp. problems, prove algs are optimal.**
- **Extends to:** Foundations of AI, optimization, computer-assisted proofs, and PDE learning.

- 
- C., “On the computation of geometric features of spectra of linear operators on Hilbert spaces,” **Found. Comput. Math.**, to appear.
  - C., Horning, Townsend “Computing spectral measures of self-adjoint operators,” **SIAM Rev.**, 2021.
  - 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: On the barriers of deep learning and Smale’s 18th problem,” **Proc. Natl. Acad. Sci. USA**, 2022.
  - C., “Computing spectral measures and spectral types,” **Comm. Math. Phys.**, 2021.
  - C., Roman, Hansen, “How to compute spectra with error control,” **Phys. Rev. Lett.**, 2019.
  - C., “Computing semigroups with error control,” **SIAM J. Numer. Anal.**, 2022.
  - Boullé, Townsend, “Learning elliptic partial differential equations with randomized linear algebra”, **Found. Comput. Math.**, 2022.
  - Boullé, Kim, Shi, Townsend, “Learning Green’s functions associated with parabolic partial differential equations”, **JMLR**, to appear.
  - Gilles, Townsend, “Continuous analogues of Krylov methods for differential operators,” **SIAM J. Numer. Anal.**, 2019.
  - Horning, Townsend, “FEAST for Differential Eigenvalue Problems,” **SIAM J. Numer. Anal.**, 2020.
  - Ben-Artzi, C., Hansen, Nevanlinna, Seidel, “On the solvability complexity index hierarchy and towers of algorithms,” arXiv, 2020.
  - Smale, “The fundamental theorem of algebra and complexity theory,” **Bull. Amer. Math. Soc.**, 1981.
  - McMullen, “Families of rational maps and iterative root-finding algorithms,” **Ann. of Math.**, 1987.

# Summary: rigorous data-driven Koopmanism!

- “Too much” or “Too little”

**Idea:** New matrix for residual  $\Rightarrow$  **ResDMD** for computing spectra.

- Continuous spectra and spectral measures:

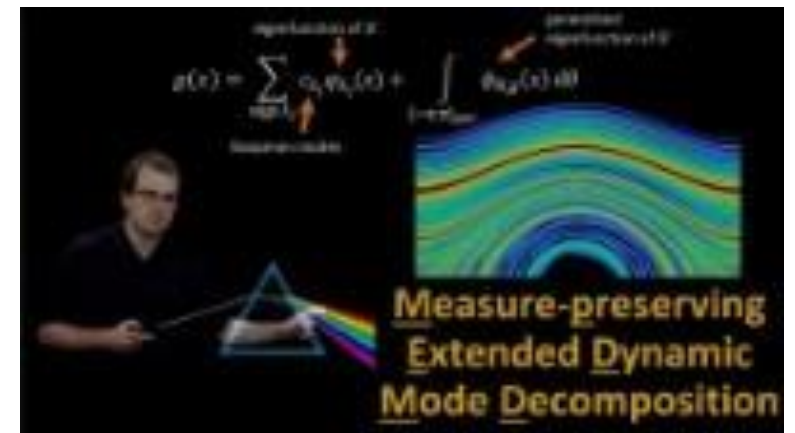
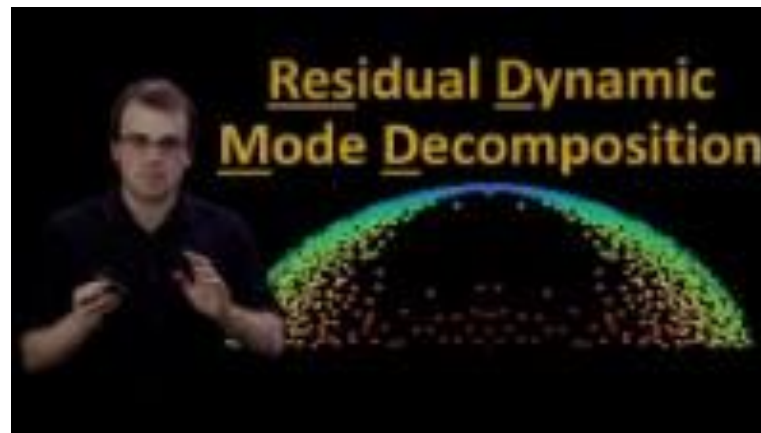
**Idea:** Convolution with rational kernels via resolvent and **ResDMD**.

- Is it right?

**Idea:** Use **ResDMD** to verify computations. E.g., learned dictionaries.

Short video summaries  
available on YouTube:

(Thanks to Steve Brunton  
for letting me use his channel!)

