

# Experimental Observation of Multifractality in Fibonacci Chains

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[arXiv:2207.13755](https://arxiv.org/abs/2207.13755)



# From the Fibonacci Sequence to Fibonacci Chains

Fibonacci numbers:

0, 1, 1, 2, 3, 5, 8, 13, 21, 34, 55, 89, ...

$$F_{n+1} = F_n + F_{n-1}$$

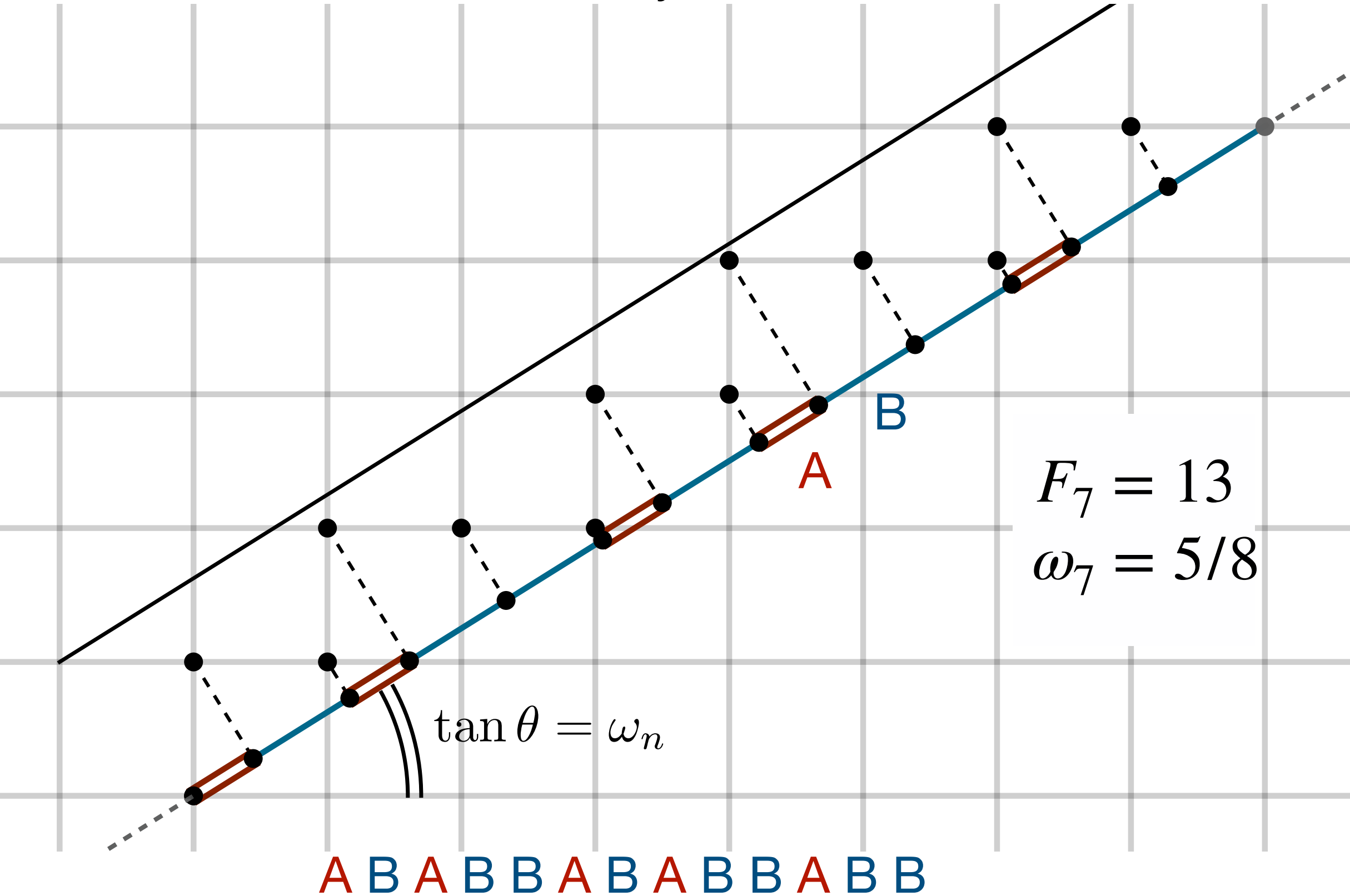
$$\omega_n = \frac{F_{n-2}}{F_{n-1}}$$

Fibonacci sequence:

$$\omega_n \xrightarrow{n \rightarrow \infty} \omega = \frac{\sqrt{5} - 1}{2}$$

*golden mean*

*Cut & Project*



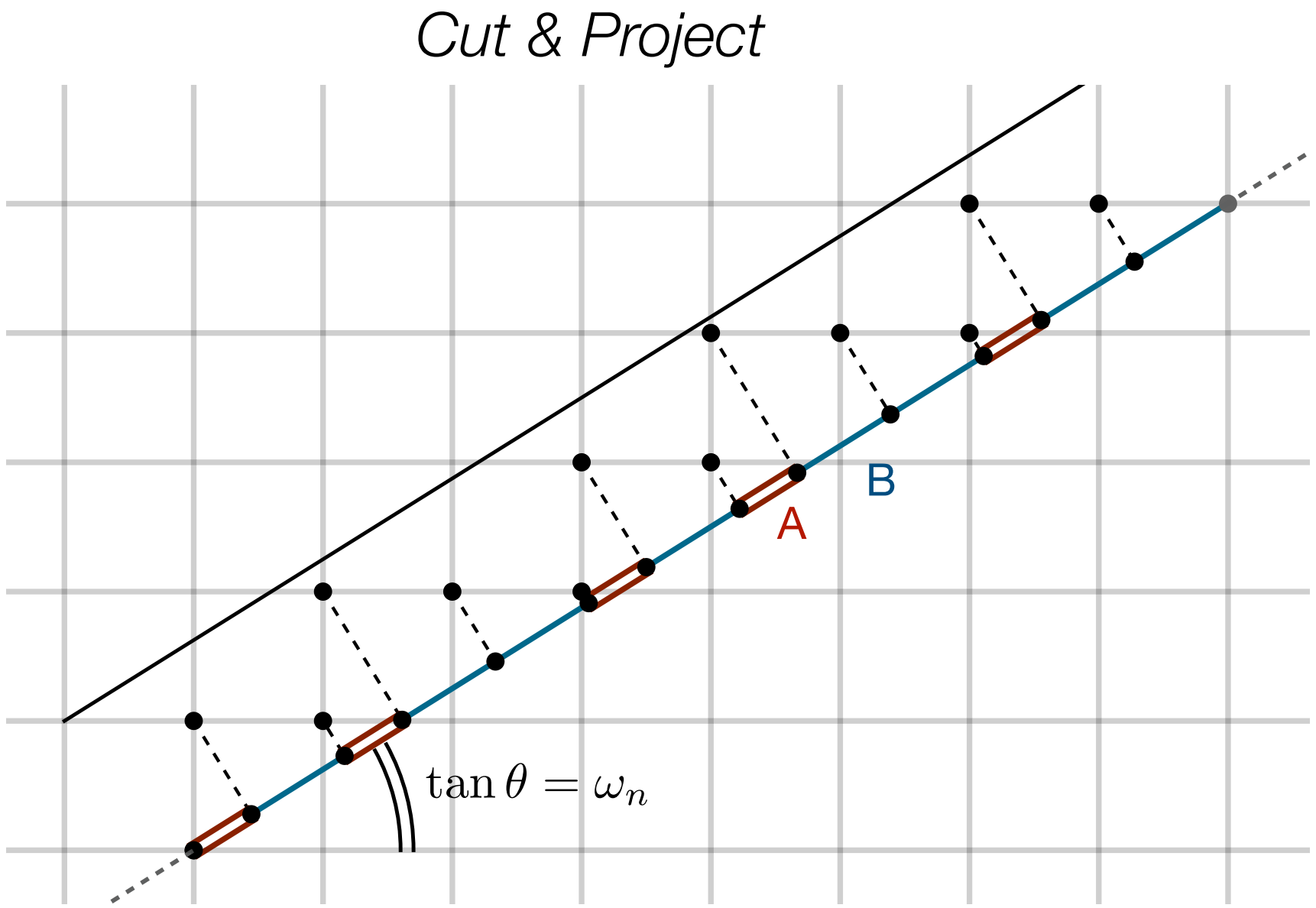
$F_7 = 13$   
 $\omega_7 = 5/8$

- If  $\tan \theta$  is rational
  - > periodic sequence with motif length  $F_n$
- $n \rightarrow \infty$ ,  $\tan \theta$  becomes irrational
  - > quasi-periodic sequence

# From the Fibonacci Sequence to Fibonacci Chains

## The Off-diagonal Fibonacci Model:

- Tight-Binding Hamiltonian  $H$  with constant on-site energies  $\nu_0$
- Assign coupling strength  $t_A$  to letter **A**,  $t_B$  to the letter **B**.
- Hopping constants follow the Fibonacci Sequence
- Important parameter  $\rho = t_B/t_A$

