Concentration in the Lanczos algorithm

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Joint work with Tom Trogdon

This material is based upon work supported by the National Science Foundation Graduate Research Fellowship Program under Grant No. DGE-1762114. Any opinions, findings, and conclusions or recommendations expressed in this material are those of the author and do not necessarily reflect the views of the National Science Foundation. We often want to evaluate $\mathbf{v}^{\mathsf{T}} f(\mathbf{A}) \mathbf{v}$ where $\mathbf{A} = \mathbf{U} \mathbf{A} \mathbf{U}^{\mathsf{T}}$ is a $n \times n$ symmetric matrix, \mathbf{v} is an arbitrary vector and f is a scalar function so that $f(\mathbf{A})$ is

 $f(\mathbf{A}) := \mathbf{U} f(\mathbf{\Lambda}) \mathbf{U}^{\mathsf{T}}.$

For instance, such expressions might arise in randomized algorithms for spectral sums since whenever $\mathbb{E}[\mathbf{v}\mathbf{v}^T] = \mathbf{I}$ we have

 $\mathbb{E}[\mathbf{v}^{\mathsf{T}}f(\mathbf{A})\mathbf{v}] = \mathrm{tr}(f(\mathbf{A})).$

A common approach to approximate $\mathbf{v}^T f(\mathbf{A})\mathbf{v}$ when \mathbf{A} is symmetric is via the Lanczos algorithm. Lanczos outputs an orthonormal basis \mathbf{Q} for Krylov subspace and a tridiagonal matrix \mathbf{T} giving the polynomial recurrence needed to construct this basis.

The Lanczos approximation is then defined as

$$\mathbf{v}^{\mathsf{T}} \mathbf{Q} f(\mathbf{T}) \mathbf{Q}^{\mathsf{T}} \mathbf{v} = \hat{\mathbf{e}}^{\mathsf{T}} f(\mathbf{T}) \hat{\mathbf{e}}$$

where $\hat{\mathbf{e}} = [1, 0, ..., 0]^{\mathsf{T}}$.

Empirical spectral measure (ESM):

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Weighted ESM:

$$\Psi[\mathbf{A},\mathbf{v}](x) = \sum_{i=1}^n ([\mathbf{U}]_{:,i}^\top \mathbf{v})^2 \mathbbm{1}[\lambda_i \leq x] = \mathbf{v}^\top \mathbbm{1}[\mathbf{A} \leq x] \mathbf{v}$$

Gaussian quadrature for μ defined as

$$[\mu]_k^{\rm gq}(x) = \sum_{i=1}^k \omega_i \mathbbm{1}[\theta_i \leq x]$$

where $\{\omega_i\}_{i=1}^k$ and $\{\theta_i\}_{i=1}^k$ are chosen so that μ and $[\mu]_k^{gq}$ share moments through degree 2k - 1.

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The k-point Gaussian quadrature rule $[\mu]_k^{gq}$ for μ is obtained from orthogonal polynomials of μ .

If $\mu = \Psi[\mathbf{A}, \mathbf{v}]$ then **T** from Lanczos gives upper left $k \times k$ block of Jacobi matrix for orthogonal polynomials.¹ Thus,

- Nodes are eigenvalues of ${\bf T}$
- Weights are squares of first components of eigenvectors of ${f T}$

Thus,

$$[\Psi[\mathbf{A},\mathbf{v}]]_k^{\mathrm{gq}} = \Psi[\mathbf{T},\hat{\mathbf{e}}]$$

The Lanczos approximation to the weighted CESM is itself a probability distribution function.

¹Golub and Meurant 2009.

It's not hard to see,

$$\mathbf{v}^{\mathsf{T}} f(\mathbf{A}) \mathbf{v} = \int f(x) \mathrm{d} \Psi[\mathbf{A}, \mathbf{v}](x), \qquad \quad \mathbf{\hat{e}}^{\mathsf{T}} f(\mathbf{T}) \mathbf{\hat{e}} = \int f(x) \mathrm{d} \Psi[\mathbf{A}, \mathbf{v}](x).$$

So to study Lanczos approximation to $\mathbf{v}^{\mathsf{T}} f(\mathbf{A}) \mathbf{v}$ we can just study Gaussian quadrature approximation of $\Psi[\mathbf{A}, \mathbf{v}]$.