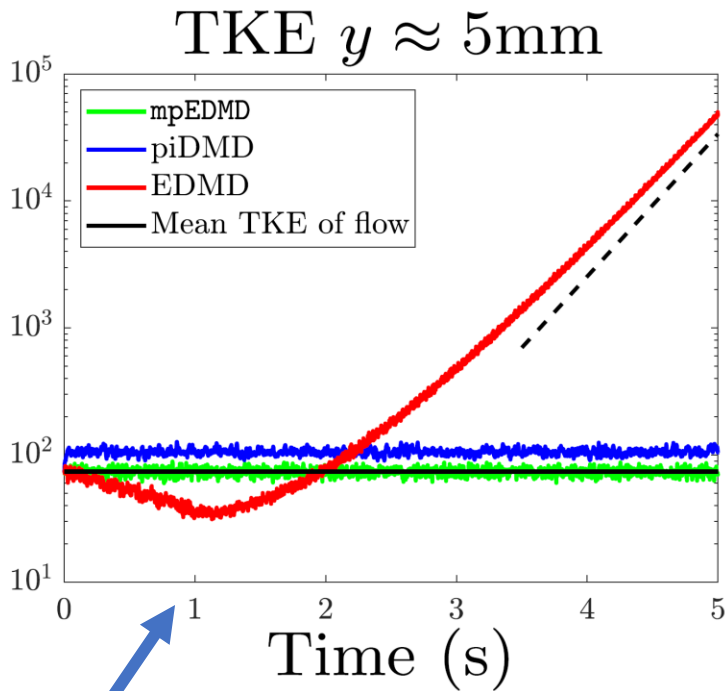


Code: <https://github.com/MColbrook/Residual-Dynamic-Mode-Decomposition>

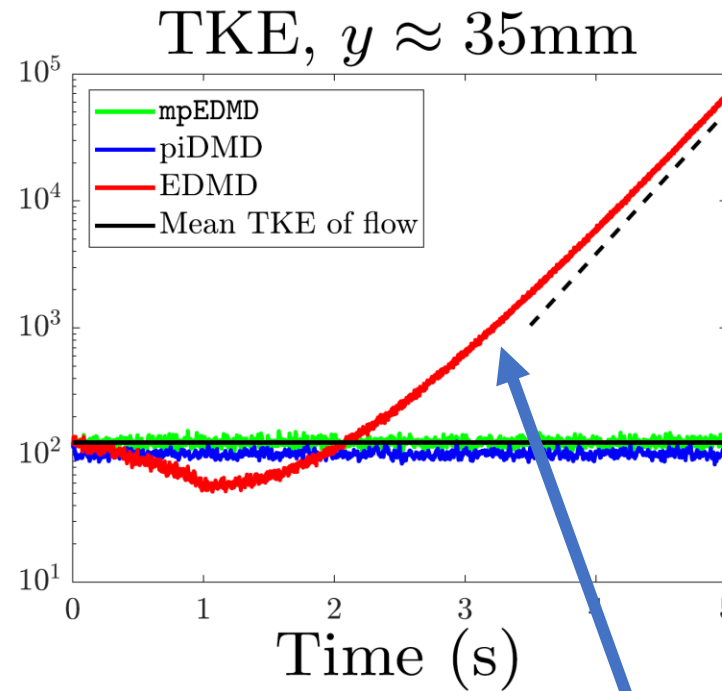
Additional slides...

# measure-preserving EDMD...

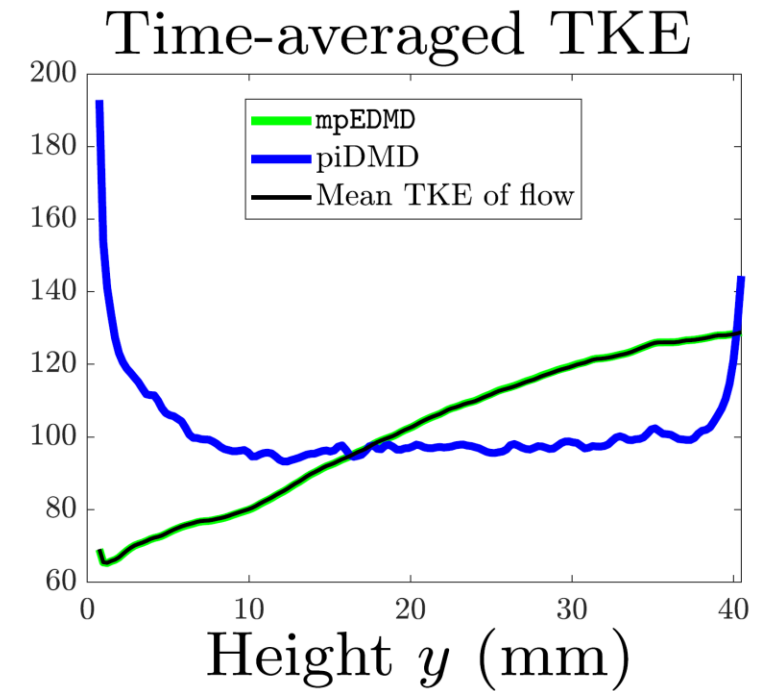
- Polar decomposition of  $\mathcal{K}$ . Easy to combine with any DMD-type method!
- Converges for spectral measures, spectra, Koopman mode decomposition.
- Measure-preserving discretization for arbitrary measure-preserving systems.