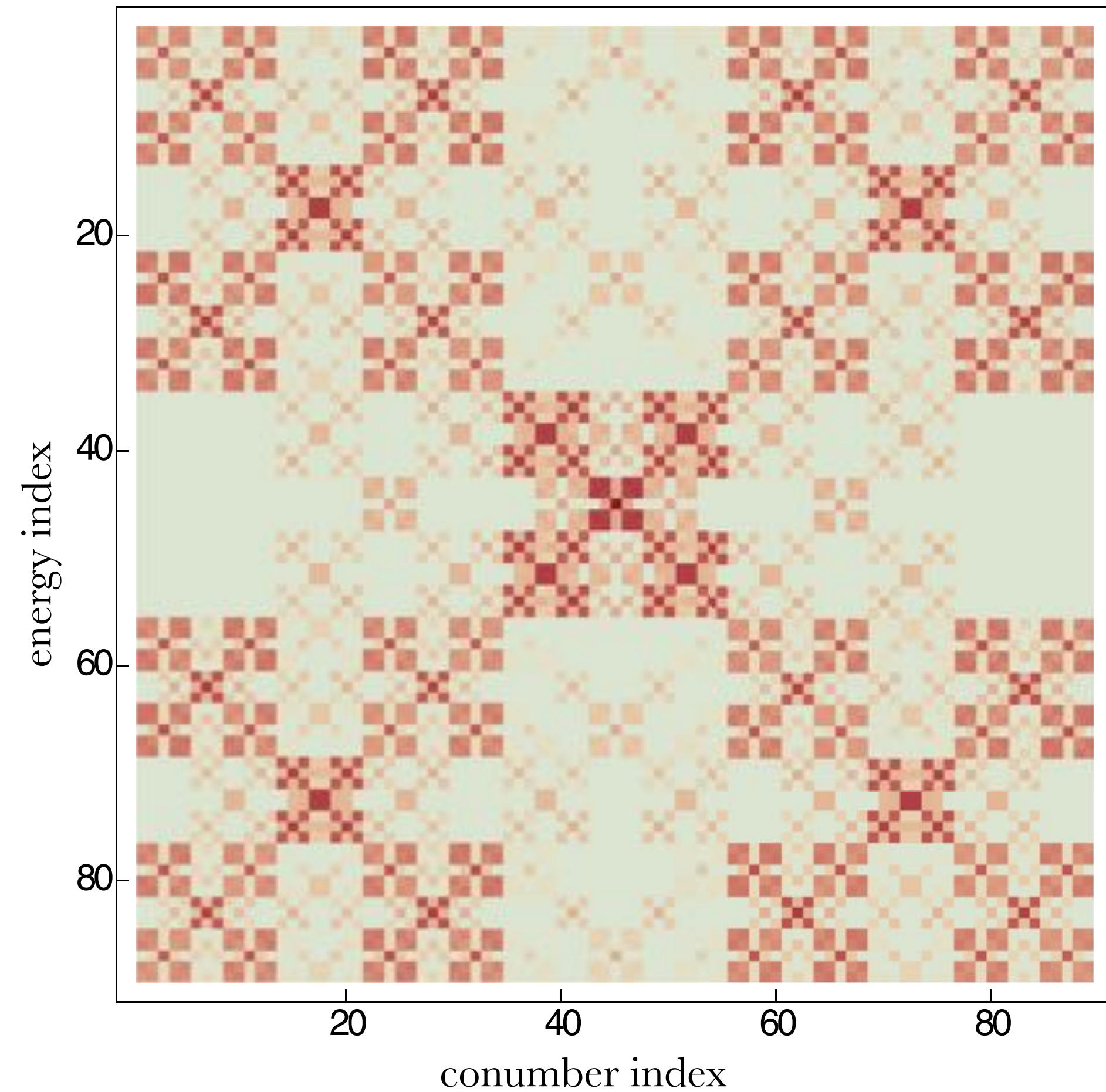


$$H = \begin{pmatrix} \nu_0 & t_A & 0 & 0 & 0 & 0 & 0 & \dots \\ t_A & \nu_0 & t_B & 0 & 0 & 0 & 0 & \dots \\ 0 & t_B & \nu_0 & t_A & 0 & 0 & 0 & \dots \\ 0 & 0 & t_A & \nu_0 & t_B & 0 & 0 & \dots \\ 0 & 0 & 0 & t_B & \nu_0 & t_B & 0 & \dots \\ 0 & 0 & 0 & 0 & t_B & \nu_0 & t_A & \dots \\ 0 & 0 & 0 & 0 & 0 & t_A & \nu_0 & \dots \\ \dots & \dots & \dots & \dots & \dots & \dots & \dots & \dots \end{pmatrix}$$

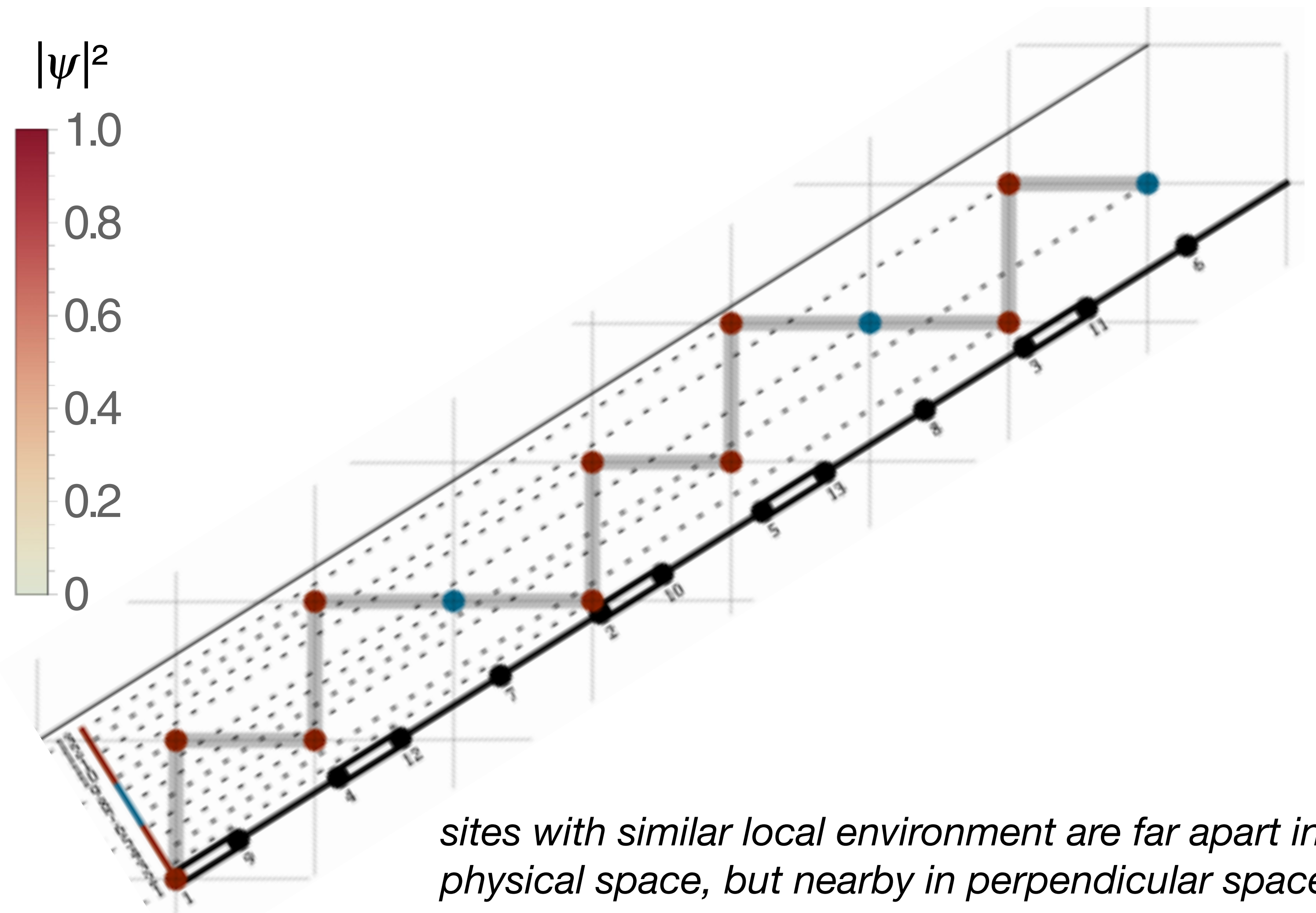
—> Check for Fractality in eigenvalues and eigenstates



# The Fibonacci Chain — a well known 1D Quasicrystal



Images taken from *Macé et al., PHYSICAL REVIEW B 93, 205153 (2016)*

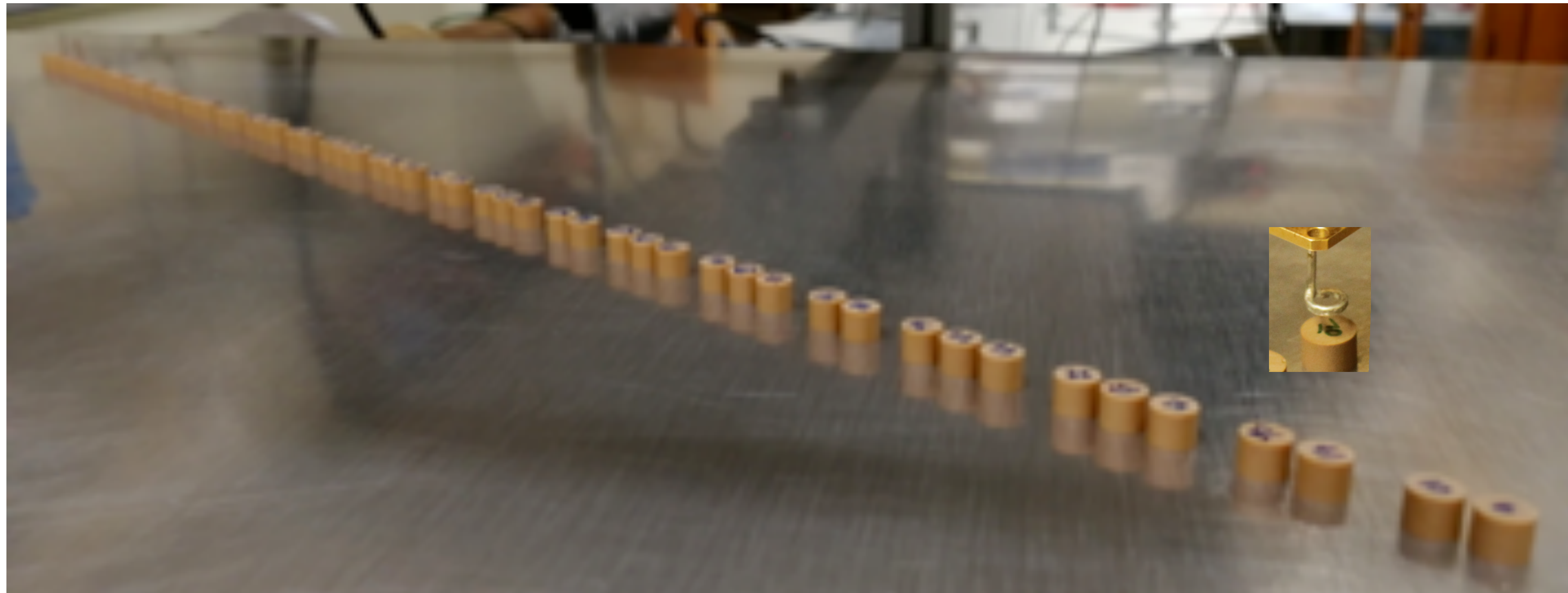
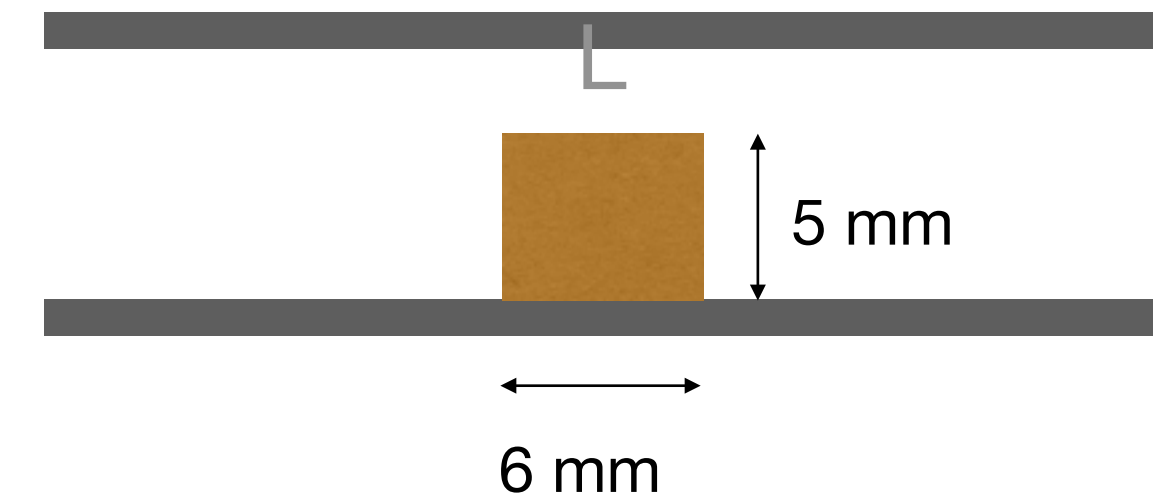


