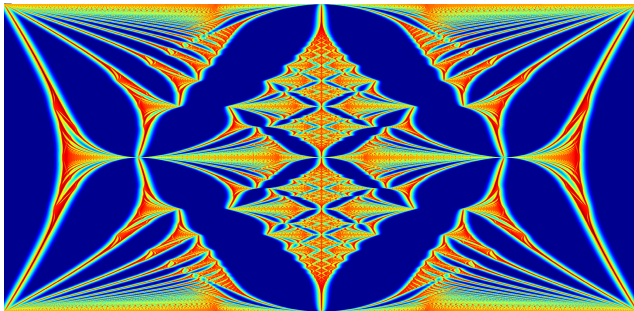


# Computing Spectral Measures

Matthew Colbrook, University of Cambridge

- Papers:** (i) C., "Computing spectral measures and spectral types", *Communications in Mathematical Physics* (2021)  
(ii) C., A. Horning, A. Townsend, "Computing spectral measures of self-adjoint operators", *SIAM Review* (to appear)

**Software:** SpecSolve available at <https://github.com/SpecSolve>



## Spectral Measures

**Finite-dimensional:**  $A \in \mathbb{C}^{n \times n}$  self-adjoint, o.n. basis of e-vectors  $\{v_j\}_{j=1}^n$

$$v = \left( \sum_{k=1}^n v_k v_k^* \right) v, \quad v \in \mathbb{C}^n \quad Av = \left( \sum_{k=1}^n \lambda_k v_k v_k^* \right) v, \quad v \in \mathbb{C}^n.$$

**Infinite-dimensional:** Self-adjoint operator  $\mathcal{L} : \mathcal{D}(\mathcal{L}) \rightarrow \mathcal{H}$  with spectrum

$$\Lambda(\mathcal{L}) = \{z \in \mathbb{C} : \mathcal{L} - z \text{ not bounded invertible}\}.$$

**Bad news:** Typically, no longer an o.n. basis of e-vectors.

**Spectral Theorem:** Projection-valued spectral measure  $\mathcal{E}$  (assigns an orthogonal projector to each Borel-measurable set) with

$$f = \left( \int_{\mathbb{R}} d\mathcal{E}(y) \right) f, \quad f \in \mathcal{H} \quad \mathcal{L}f = \left( \int_{\mathbb{R}} y d\mathcal{E}(y) \right) f, \quad f \in \mathcal{D}(\mathcal{L}).$$

**Intuition:** Diagonalises an infinite-dimensional operator.

**GOAL:** Compute (scalar versions of)  $\mathcal{E}$ .

# Motivation

**Scalar-valued measures (action of projections):**

$$\mu_f(\Omega) = \langle \mathcal{E}(\Omega)f, f \rangle$$

**Lebesgue decomposition theorem:**

$$d\mu_f(y) = \underbrace{\sum_{\lambda \in \Lambda^p} \langle \mathcal{P}_\lambda f, f \rangle \delta(y - \lambda) dy}_{\text{discrete part}} + \underbrace{\rho_f(y) dy + d\mu_f^{(\text{sc})}(y)}_{\text{continuous part}}.$$

**Crucial in:** quantum mechanics, scattering in particle physics, correlation in stochastic processes/signal-processing, fluid stability, resonances, density-of-states in materials science, orthogonal polynomials, random matrix theory, evolution PDEs,...

**Example:** in quantum mechanics,  $\mu_f$  describes the likelihood of different outcomes when the observable  $\mathcal{L}$  is measured. Can also solve SE

$$i \frac{df}{dt} = \mathcal{L}f, \quad f(0) = f_0, \quad \text{via} \quad f(t) = \left( \int_{\mathbb{R}} \exp(-ity) d\mathcal{E}(y) \right) f_0.$$

## A Hard Problem!

*"Most operators that arise in practice are not presented in a representation in which they are diagonalized... this raises the question of how to implement the methods of finite dimensional numerical linear algebra to compute the spectra of infinite dimensional operators. Unfortunately, there is a dearth of literature on this basic problem and, so far as we have been able to tell, there are no proven techniques."* W. Arveson, Berkeley (1994)

Some methods do exist, but treat cases with a lot of structure (e.g. compact perturbations of tridiagonal Toeplitz, some classes of singular Sturm–Liouville operators, etc.)