Snapshots collected over 1s




EDMD unstable!





# Solvability Complexity Index Hierarchy

Class  $\Omega \ni A$ , want to compute  $\Xi: \Omega \rightarrow (\mathcal{M}, d)$   metric space

- $\Delta_0$ : Problems solved in finite time (v. rare for cts problems).

- $\Delta_1$ : Problems solved in “one limit” with full error control:

$$d(\Gamma_n(A), \Xi(A)) \leq 2^{-n}$$

- $\Delta_2$ : Problems solved in “one limit”:

$$\lim_{n \rightarrow \infty} \Gamma_n(A) = \Xi(A)$$

- $\Delta_3$ : Problems solved in “two successive limits”:

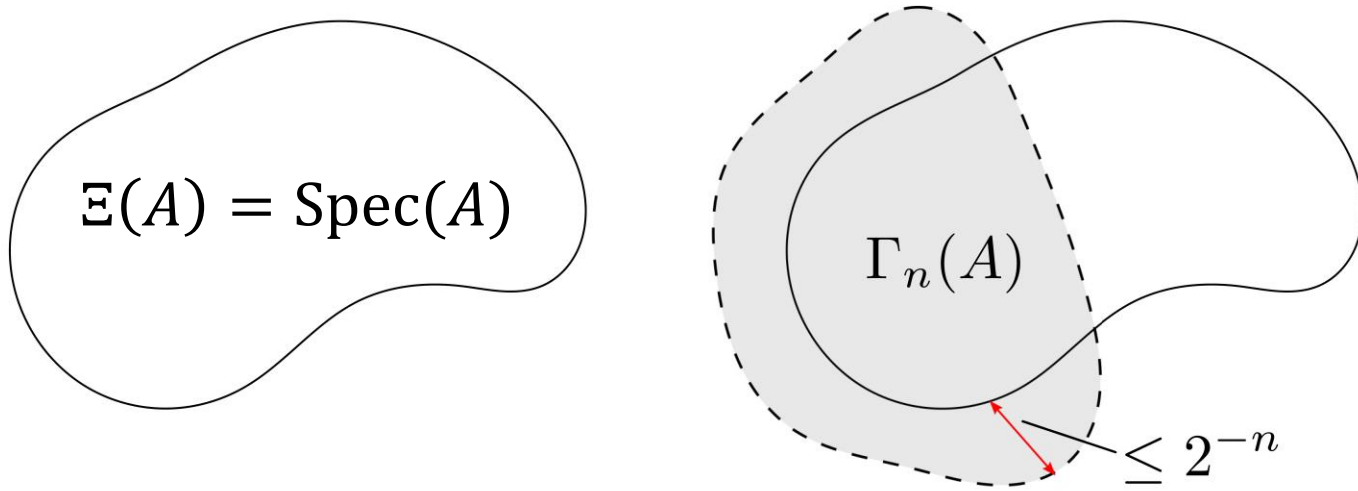
$$\lim_{n \rightarrow \infty} \lim_{m \rightarrow \infty} \Gamma_{n,m}(A) = \Xi(A)$$

⋮

- 
- Ben-Artzi, C., Hansen, Nevanlinna, Seidel, “*On the solvability complexity index hierarchy and towers of algorithms*,” preprint.
  - Hansen, “*On the solvability complexity index, the  $n$ -pseudospectrum and approximations of spectra of operators*,” **J. Amer. Math. Soc.**, 2011.
  - McMullen, “*Families of rational maps and iterative root-finding algorithms*,” **Ann. of Math.**, 1987.
  - Doyle, McMullen, “*Solving the quintic by iteration*,” **Acta Math.**, 1989.
  - Smale, “*The fundamental theorem of algebra and complexity theory*,” **Bull. Amer. Math. Soc.**, 1981.

