

Why is there only one definition of eigenvalue?

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Observation

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This is directly related to the fact that there is only one notion of graph quasirandomness.

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We usually define the eigenvalues of a matrix M to be the value λ so there is a vector v such that $Mv = \lambda v$. But this definition doesn't have a clear analog for a tensor: if v is a vector, Tv is a matrix, while if M is a matrix, $T \cdot M$ is a vector.

But there is an alternative definition of the eigenvalue of a matrix:
an eigenvalue is a local maximum of the inner product

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This definition does make sense: we can define a *vector eigenvalue*
of T to be a local maximum of

$$\lambda = \sum_{i,j,k} t_{ijk} v_i v_j v_k$$

over unit vectors v .

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That is, suppose M is a large, square, symmetric, dense matrix:

- M is an $n \times n$ matrix with n big,
- $m_{ij} = m_{ji}$,
- the entries $m_{ij} \in [0, 1]$,
- the average value $\sum_{i,j} m_{ij} = pn^2$.

Then the first eigenvalue measures the even distribution of M :

Lemma

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Suppose $\lambda_1 \approx np$. We can then look at λ_2 .

Suppose we generate M randomly: for each ij , we flip a fair coin independently to decide whether m_{ij} is 0 or 1. Then, with high probability, $\lambda_1 \approx \frac{n}{2}$ and λ_2 is close to 0 (and therefore all other eigenvalues are also close to 0).

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For most $ijkl$, these are four different pairs, so each quadruple has a $1/2^4$ chance of being present, so this sum is $\frac{1}{2^4} n^4$. If we think of M as the adjacency matrix of a graph, this amounts to counting the number of cycles of length 4.

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Since $\lambda_1 \approx \frac{n}{2}$, there is no room left for the other eigenvalues: $\sum_{1 < i \leq n} \lambda_i^4$ must be small.

Furthermore, this is, suitably interpreted, an equivalence:

Theorem

There is a language \mathcal{L} for probabilistic matrices such that, for each sentence σ and each $p \in (0, 1)$, if σ is (with high probability) true in the random matrix with density p then there is an ϵ so that σ holds in any matrix with n large enough, $\lambda_1 - np < \epsilon$, and $|\lambda_2| < \epsilon$.

For tensors, this doesn't work. λ_1 still measures even distribution:

Lemma

- $\lambda_1 \geq n^{3/2}p$, and
- $\lambda_1 = n^{3/2}p$ if and only if, for each $i \leq n$, $\sum_{j,k} a_{ijk} = n^2p$

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but consider the tensor generated as follows:

- Generate a random *matrix* $B = [b_{ij}]$ by flipping a coin for each pair i, j .
- Set $t_{ijk} = b_{ij} + b_{ik} + b_{jk} \pmod 2$.

Then T is evenly distributed and has a small λ_2 , but is very different from a random tensor.

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Then T is evenly distributed and has a small λ_2 , but is very different from a random tensor. For instance, in a random tensor we would have

$$\frac{1}{n^4} \sum_{ijkl} r_{ijk} r_{ijl} r_{ikl} r_{jkl} \approx 1/2^4$$

while

$$\frac{1}{n^4} \sum_{ijkl} t_{ijk} t_{ijl} t_{ikl} t_{jkl} \approx 1/2^3.$$

We could have defined a different kind of eigenvalue:

Definition

When T is a symmetric tensor, a *matrix eigenvalue* of T is a value λ which is a local maximum of

$$\sum_{i,j,k} t_{ijk} m_{ij} m_{ik} m_{jk}$$

among unit matrices M . We call such an M an *eigenmatrix* of T .

The first eigenvalue measures a stronger kind of even distribution:

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The second *matrix* eigenvalue does actually measure deviation from quasirandomness.

But there are other notions:

- eigenvalues are local maxima of $\sum_{i,j,k} t_{ijk} m_{ij} v_k$ among unit matrices M and unit vectors v (this is the notion related to Tao's *slice rank*),

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- eigenvalues are local maxima of $\sum_{i,j,k} t_{ijk} m_{ij} m_{ik}$ among unit matrices M
- etc.

Suppose we have a d -tensor. There is a notion of eigenvalue corresponding to every antichain of proper subsets of $\{1, \dots, d\}$, identifying antichains under permutations of $\{1, \dots, d\}$.

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So for 3-tensors, we have antichains like:

- $\{\{1\}, \{2\}, \{3\}\}$ corresponding to vector eigenvalues,
- $\{\{1, 2\}, \{1, 3\}, \{2, 3\}\}$ corresponding to matrix eigenvalues,
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- etc.

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The end.