In contrast, want a **general** method to resolve spectral measures of  $\mathcal{L}$  (e.g. PDEs, integral operators, infinite matrices,...) and not an underlying discretisation or truncation.

finite-dimensional NLA  $\Rightarrow$  infinite-dimensional NLA



## Ideas from OPs: Computational Favard Theorem

$a_j, b_j \in \mathbb{R}$  and  $a_j > 0$ :

$$J = \begin{pmatrix} b_1 & a_1 & & \\ a_1 & b_2 & a_2 & \\ & a_2 & b_3 & \ddots \\ & & \ddots & \ddots \end{pmatrix}$$

OPs orthogonal w.r.t.  $\mu_J$ . For  $z = x + i\epsilon$ , define

$$G(z) := \langle (J - z)^{-1} \mathbf{e}_1, \mathbf{e}_1 \rangle = \int_{\mathbb{R}} \frac{d\mu_J(x)}{x - z} = \frac{1}{-z + b_1 - \frac{a_1^2}{-z + b_2 - \dots}}.$$

Then have weak convergence:

$$\lim_{\epsilon \downarrow 0} \frac{\operatorname{Im}(G(z))}{2\pi} = \mu_J.$$

# Ideas from Physics: Smoothed Measures

**Idea:** For  $z = x + i\epsilon$ , use

$$\mu_f^\epsilon(x) = \left\langle \frac{(\mathcal{L} - z)^{-1} - (\mathcal{L} - \bar{z})^{-1}}{2\pi i} f, f \right\rangle = \frac{1}{\pi} \int_{\Lambda(\mathcal{L})} \frac{\epsilon}{(x - \lambda)^2 + \epsilon^2} d\mu_f(\lambda).$$

Convolution with Poisson kernel: smoothed measure.

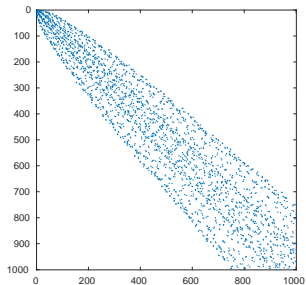
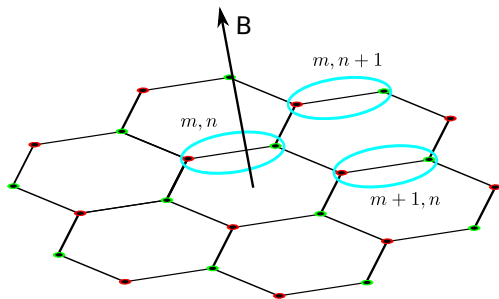
Converges weakly to measure as  $\epsilon \downarrow 0$ :

$$\int_{\mathbb{R}} \phi(y) \mu_f^\epsilon(y) dy \rightarrow \int_{\mathbb{R}} \phi(y) d\mu_f(y), \quad \text{as } \epsilon \downarrow 0,$$

for any bounded, continuous function  $\phi$ .

Approximate  $\mu_f^\epsilon$  via  $\mu_{f,N}^\epsilon$  ( $N = \text{truncation parameter}$ ).

# Numerical Balancing Act: Magnetic Graphene



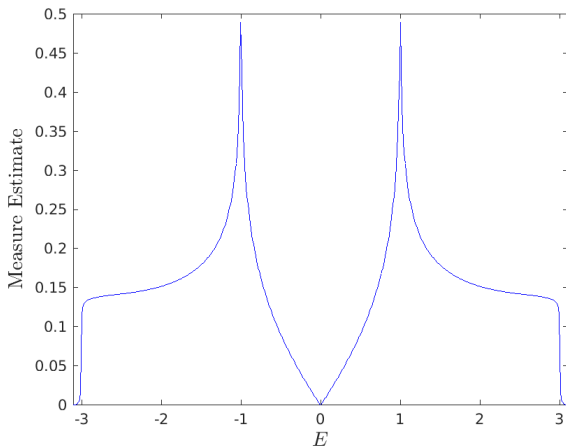
# Numerical Balancing Act: Magnetic Graphene