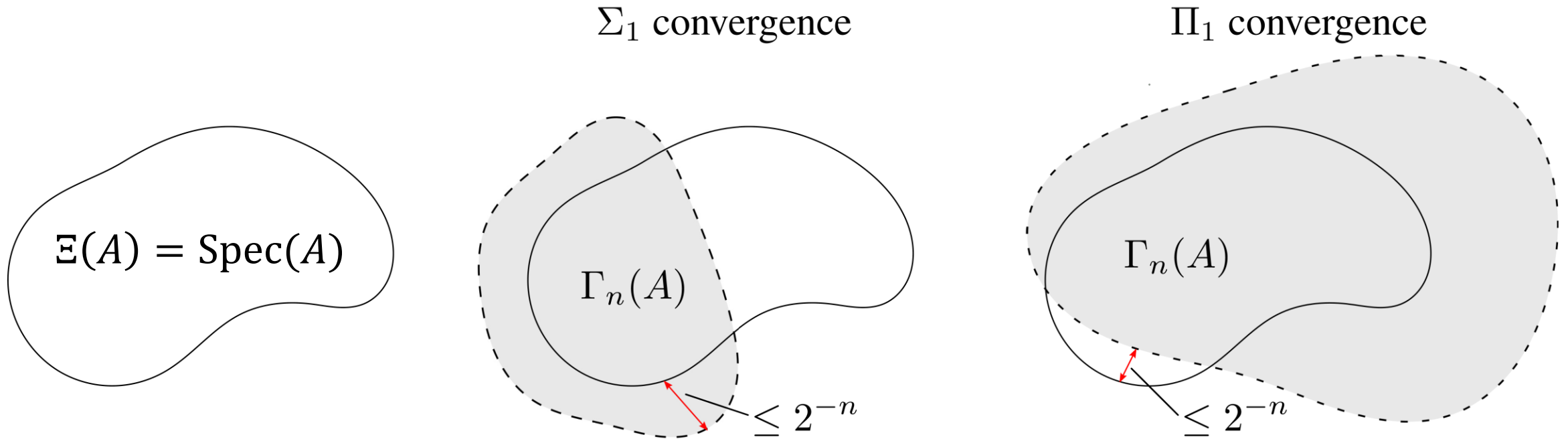
# Error control for spectral problems

$\Sigma_1$  convergence



- $\Sigma_1: \exists \text{ alg. } \{\Gamma_n\} \text{ s.t. } \lim_{n \rightarrow \infty} \Gamma_n(A) = \Xi(A), \max_{z \in \Gamma_n(A)} \text{dist}(z, \Xi(A)) \leq 2^{-n}$

# Error control for spectral problems



- $\Sigma_1: \exists$  alg.  $\{\Gamma_n\}$  s.t.  $\lim_{n \rightarrow \infty} \Gamma_n(A) = \Xi(A)$ ,  $\max_{z \in \Gamma_n(A)} \text{dist}(z, \Xi(A)) \leq 2^{-n}$
- $\Pi_1: \exists$  alg.  $\{\Gamma_n\}$  s.t.  $\lim_{n \rightarrow \infty} \Gamma_n(A) = \Xi(A)$ ,  $\max_{z \in \Xi(A)} \text{dist}(z, \Gamma_n(A)) \leq 2^{-n}$