*sites with similar local environment are far apart in physical space, but nearby in perpendicular space*

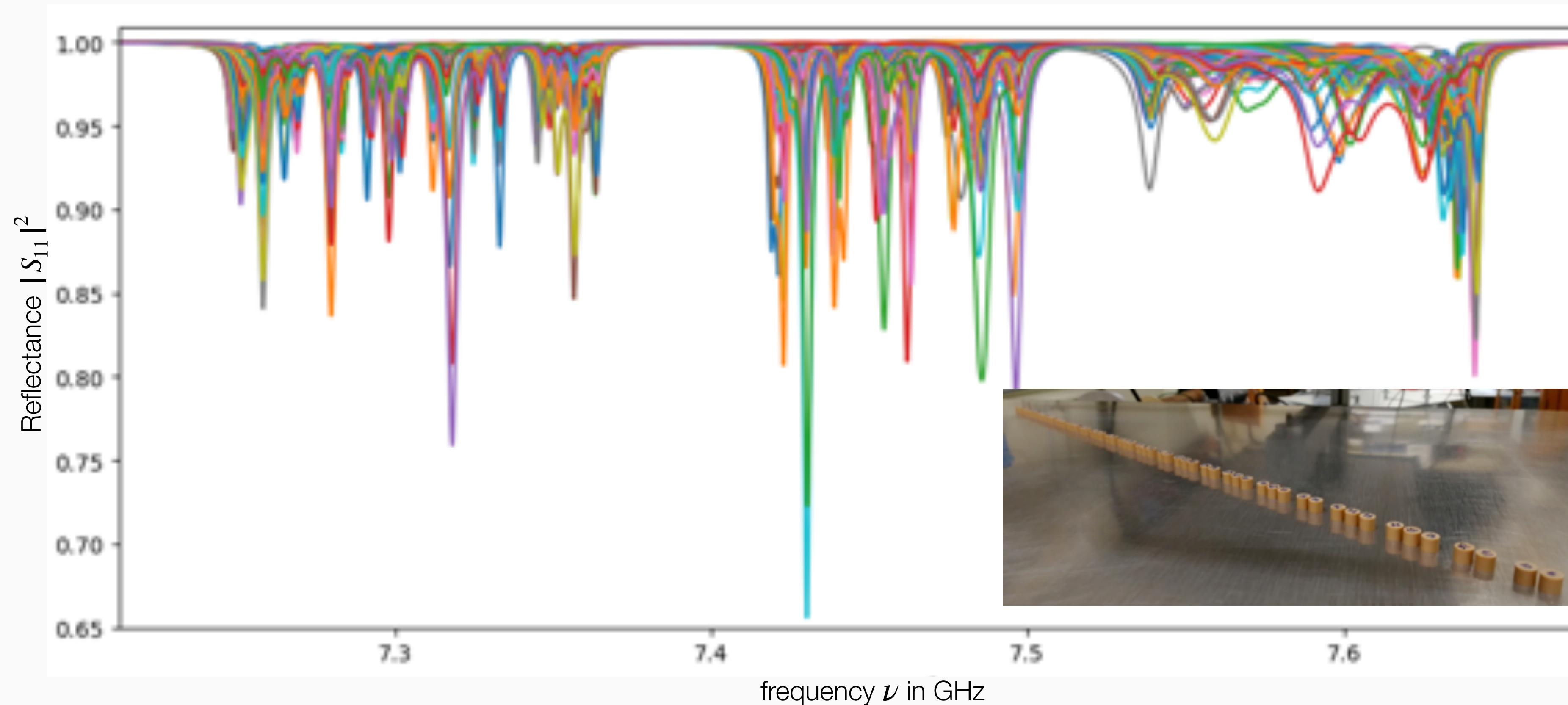


# The Experimental Fibonacci Chain

- coupled dielectric resonators,  $\nu_0 \approx 7.45$  GHz
- evanescent coupling  $\rightarrow d_A = 7$  mm ( $t_A = 126$  MHz) and  $d_B = 8$  mm ( $t_B = 81$  MHz)
- measure the reflection spectrum above each resonator  $\rightarrow$  local density of states



# The Experimental Fibonacci Chain



- measure  $S_{11}$  above all sites
- extract complex resonance frequencies and amplitudes for all peaks for all positions
- access to all eigenvalues  $\{\nu_j\}$  and eigenstates  $\{\Psi_j(i)\}$

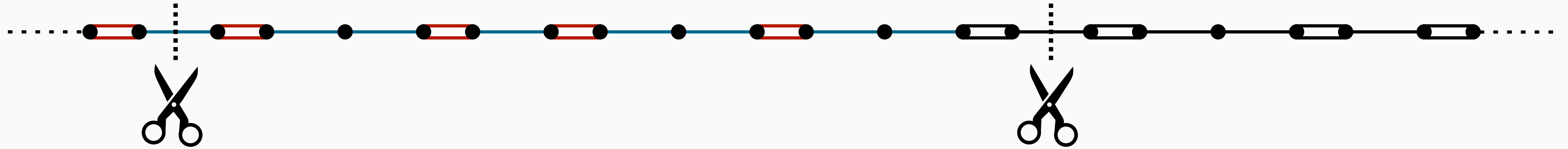
# Average over Different Permutations



- one basic motif ( $F_7 = 13$ )
- Different permutations of basic motif ( $F_7 = 13$ )
- “cut” on (neglect) a weak coupling
- chains end on strong coupling  $\rightarrow$  no edge states



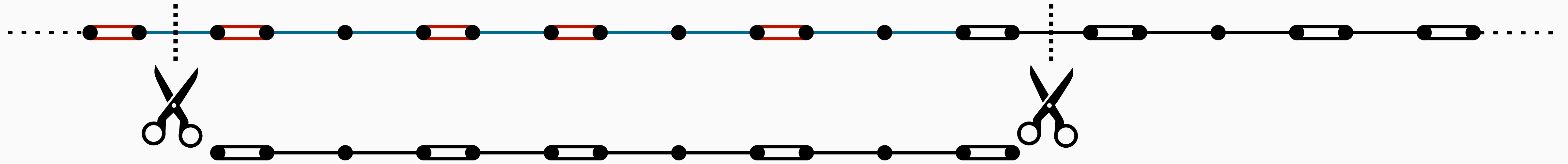
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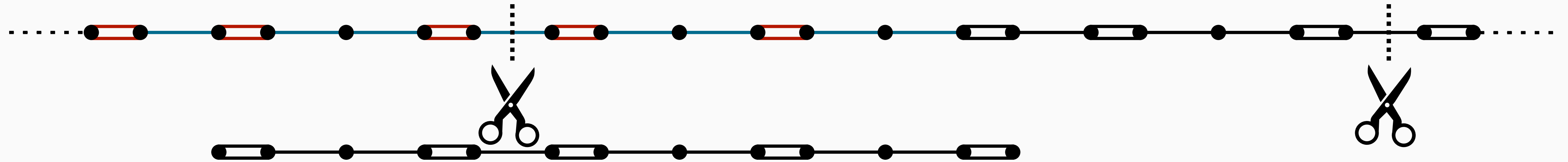


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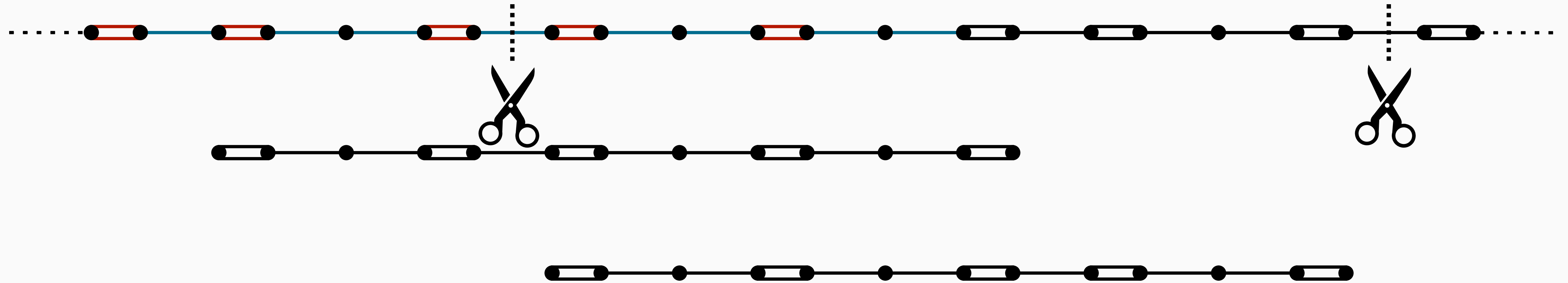
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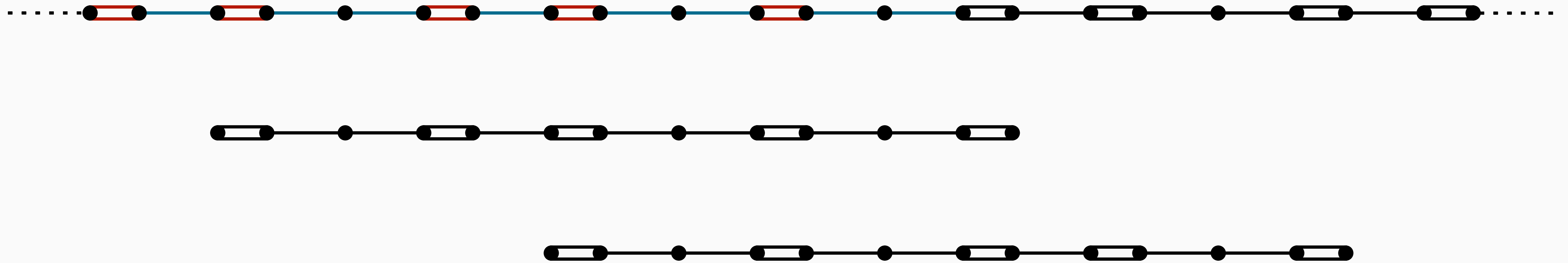
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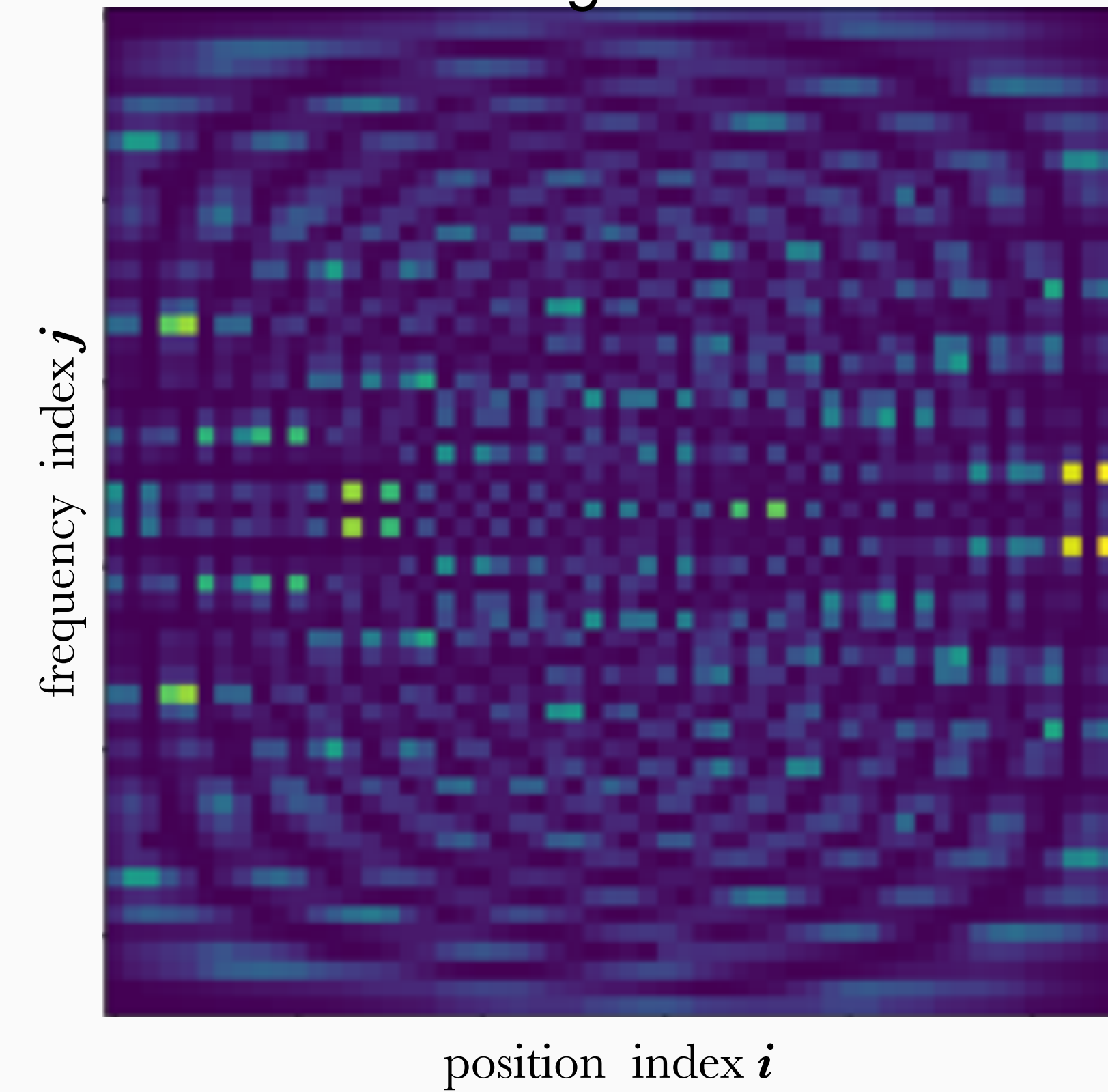