# Numerical Balancing Act: Magnetic Graphene

## Theorem (C. (2021))

*If we know rate of off-diagonal decay of infinite matrix, can compute measure in one limit. Extends to other operators such as PDEs.*

This is through a rectangular least squares type problem that computes  $(\mathcal{L} - z)^{-1}f$  with (asymptotic) error control.  $N(\epsilon)$  chosen **adaptively**.

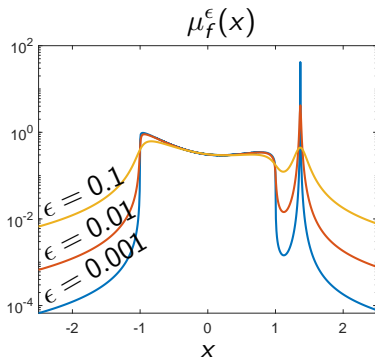


## Example: Integral Operator

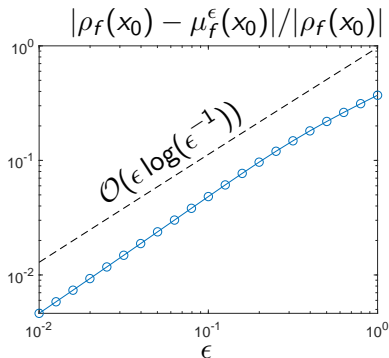
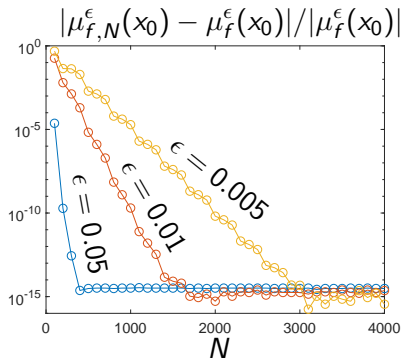
$$\mathcal{L}u(x) = xu(x) + \int_{-1}^1 e^{-(x^2+y^2)} u(y) dy, \quad x \in [-1, 1].$$

Discretise using adaptive Chebyshev collocation method.

Look at  $\mu_f$  with  $f(x) = \sqrt{3/2}x$ .



## Example: Integral Operator



$|\rho_f(x_0) - \mu_f^\epsilon(x_0)| = \mathcal{O}(\epsilon \log(\epsilon^{-1}))$  and need  $N \approx 20/\epsilon$ .

$\Rightarrow$  Infeasible to get more than five or six digits!

Q: Can we do better?



# Accelerating Convergence

Let  $m \in \mathbb{N}$ ,  $K \in L^1(\mathbb{R})$ . We say  $K$  is an  $m$ th order kernel if:

- (i) Normalized:  $\int_{\mathbb{R}} K(x) dx = 1$ ,
- (ii) Zero moments:  $K(x)x^j$  integrable,  $\int_{\mathbb{R}} K(x)x^j dx = 0$  for  $0 < j < m$ ,
- (iii) Decay at  $\pm\infty$ : There is a constant  $C_K$ , independent of  $x$ , such that

$$|K(x)| \leq C_K(1 + |x|)^{-(m+1)}, \quad x \in \mathbb{R}.$$

## Theorem (C., Horning, Townsend (2021))

If  $K$  is  $m$ th order,  $K_{\epsilon}(x) = \epsilon^{-1}K(x\epsilon^{-1})$  and  $\mu_f$  locally absolutely continuous near  $x_0$  with density  $\rho_f$  then

- Pointwise: If  $\rho_f$  locally  $\mathcal{C}^{n,\alpha}$  near  $x_0$  then

$$|[K_{\epsilon} * \mu_f](x_0) - \rho_f(x_0)| = \mathcal{O}(\epsilon^{n+\alpha}) + \mathcal{O}(\epsilon^m \log(\epsilon^{-1}))$$