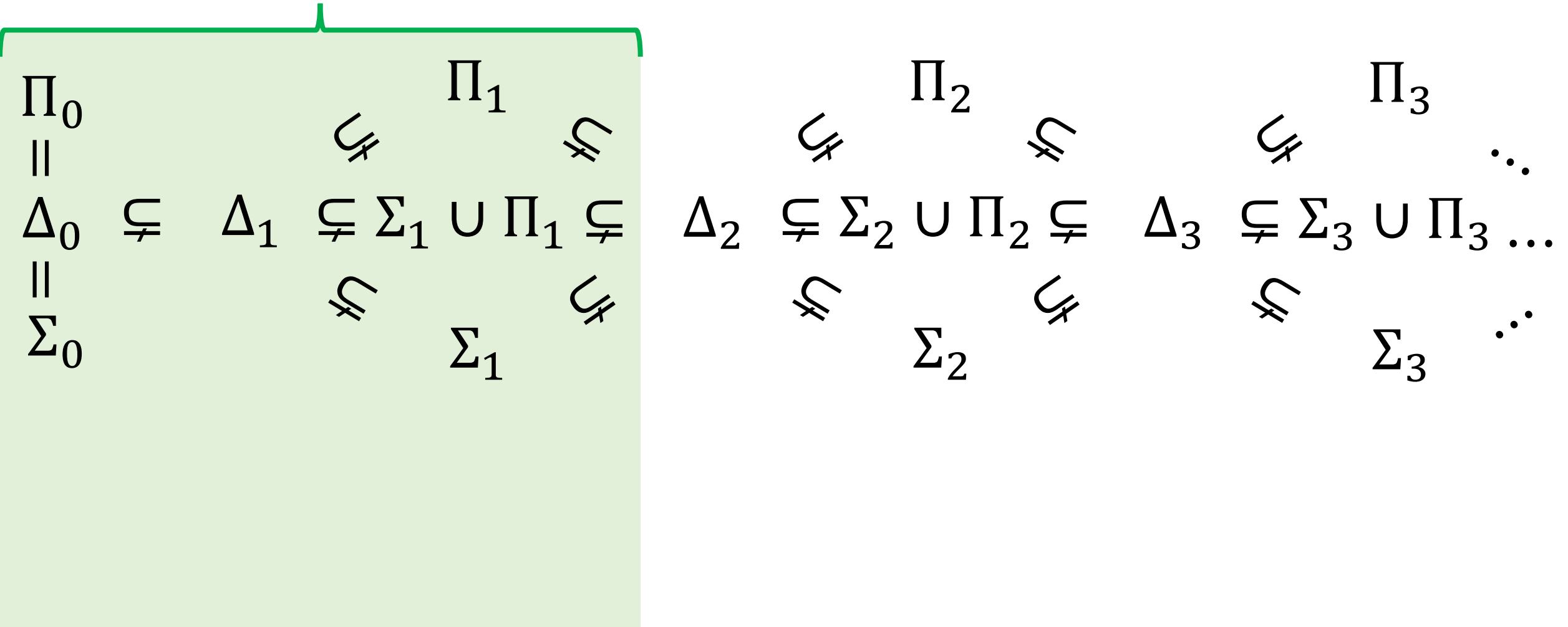
**Such problems can be used in a proof!**