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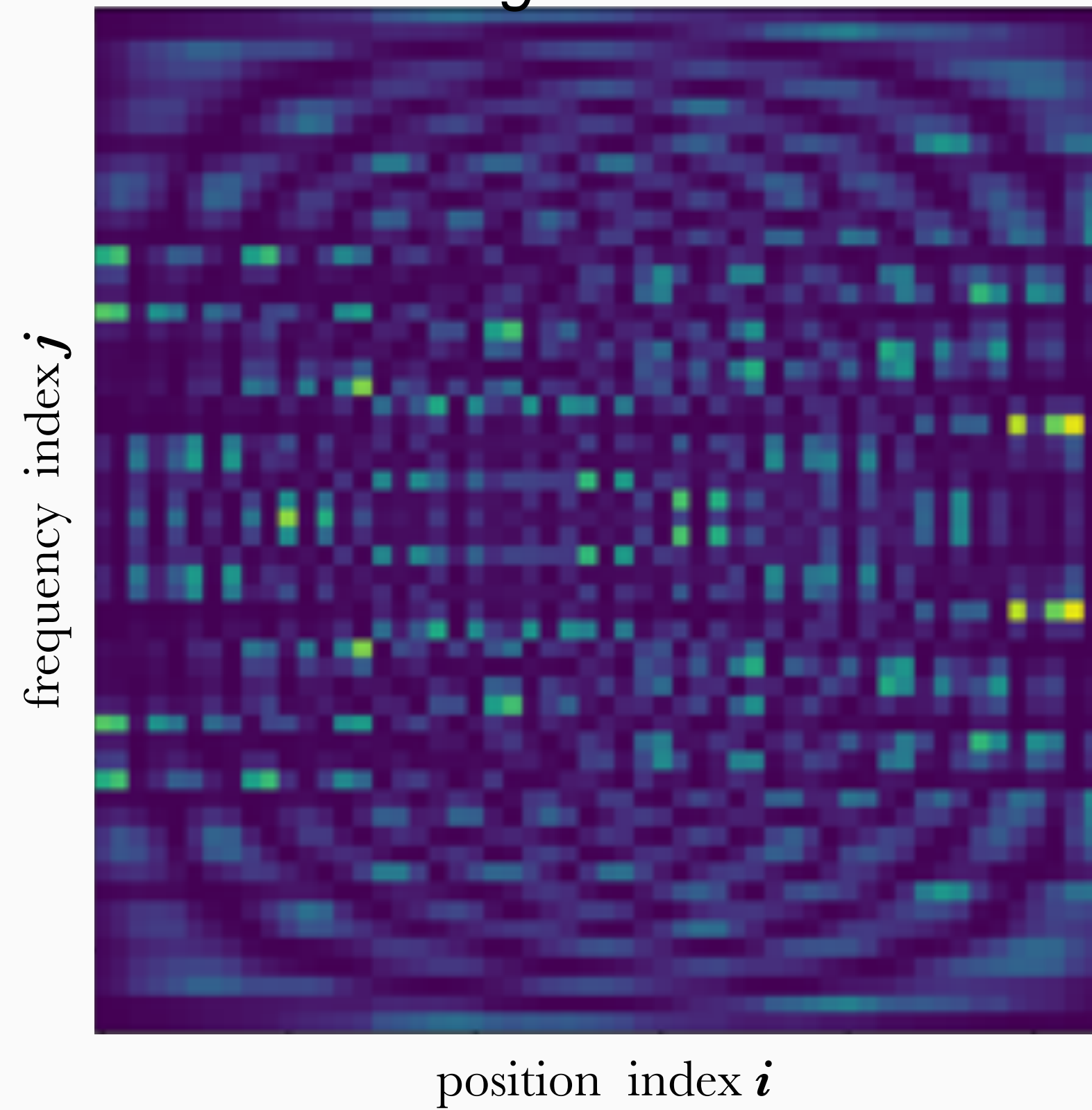
# Average over Different Permutations