- $L^p$ : If  $\rho_f$  locally  $\mathcal{W}^{n,p}$  near  $x_0$  ( $1 \leq p < \infty$ ) then

$$\|[K_{\epsilon} * \mu_f] - \rho_f\|_{L^p_{loc}} = \mathcal{O}(\epsilon^n) + \mathcal{O}(\epsilon^m \log(\epsilon^{-1}))$$

## Rational Kernels

**Idea:** Replace Poisson kernel with rational kernel

$$K(x) = \frac{1}{2\pi i} \sum_{j=1}^m \frac{\alpha_j}{x - a_j} - \frac{1}{2\pi i} \sum_{j=1}^m \frac{\beta_j}{x - b_j}.$$

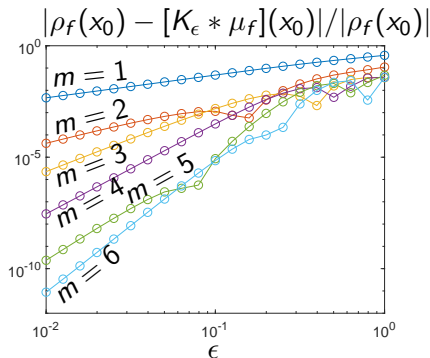
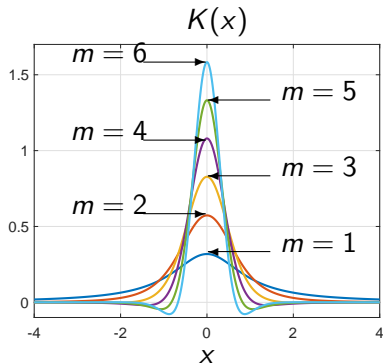
Can compute convolution with error control using resolvent

$$\begin{aligned} & [K_\epsilon * \mu_f](x) \\ &= \frac{-1}{2\pi i} \left[ \sum_{j=1}^m \alpha_j \langle (\mathcal{L} - (x - \epsilon a_j))^{-1} f, f \rangle - \sum_{j=1}^m \beta_j \langle (\mathcal{L} - (x - \epsilon b_j))^{-1} f, f \rangle \right]. \end{aligned}$$

Fix  $a_j$  in UHP,  $b_j$  in LHP  $\Rightarrow$  unique  $\{\alpha_j, \beta_j\}$  s.t.  $K$  an  $m$ th order kernel.

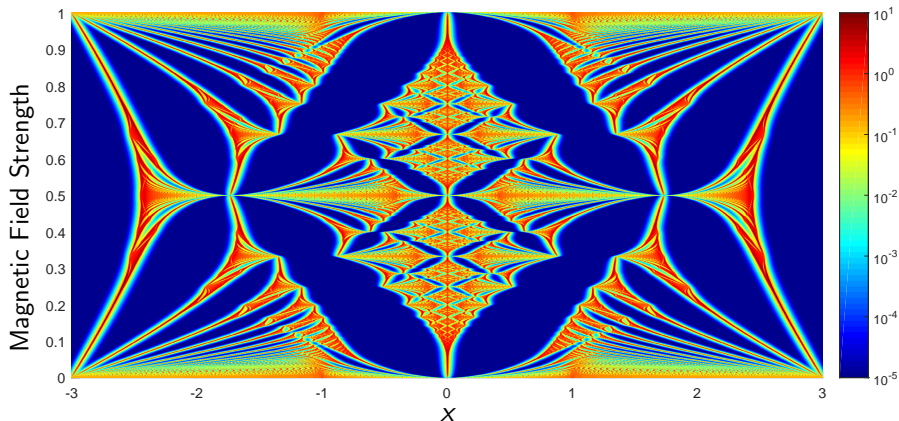
**NB:** At moment recommend  $\{a_j = \overline{b_j}\}$  equally spaced along  $\{\text{Im}(z) = 1\}$ .

# Integral Operator Revisited



See paper for general differential (even PDEs), integral and lattice operator examples - use sparse spectral methods for discretisation.