# Small sample of classification theorems

Increasing difficulty



Error control

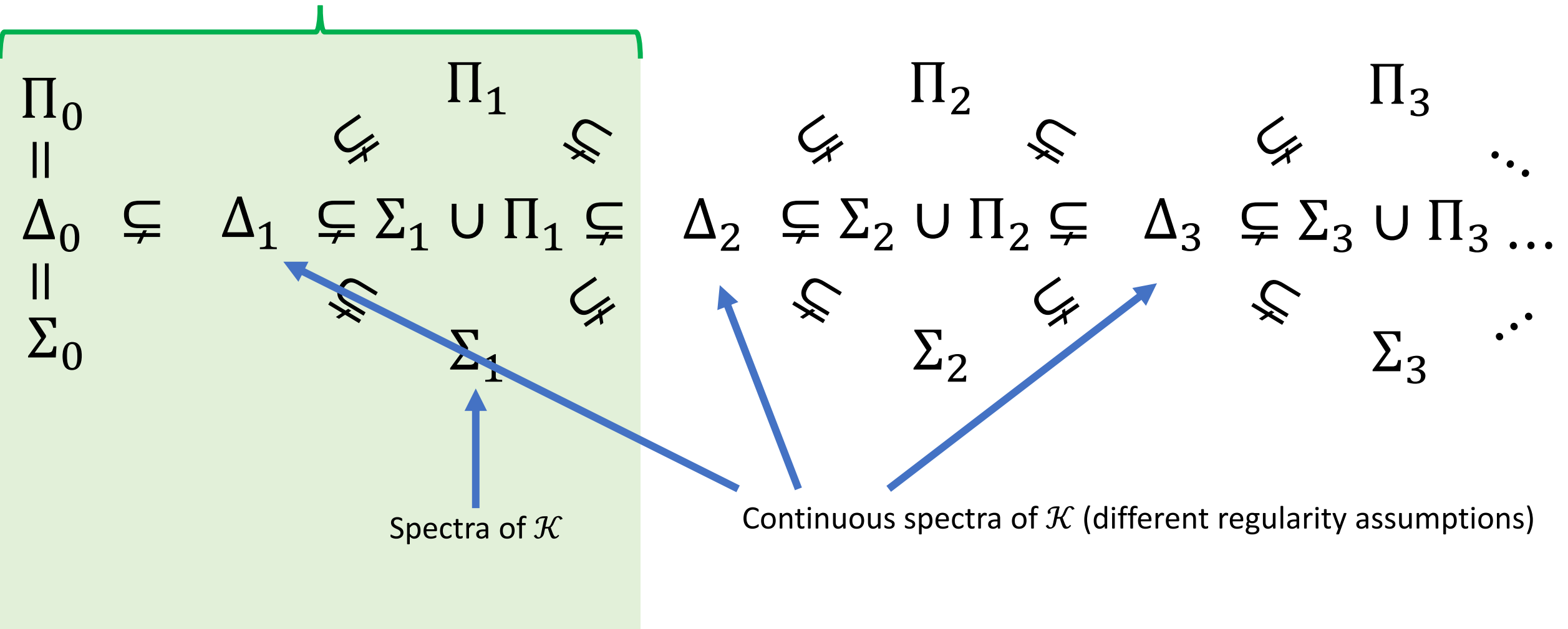


# Small sample of classification theorems

Increasing difficulty



Error control



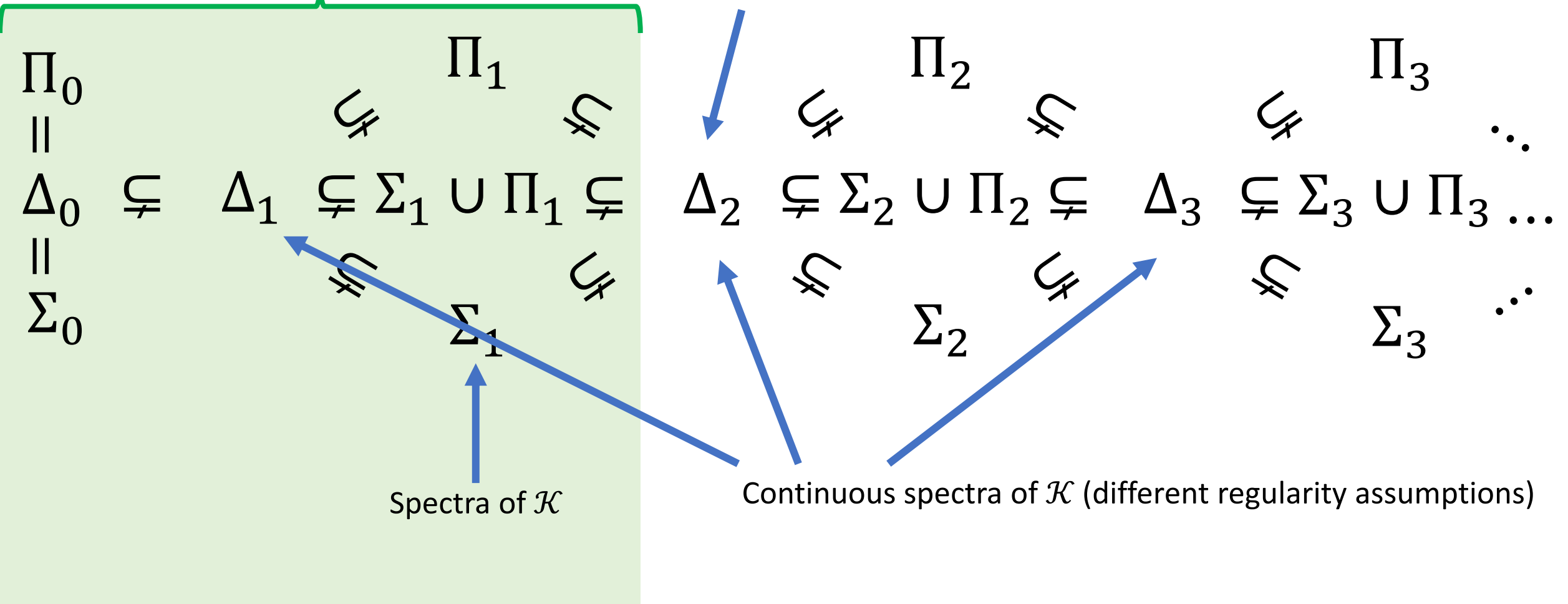
# Small sample of classification theorems