- $F_{10} = 55$  resonators
- 8 different permutations

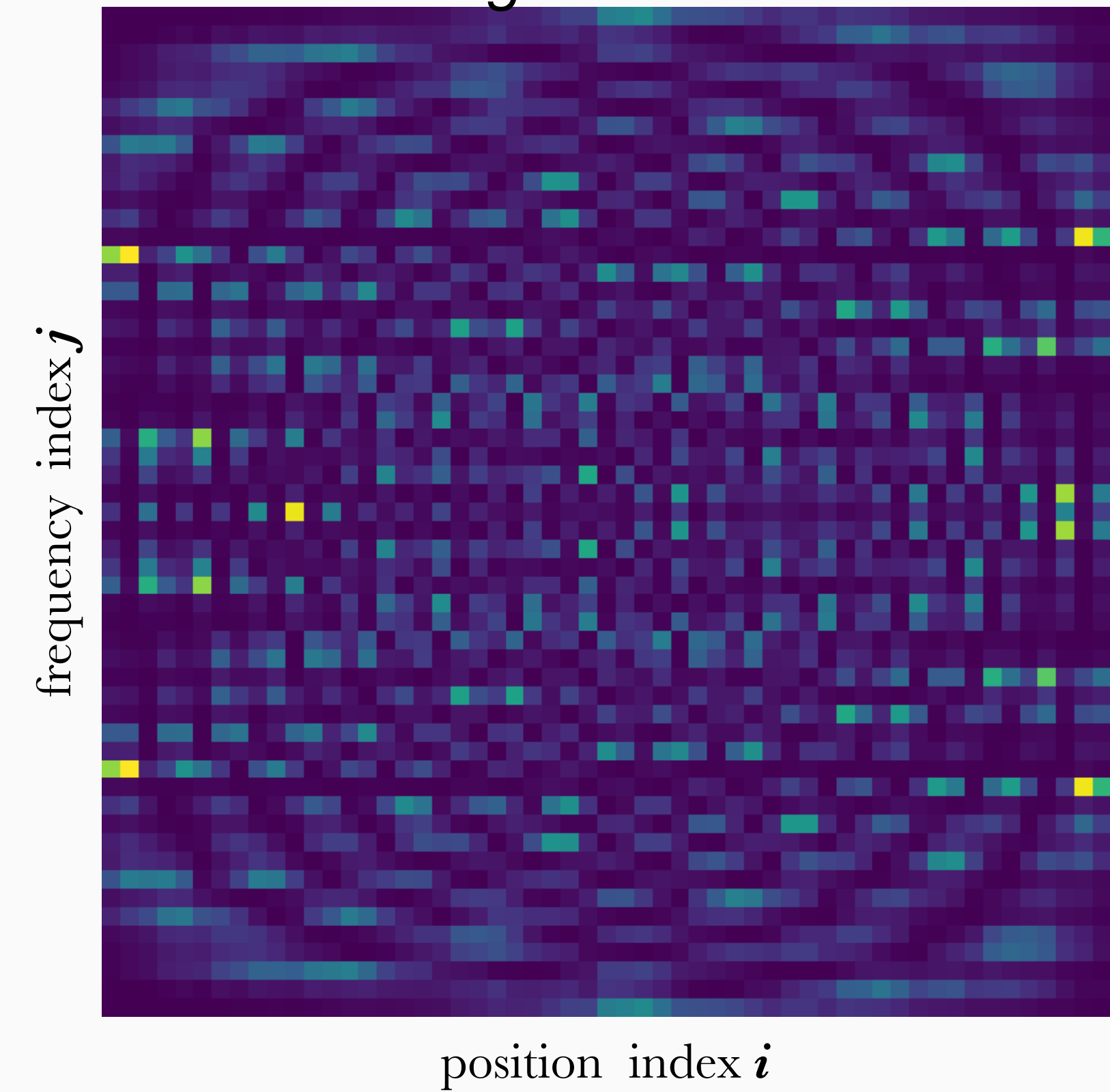
*configuration 1*



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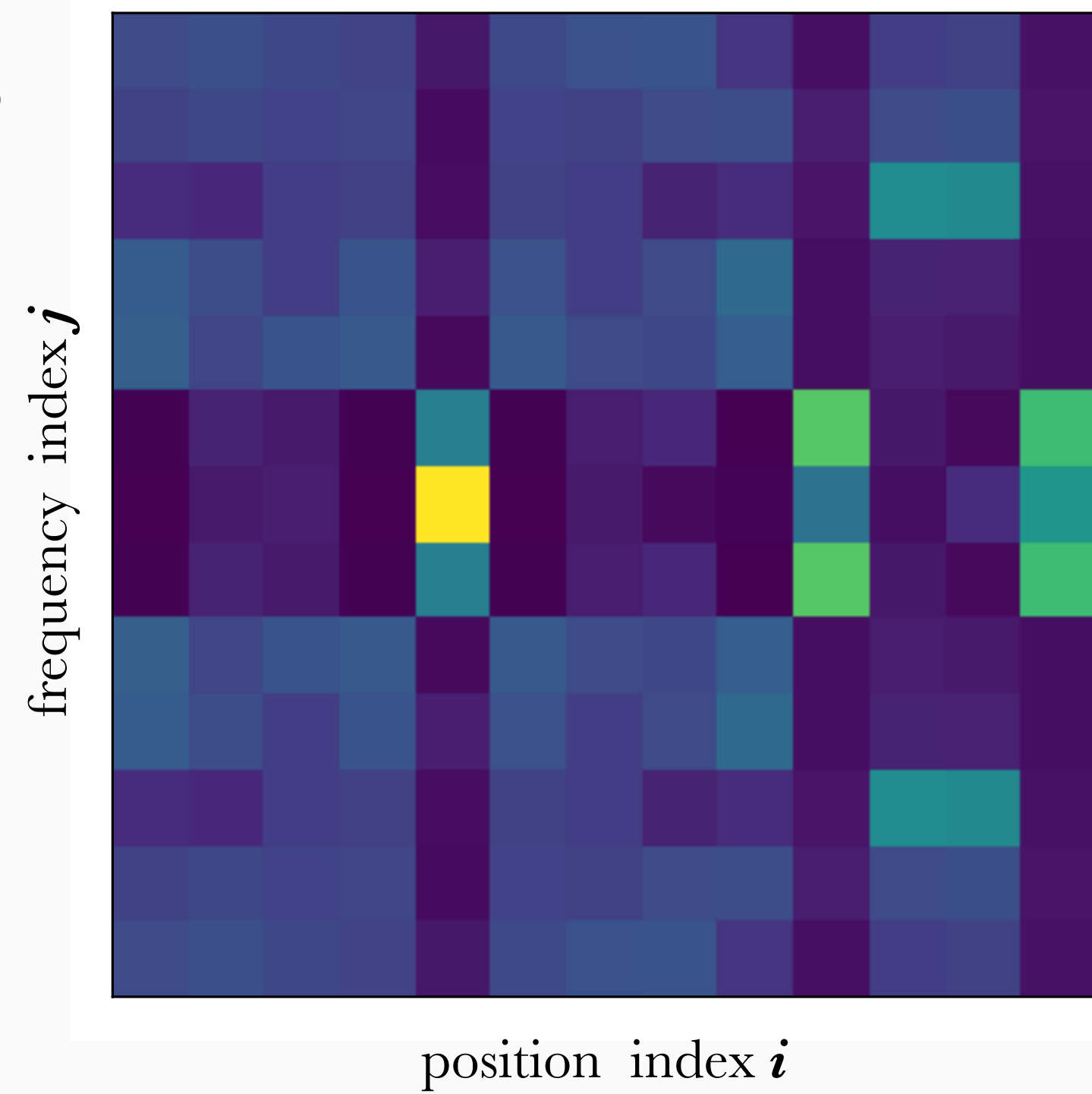


*configuration 3*



# Revealing the Fractality via Conumbers

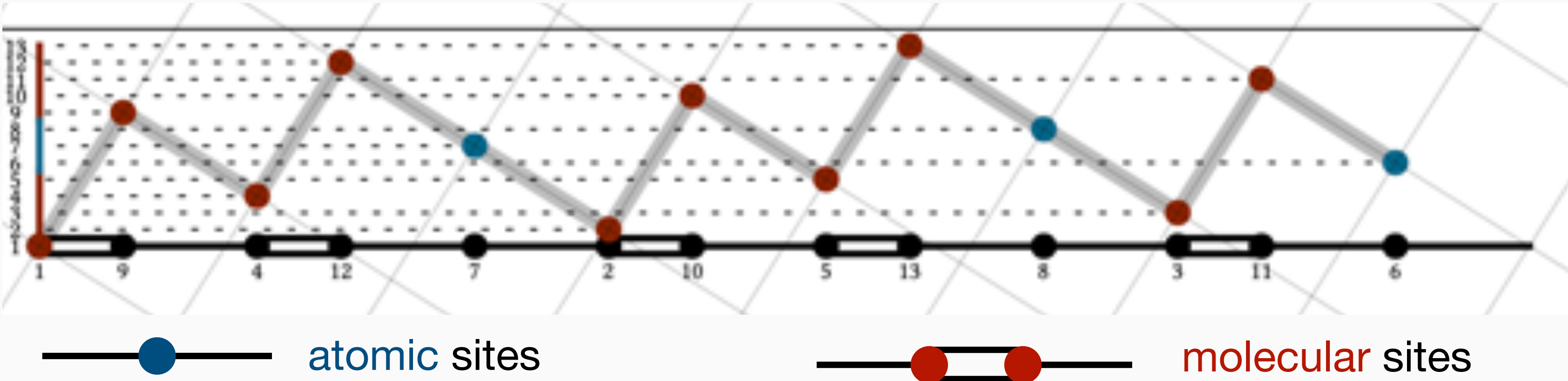
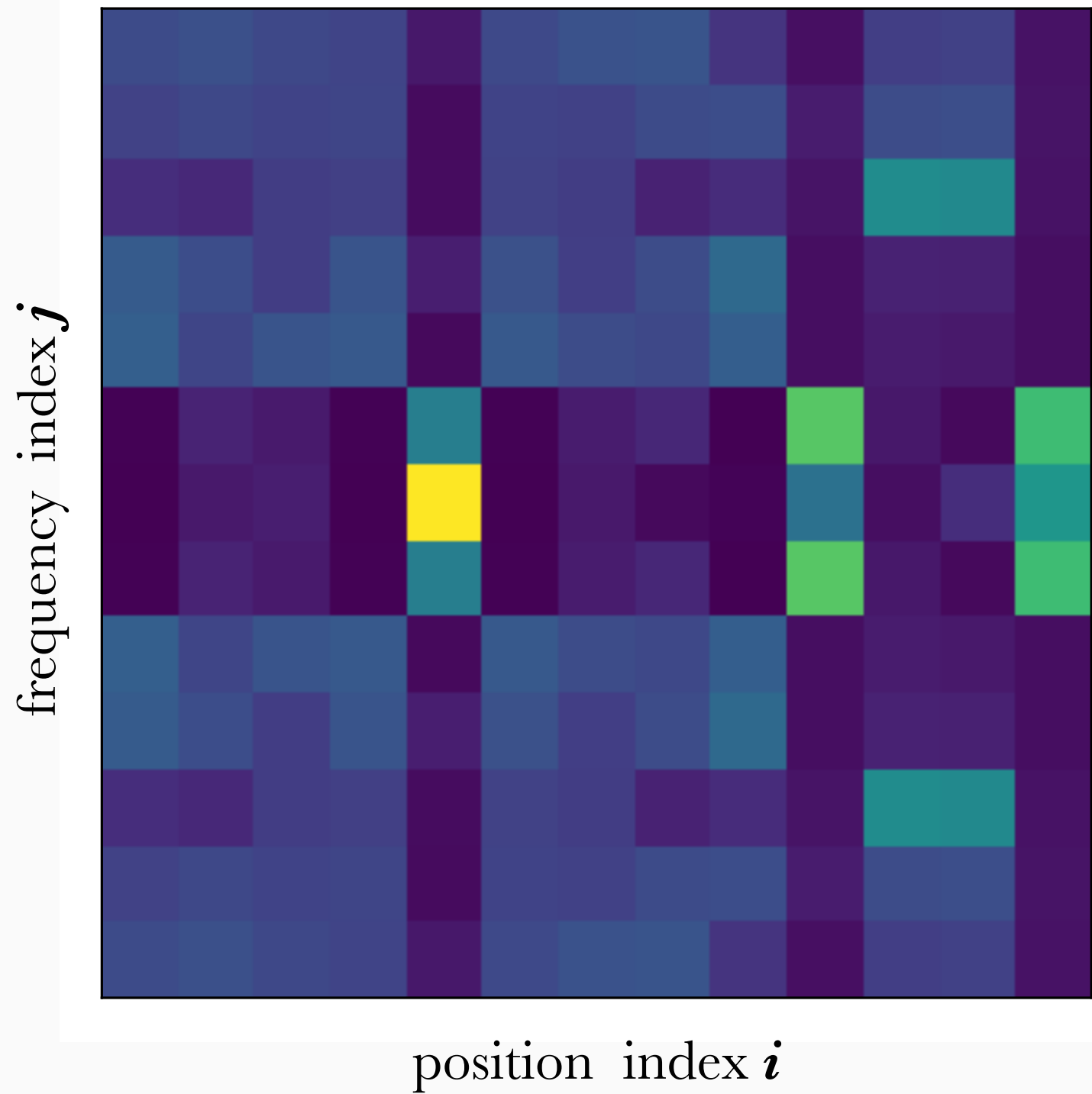
$$F_7 = 13$$