We will run the Lanczos algorithm on $\mathbf{A}, \hat{\mathbf{e}}$ for k iterations to construct a Gaussian quadrature rule for $\Psi[\mathbf{A}, \hat{\mathbf{e}}]$ where $\mathbf{A} \sim \text{GOE}(n)$ and $\hat{\mathbf{e}} = [1, 0, ..., 0]^{\mathsf{T}}$.

To generate ${f A}$ can generate ${f X}$ with i.i.d. standard normal entries and then define

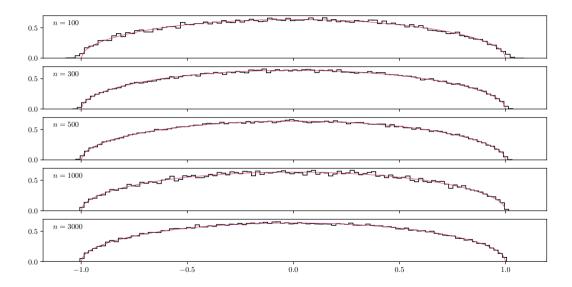
$$\mathbf{A} = \frac{\mathbf{X} + \mathbf{X}^{\mathsf{T}}}{2\sqrt{2n}}.$$

Equivalently, for $i \leq j$ entries of **A** are independent with distribution

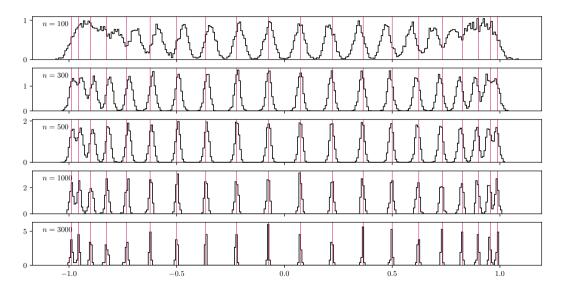
$$2\sqrt{2n}[\mathbf{A}]_{i,i} \sim \mathcal{N}(0,2), \qquad \qquad 2\sqrt{2n}[\mathbf{A}]_{i,j} \sim \mathcal{N}(0,1)$$

Note that A is unitarily invariant and the eigenvalues eventually lie between [-1,1] with high probability.

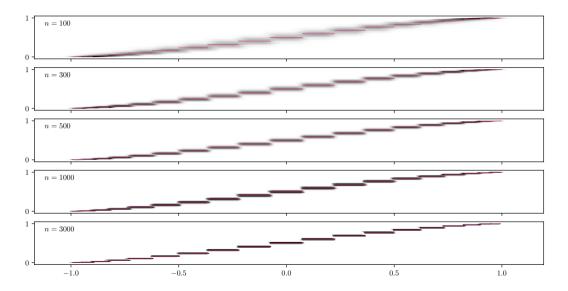
Weighted empirical spectral measure



Gaussian quadrature node (Ritz values)



Gaussian quadrature rule



As $n \to \infty$ we see "deterministic behavior"

- What is the limit?
- How fast does it converge?
- What do the fluctuations look like?

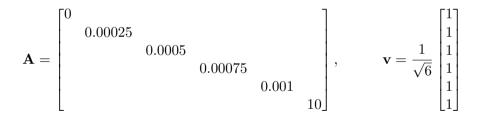
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These examples were computed in finite precision arithmetic without reorthogonalization

- isn't the Lanczos algorithm unstable?

In finite precision arithmetic, the Lanczos algorithm might behave extremely differently than in exact arithmetic.