Increasing difficulty



Error control

Spectra of compact operators

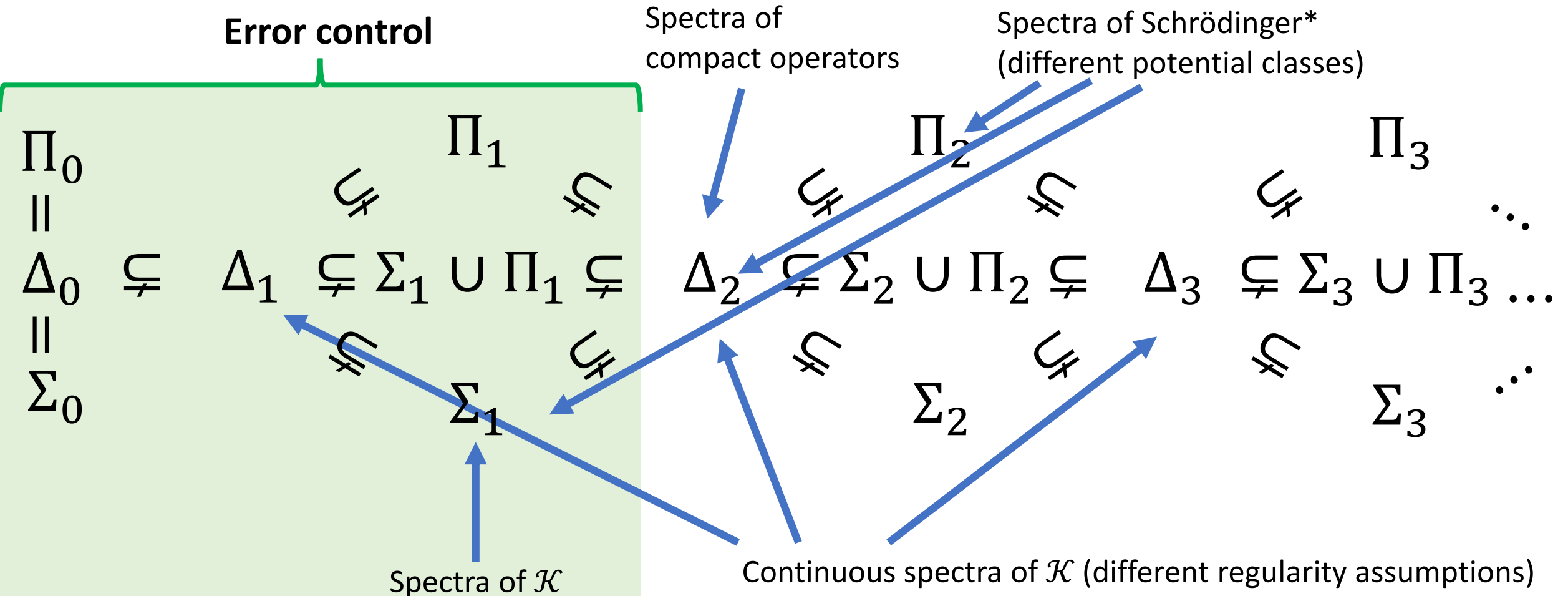


# Small sample of classification theorems

Increasing difficulty



Error control



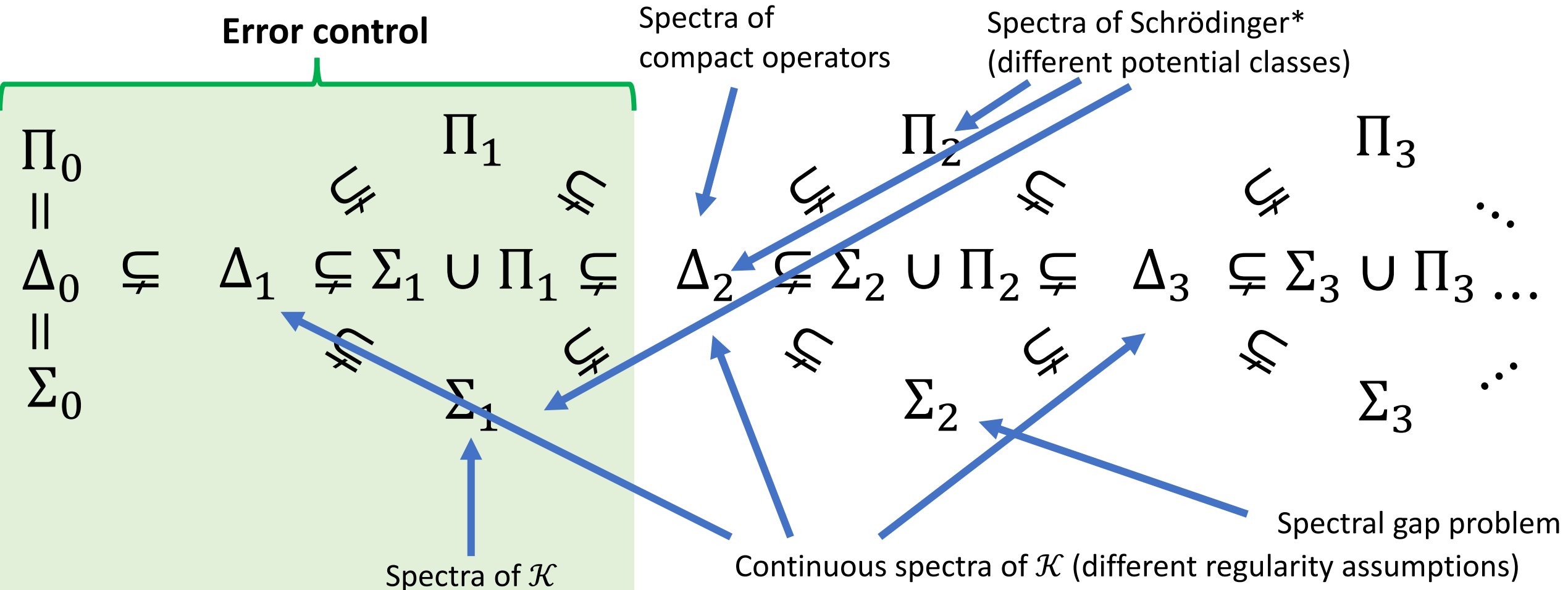
\*Open problem of Schwinger: "The special canonical group," "Unitary operator bases," PNAS, 1960.