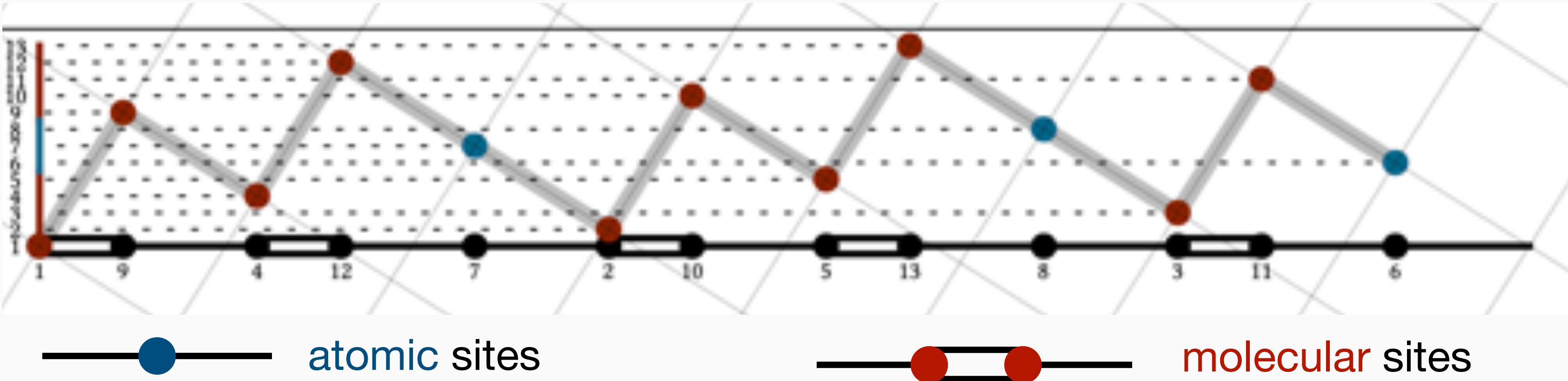
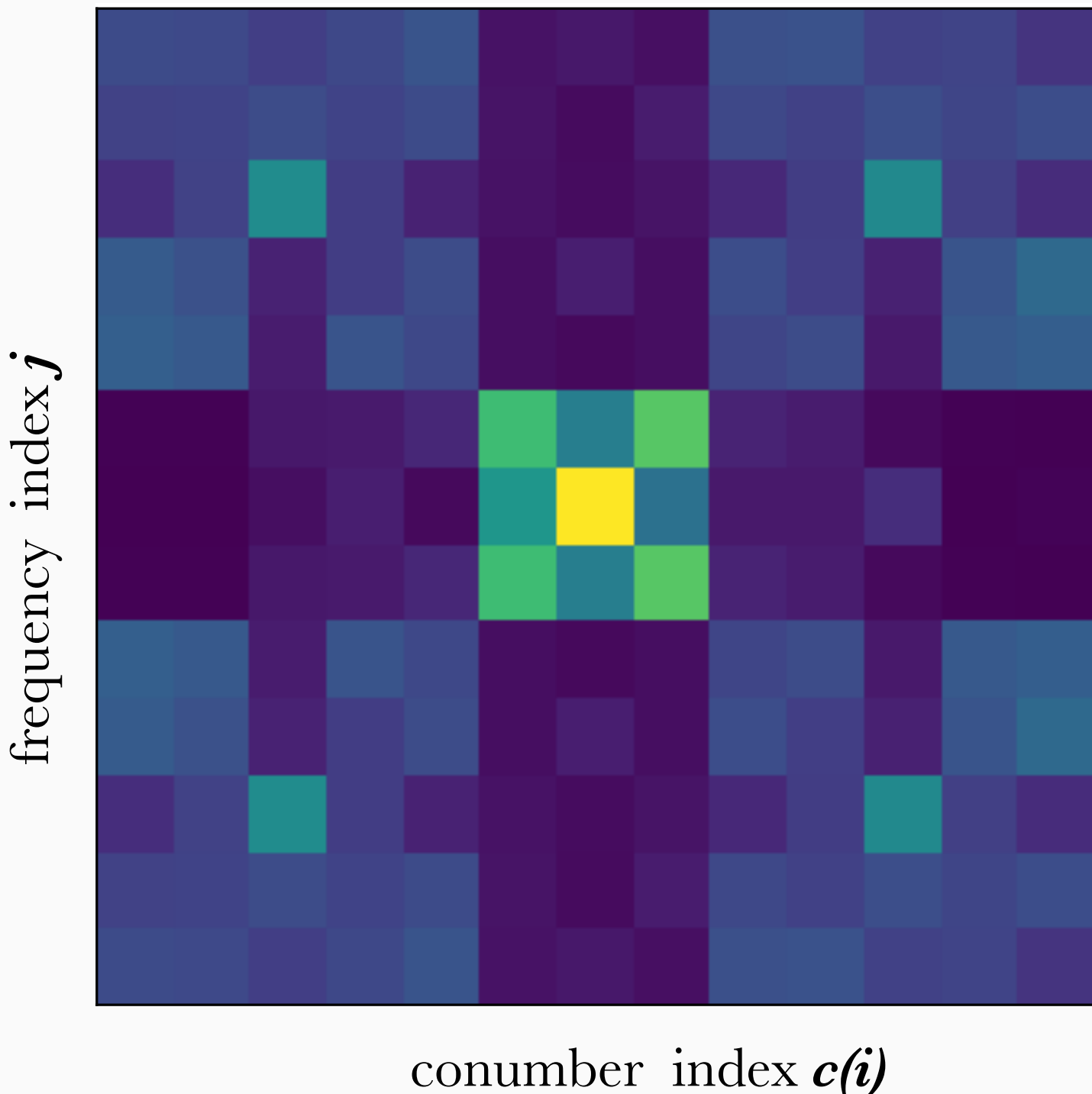
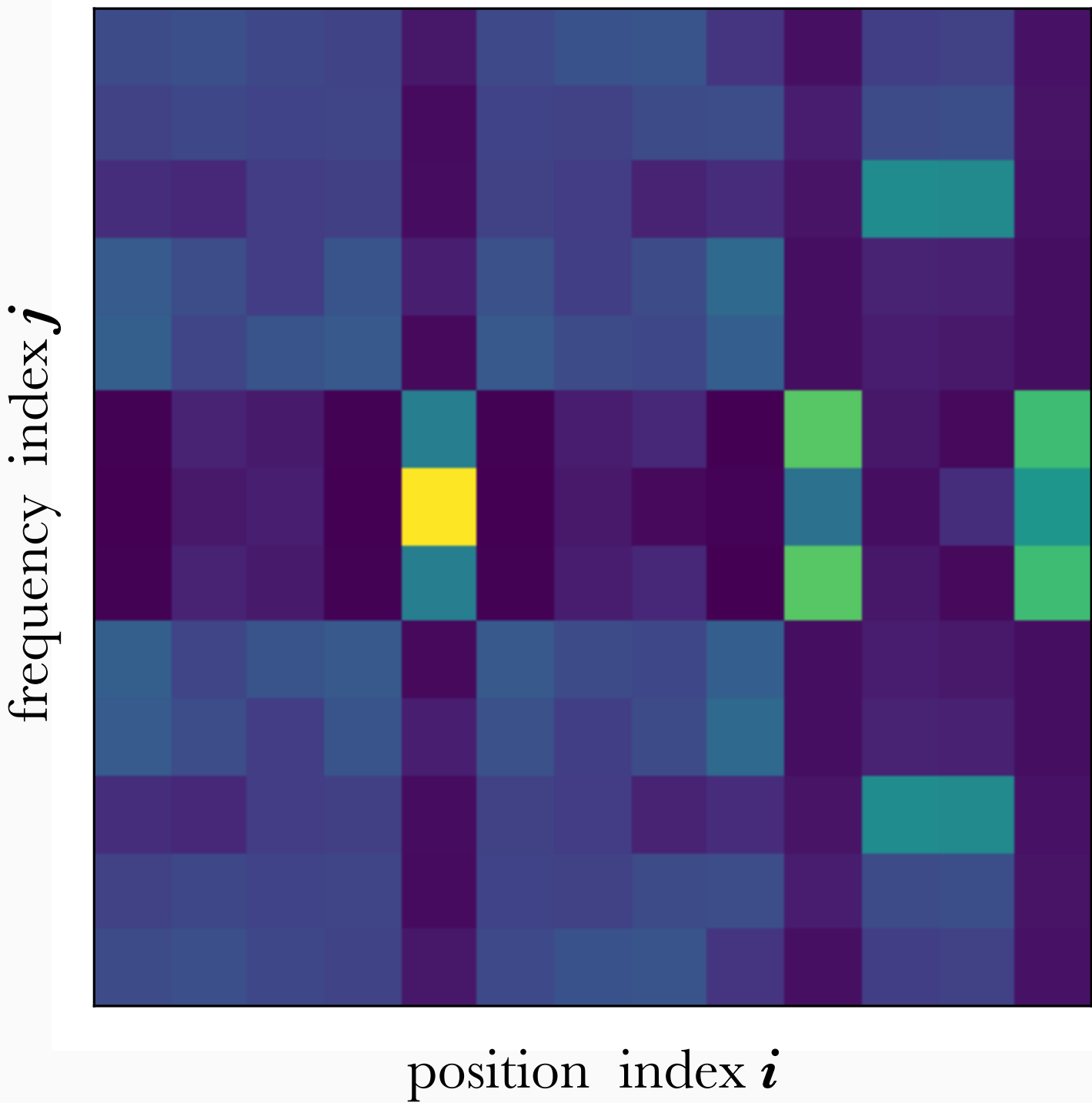
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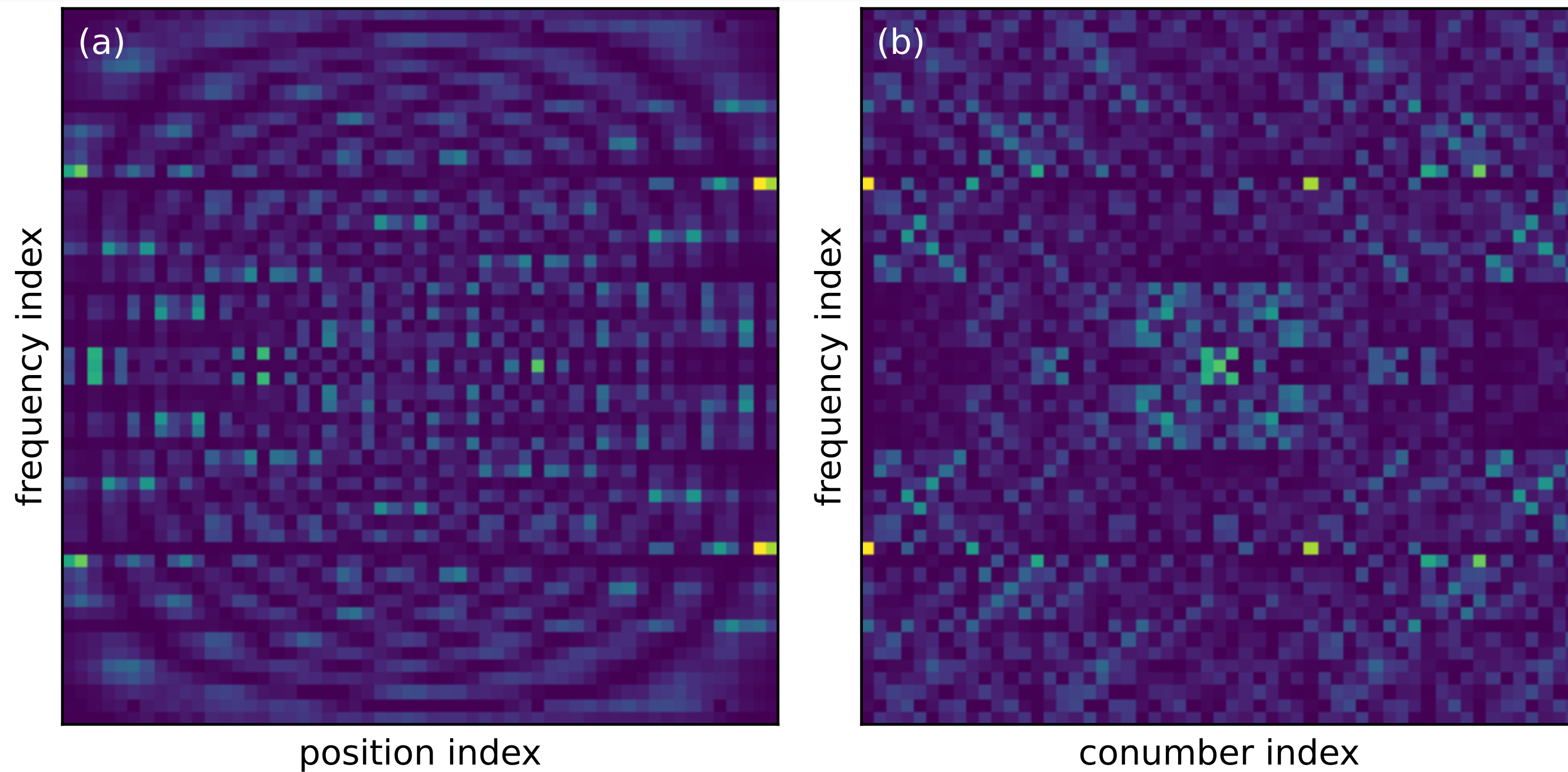
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# Revealing the Fractality via Conumbers