Denote by \mathbf{T}, \mathbf{Q} the exact arithmetic output and $\tilde{\mathbf{T}}, \tilde{\mathbf{Q}}$ the finite precision output. How many digits of accuracy do we have for the following quantities:

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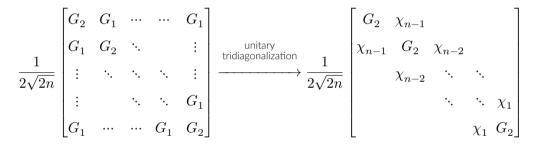
	$ ilde{\mathbf{Q}}-\mathbf{Q}$						$ ilde{\mathbf{T}}-\mathbf{T}$						$ ilde{\mathbf{Q}}^{ op} ilde{\mathbf{Q}} - \mathbf{I}$					
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Even for a very small example without any super extreme numbers, the Lanczos algorithm is not at all forward stable.

There is a lot of theory about Lanczos in finite precision (although no real forward analysis)³

³ Paige 1971; Paige 1976; Paige 1980; Grcar 1981; Simon 1982; Greenbaum 1989; Meurant 2006.

It is well known⁴ that GOE can be tridiagonalized:



The transform does not change the first entry of a vector so Lanczos on $\mathbf{A}, \hat{\mathbf{e}}$ will produce this tridiagonal matrix (in distribution).

⁴Trotter 1984; Dumitriu and Edelman 2002.

Let's look at the top-left $k \times k$ block as $n \to \infty$.

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$$\lim_{n \to \infty} \frac{1}{2\sqrt{2n}} \begin{bmatrix} G_2 & \chi_{n-1} & & & \\ \chi_{n-1} & G_2 & \chi_{n-2} & & \\ & \chi_{n-2} & \ddots & \ddots & \\ & & \ddots & \ddots & \chi_{n-k} \\ & & & \chi_{n-k} & G_2 \end{bmatrix} = \frac{1}{2} \begin{bmatrix} 0 & 1 & & & \\ 1 & 0 & 1 & & \\ & 1 & \ddots & \ddots & \\ & & \ddots & \ddots & 1 \\ & & & 1 & 0 \end{bmatrix}$$

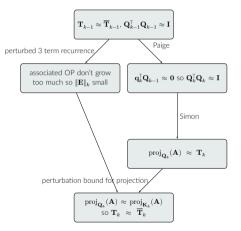
We can then analyze $\hat{\mathbf{e}}^{\mathsf{T}} f(\mathbf{T}) \hat{\mathbf{e}}$ using that

$$\hat{\mathbf{e}}^{\mathsf{T}} f(\mathbf{T}) \hat{\mathbf{e}} = \int f(x) \mathrm{d} \Psi[\mathbf{T}, \hat{\mathbf{e}}](x).$$

To do this we do a perturbation analysis for this integral based on perturbations of tridiagonal matrices.

Notation:

- $\mathbf{T}_k, \mathbf{Q}_k$ output of finite precision Lanczos
- $\,\overline{\mathbf{T}}_k$ limiting tridiagonal matrix
- $\mathbf{K}_k = [p_0(\mathbf{A})\mathbf{v}, \dots, p_{k-1}(\mathbf{A})\mathbf{v}]$ (these are polynomials of \mathbf{T}_k)
- $\mathbf{E}_k = \mathbf{Q}_k \mathbf{K}_k$ (can write in terms of associated polynomials of \mathbf{T}_k)



- For fixed k, the tridiagonal matrix output by the Lanczos algorithm run on a GOE matrix of size n concentrates rapidly as $n \to \infty$, and we can study the "average case" behavior of Lanczos as well as the fluctuations of Lanczos about this average case.
 - For any matrix and any ball of nonzero radius centered at this matrix, there is a non-zero probability of sampling a GOE matrix from within that ball
- We observe that Lanczos is (whp) forward stable for sufficiently large matrices⁵
 - We think we can prove this rigorously (probably need $\epsilon = O(1/n)$)
 - This would give (maybe first true) forward analysis result on Lanczos

⁵testing very big dense matrices is prohibitively expensive so we haven't done super big tests yet

Quote from Edelman and Rao^6 :

It is a mistake to link psychologically a random matrix with the intuitive notion of a 'typical' matrix or the vague concept of 'any old matrix'.

⁶Edelman and Rao 2005.

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