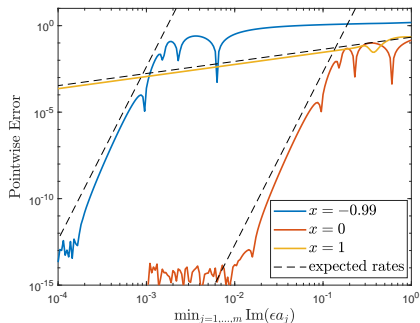
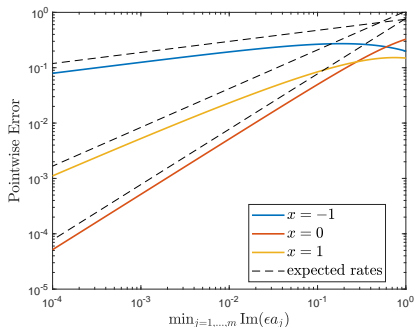
# Beautiful Fractal Structure!



Spectral measure of magnetic graphene, computed to high precision (see log scale) using  $m = 4$  kernel.

# OP Example: Jacobi Polynomials

$$d\mu_J = \frac{(1-x)^\alpha(1+x)^\beta}{N(\alpha,\beta)} dx = f_{\alpha,\beta}(x) dx,$$



Left: Pointwise errors for  $x = -1, 0, 1$  for  $m = 1$  and  $\alpha = 0.7$ ,  $\beta = 0.3$ . Right: Pointwise errors for  $x = -0.99, 0, 1$  for  $m = 10$  and  $\alpha = 0.7$ ,  $\beta = -0.3$ .

## ODEs Matlab Example

$$[\mathcal{L}u](x) = c_p(x) \frac{d^p u}{dx^p}(x) + \cdots + c_1(x) \frac{du}{dx}(x) + c_0(x)u(x), \quad p \geq 0,$$

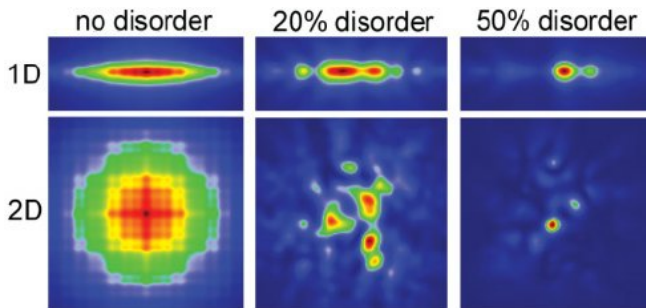
$$[\mathcal{L}u](x) = -\frac{d^2 u}{dx^2}(x) + \frac{x^2}{1+x^6}u(x), \quad f(x) = \sqrt{9/\pi} \cdot x^2/(1+x^6).$$

```
xi= linspace(0,5,101);           % Evaluation pts
f=@(x) x.^2./(1+x.^6)*sqrt(9/pi); % Measure wrt f(x)
c={@(x) x.^2./(1+x.^6),@(x) 0, @(x) -1}; % Schrodinger op
mu=diffMeas(c, f, xi, 0.1, 'order', 1); % epsilon=0.1, m=1
```

SpecSolve currently has capabilities for ODEs on real line & half-line, integral operators, and discrete operators.

# Physics Example: Background

Periodic systems have extended states (not localised), but add disorder...



Left: P. Anderson, **Nobel Prize in Phys. 1977** for discovering Anderson localisation. Right: Examples in 1D and 2D photonic lattices.