$$F_7 = 55$$

*one single configuration*



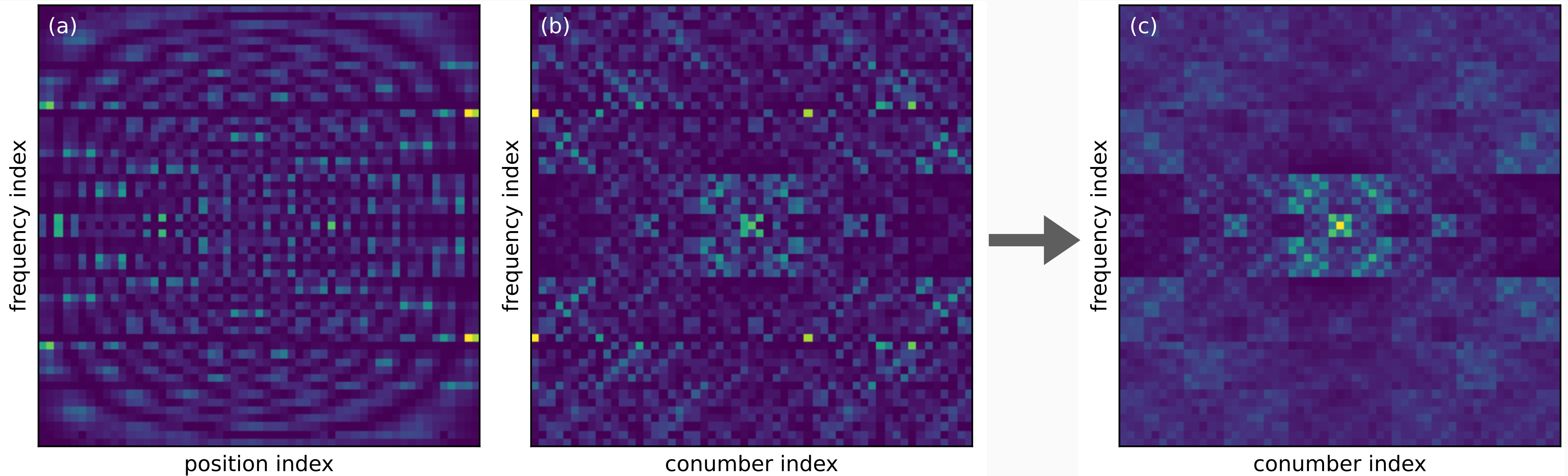


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*one single configuration*

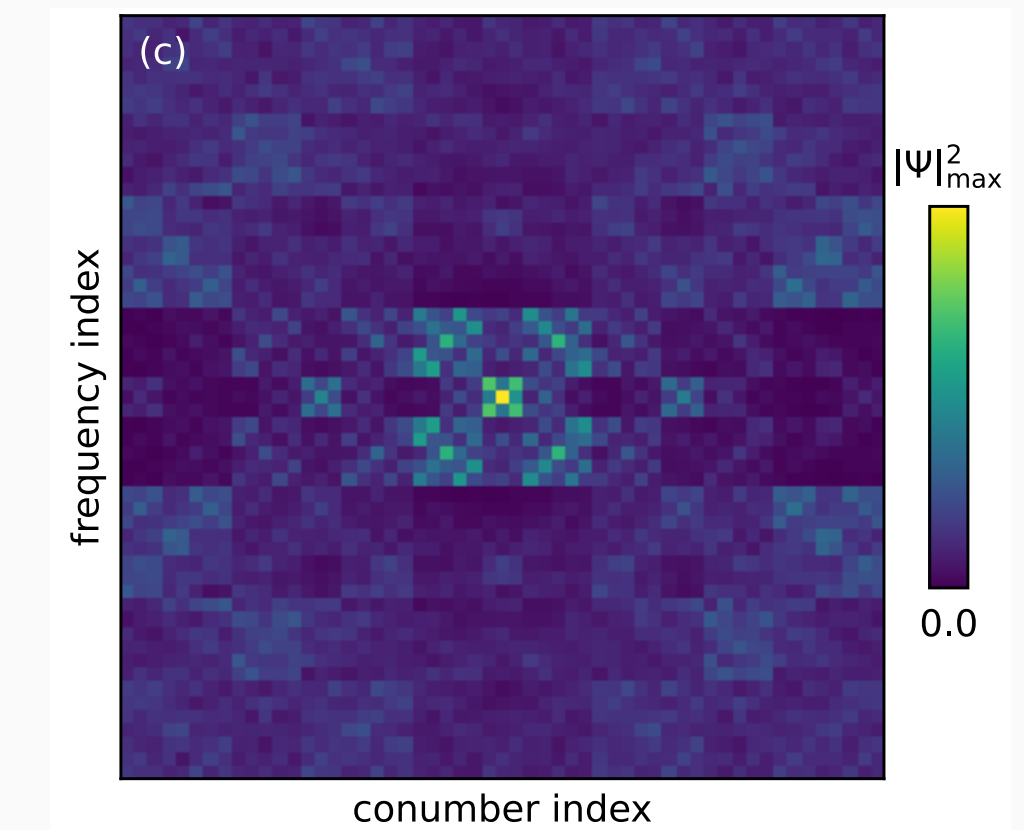
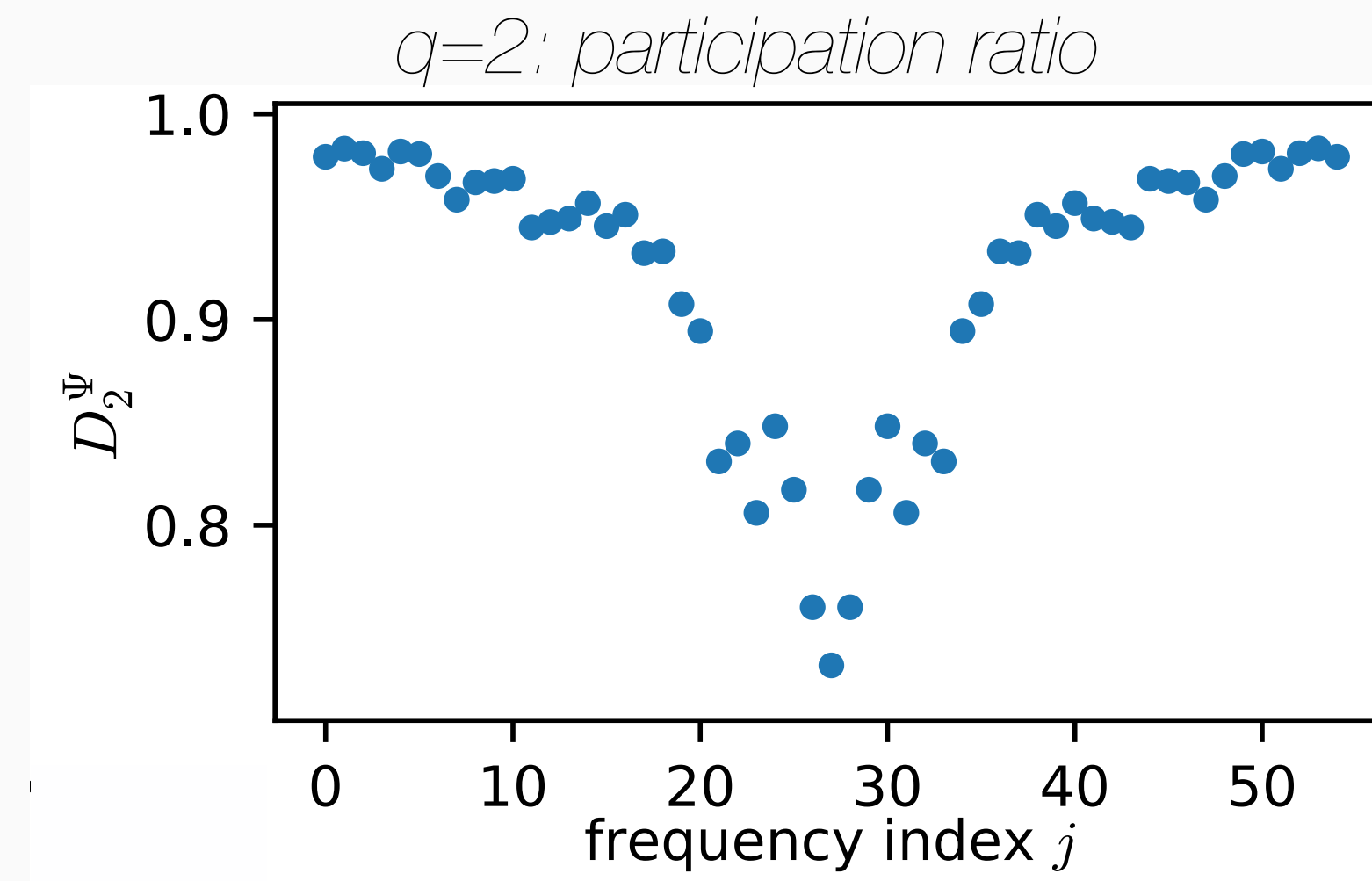
*average over all permutations*



# Fractal Dimensions of the Wavefunctions

Fractal Dimension  $D_q^\psi$  of state  $j$  as a function of  $q$

$$\chi_q^{(n)}(j) = \sum_i |\psi_j^{(n)}(i)|^{2q} \underset{n \rightarrow \infty}{\sim} F_n^{-(q-1)D_q^\psi(j)}$$