# Small sample of classification theorems

Increasing difficulty



Error control

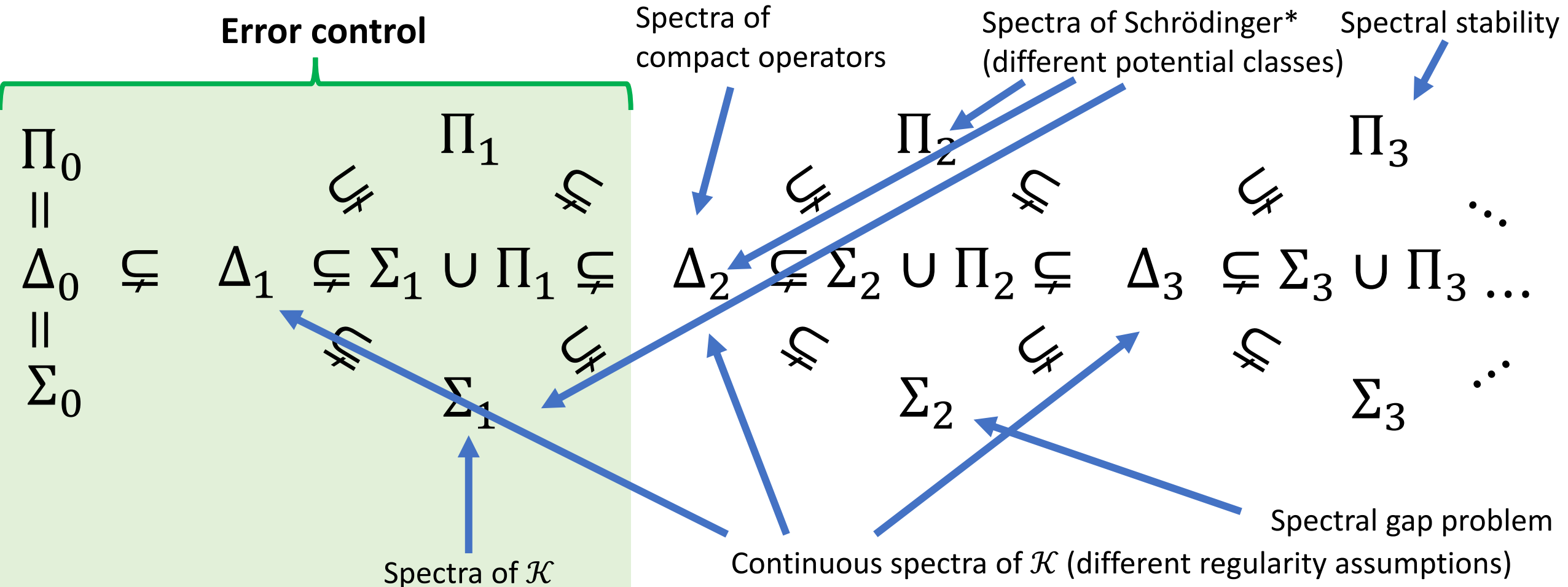


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# Small sample of classification theorems

Increasing difficulty



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