**What happens in aperiodic systems? Do we need disorder?**

# Bulk Localised Transport States: A new state for quasicrystals

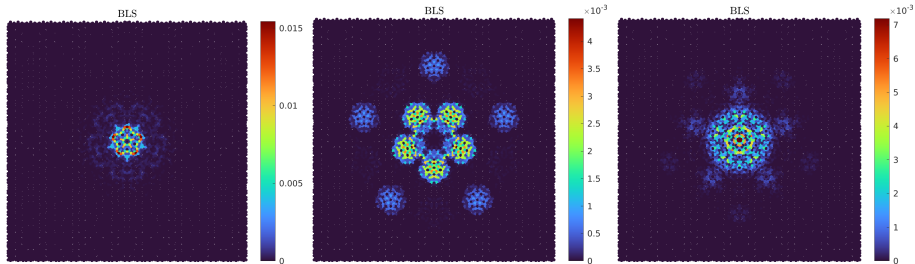
- **Bulk Localised Transport States (BLTs):** New states for magnetic quasicrystals
  - localised
  - “in-gap” (confirmed via comp. of inf-dim (topological) Chern numbers)
  - support transport
- **Cause (also confirmed with toy models):** Interplay of magnetic field with incommensurate areas of building blocks of quasicrystal.
- Not due to an internal edge, impurity or defect in the system.

⇒ NEW EXCITING PHYSICS!

D. Johnstone, M.J. Colbrook, A.E. Nielsen, P. Ohberg, C.W. Duncan.  
“Bulk Localised Transport States in Infinite and Finite Quasicrystals via Magnetic Aperiodicity.” *arXiv preprint*.



Transport: Error control allows us to be certain of this phenomenon.



## Example: Chern numbers

### Finite dimensions

$$\hat{P}^n = \sum_{m=1}^n |m\rangle \langle m|, \quad \hat{Q}^n = I - \hat{P}^n$$

$$\hat{x}^n = \hat{Q}^n \hat{x} \hat{P}^n, \quad \hat{y}^n = \hat{P}^n \hat{y} \hat{Q}^n$$

$$\mathcal{C}_i^n = -\frac{4\pi}{A_c^2} \text{Im} \{ \langle i | \hat{x}^n \hat{y}^n | i \rangle \}$$

### Infinite dimensions

$$\hat{P}^E = \int_{(-\infty, E]} d\mathcal{E}(\lambda)$$

$$\hat{P}_\epsilon^E = \int_{-\infty}^E [K_\epsilon * \mathcal{E}](\lambda) d\lambda, \quad \hat{Q}_\epsilon^E = I - \hat{P}_\epsilon^E$$

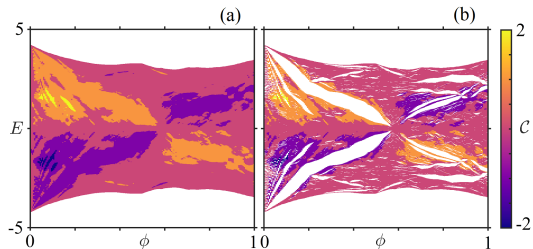
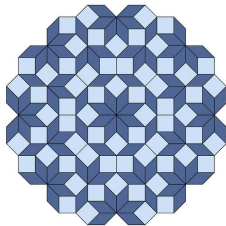
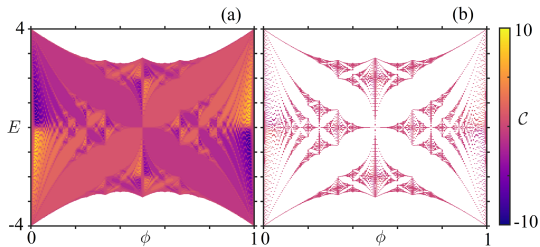
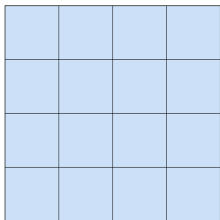
$$\hat{x}_\epsilon^E = \hat{Q}_\epsilon^E \hat{x} \hat{P}_\epsilon^E, \quad \hat{y}_\epsilon^E = \hat{P}_\epsilon^E \hat{y} \hat{Q}_\epsilon^E$$

$$\mathcal{C}_i^E = \frac{-4\pi}{A_c^2} \text{Im} \{ \langle i | \hat{x}_\epsilon^E \hat{y}_\epsilon^E | i \rangle \}$$

Round and take maximal count over site  $i$ .

Intuition: Topological index to detect in-gap (conducting) state.

# Example: Chern numbers



# Eigenvalue Hunting

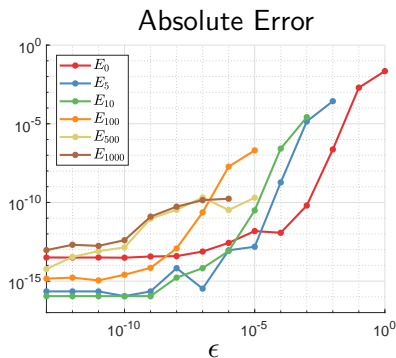
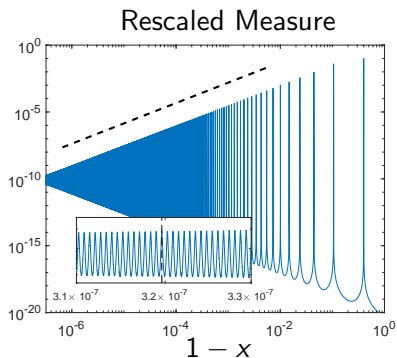
**Example:** Dirac operator.

- Describes the motion of a relativistic electron.
- Essential spectrum given by  $\mathbb{R} \setminus (-1, 1) \Rightarrow$  spectral pollution!
- Consider radially symmetric potential, coupled system on half-line:

$$\mathcal{D}_V = \begin{pmatrix} 1 + V(r) & -\frac{d}{dr} + \frac{\kappa}{r} \\ \frac{d}{dr} + \frac{\kappa}{r} & -1 + V(r) \end{pmatrix}.$$

- Map to  $[-1, 1]$  and solve shifted linear systems using sparse spectral methods.

# Eigenvalue Hunting



**NB:** Previous state-of-the-art achieves a few digits for a few excited states.

# Programme: Foundations of Infinite-Dimensional Spectral Computations

**Key Question: What is possible in infinite-dimensional NLA?**

**How:** Deal with operators directly, instead of previous 'truncate-then-solve'

⇒ Compute many spectral properties for the first time.

**Framework:** Classify problems in a computational hierarchy measuring their intrinsic difficulty and the optimality of algorithms.<sup>1</sup>

⇒ Algorithms that realise the boundaries of what computers can achieve.

**Have foundations for:** spectra with error control, spectral type (pure point, absolutely continuous, singularly continuous), Lebesgue measure and fractal dimensions of spectra, discrete spectra, essential spectra, eigenvectors + multiplicity, spectral radii, essential numerical ranges, geometric features of spectrum (e.g. capacity), spectral gap problem, ...

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<sup>1</sup>Holds regardless of model of computation (Turing, analog,...).

## Concluding Remarks

- **DIAGONALISATION:** General framework for computing spectral measures of self-adjoint operators.
- **Convolution with RATIONAL KERNELS:**
  - Can be evaluated using resolvent. ALL you need to be able to do is solve linear systems and compute inner products.
  - High-order kernels  $\Rightarrow$  high-order convergence.
  - Generalises to normal operators for local spectral regions on curves.
- Fast, local and parallelisable  $\Rightarrow$  State-of-the-art results for PDEs, integral operators and discrete operators.
- Forms part of a PROGRAMME for foundations of infinite-dimensional spectral computations.

**Ongoing and future work:** foundations of computational PDEs, foundations of (stable) neural networks, and computer-assisted proofs.

**Code:** <https://github.com/SpecSolve> (written with Andrew Horning).

## References

- M.J. Colbrook. “Computing spectral measures and spectral types.” *Communications in Mathematical Physics*, 2021.
- M.J. Colbrook, A. Horning, and A. Townsend. “Computing spectral measures of self-adjoint operators.” *SIAM Review*, to appear.
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- M.J. Colbrook, A.C. Hansen. “The foundations of spectral computations via the Solvability Complexity Index hierarchy: Part I.” *arXiv preprint*.
- M.J. Colbrook. “The foundations of spectral computations via the Solvability Complexity Index hierarchy: Part II.” *arXiv preprint*.

For further papers in this program and numerical code:

<http://www.damtp.cam.ac.uk/user/mjc249/home.html>

If you have further ideas or problems for collaboration, please get in touch!