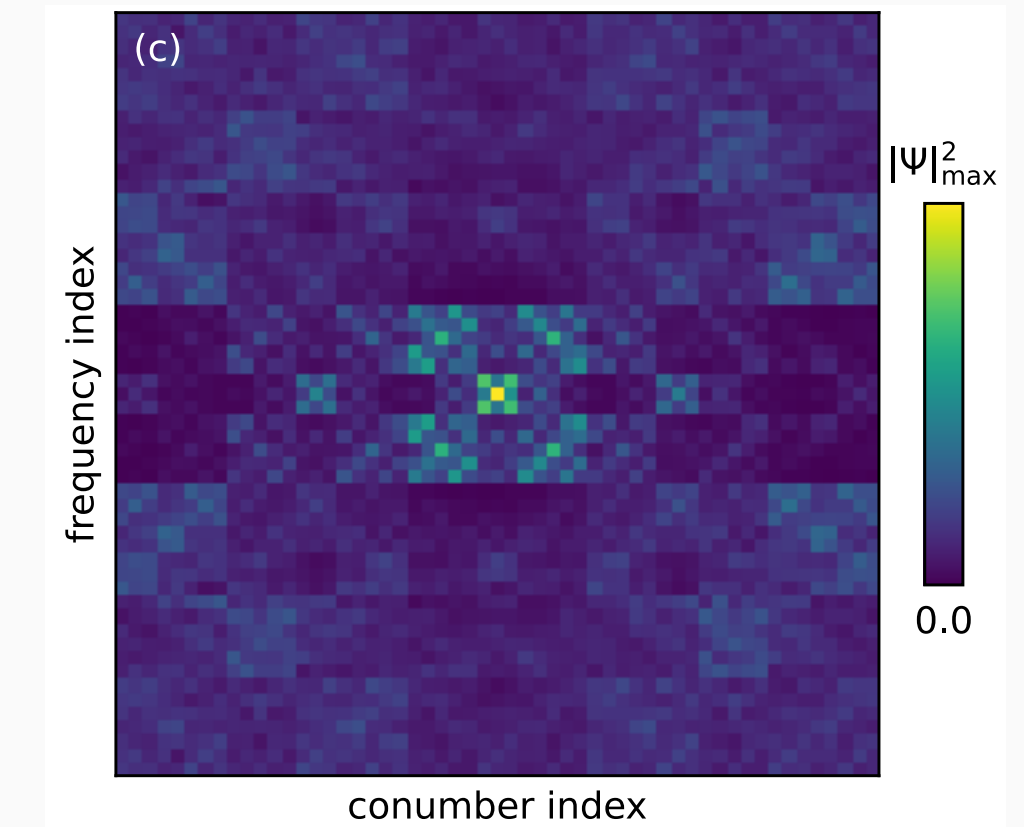
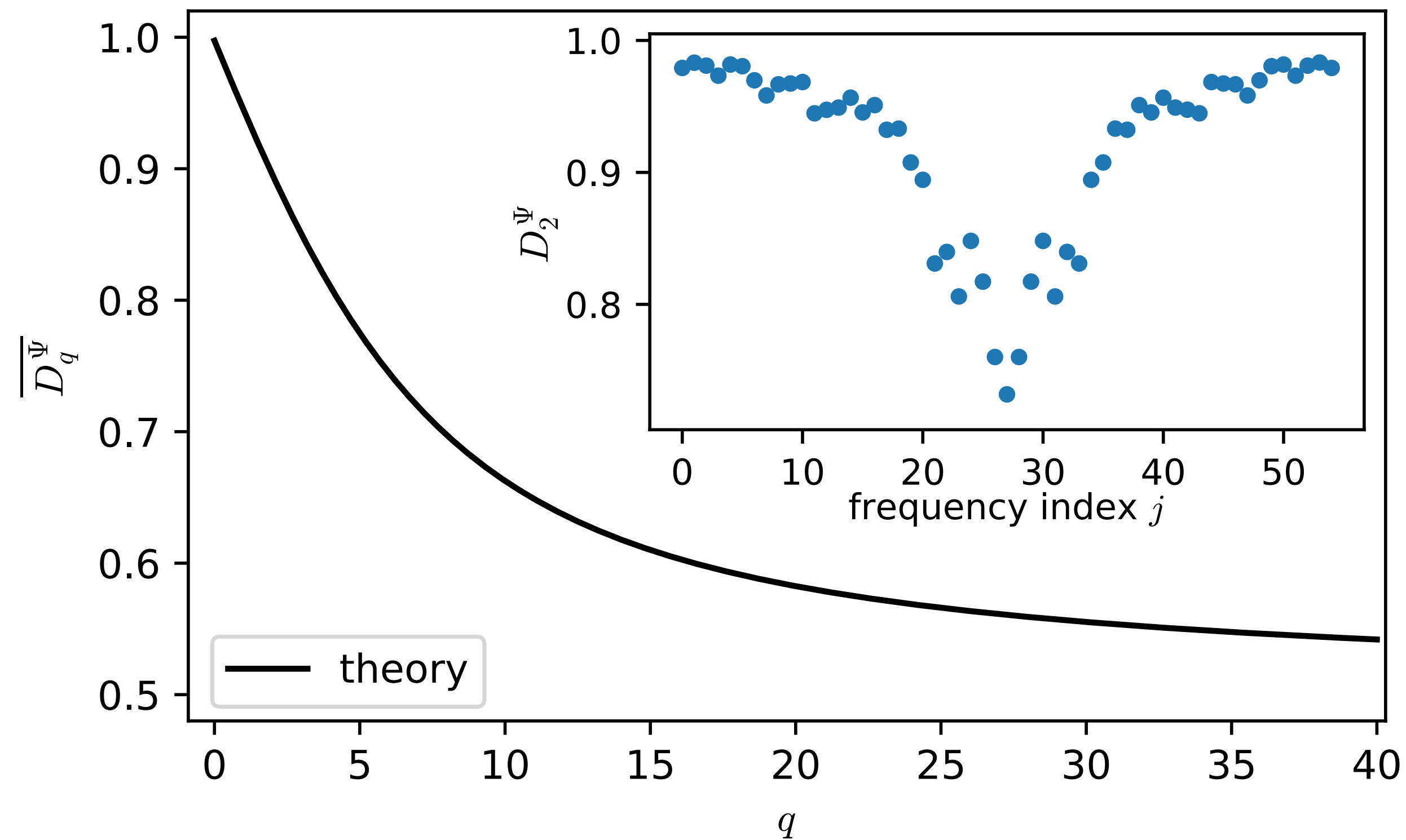
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Averaged Fractal Dimension  $\overline{D_q^\psi}$

$$\langle \chi_q^n(j) \rangle_j = \frac{1}{F_n} \sum_j \chi_q^{(n)}(j) \underset{n \rightarrow \infty}{\sim} F_n^{-(q-1)\overline{D_q^\psi}}$$





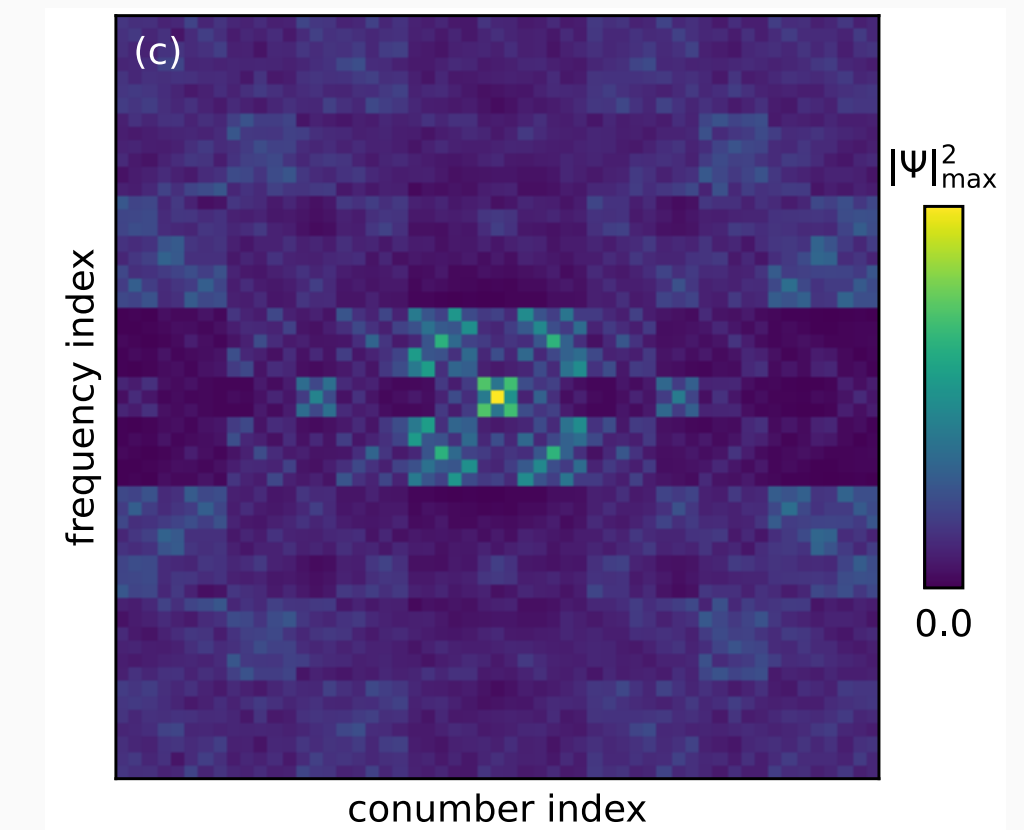
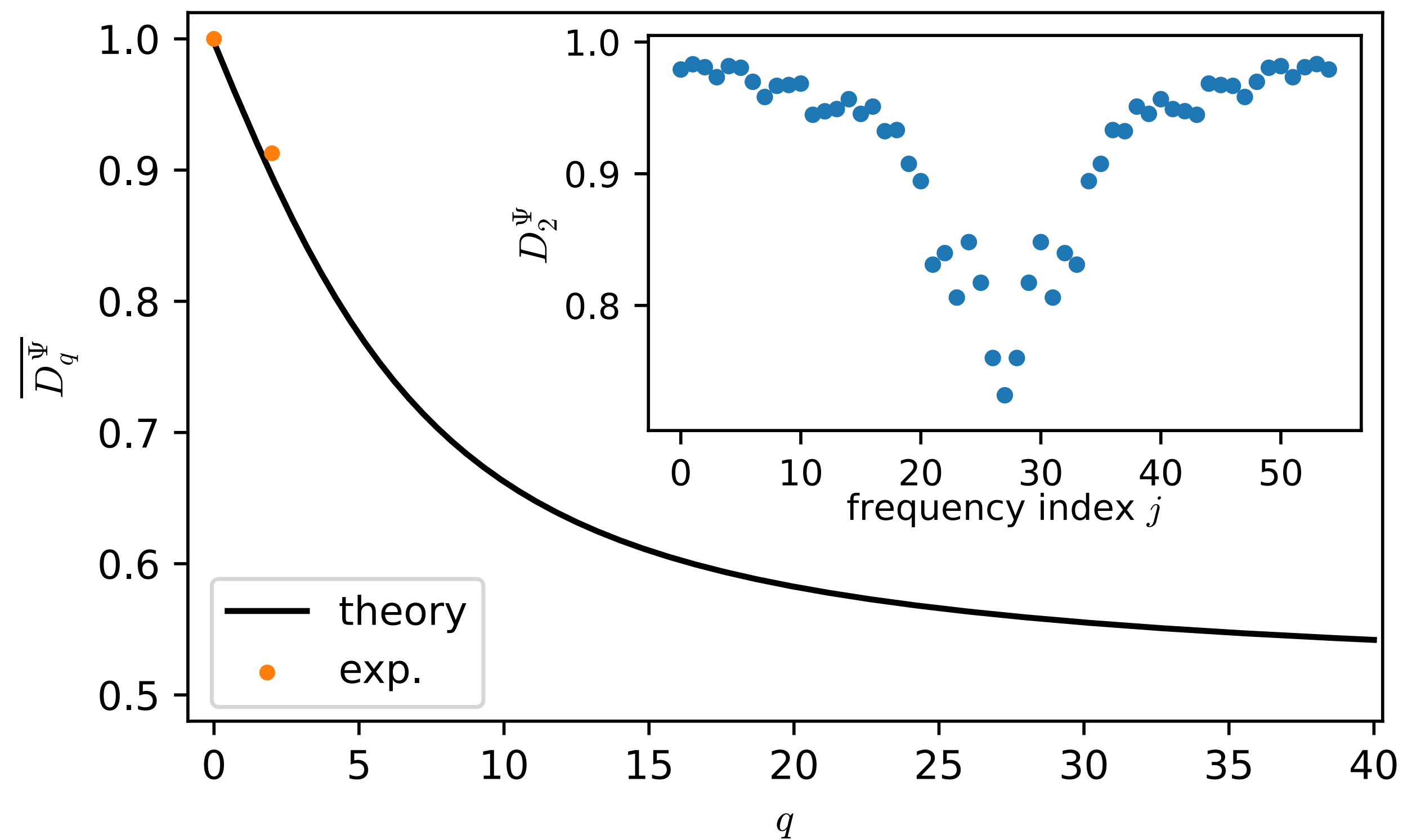
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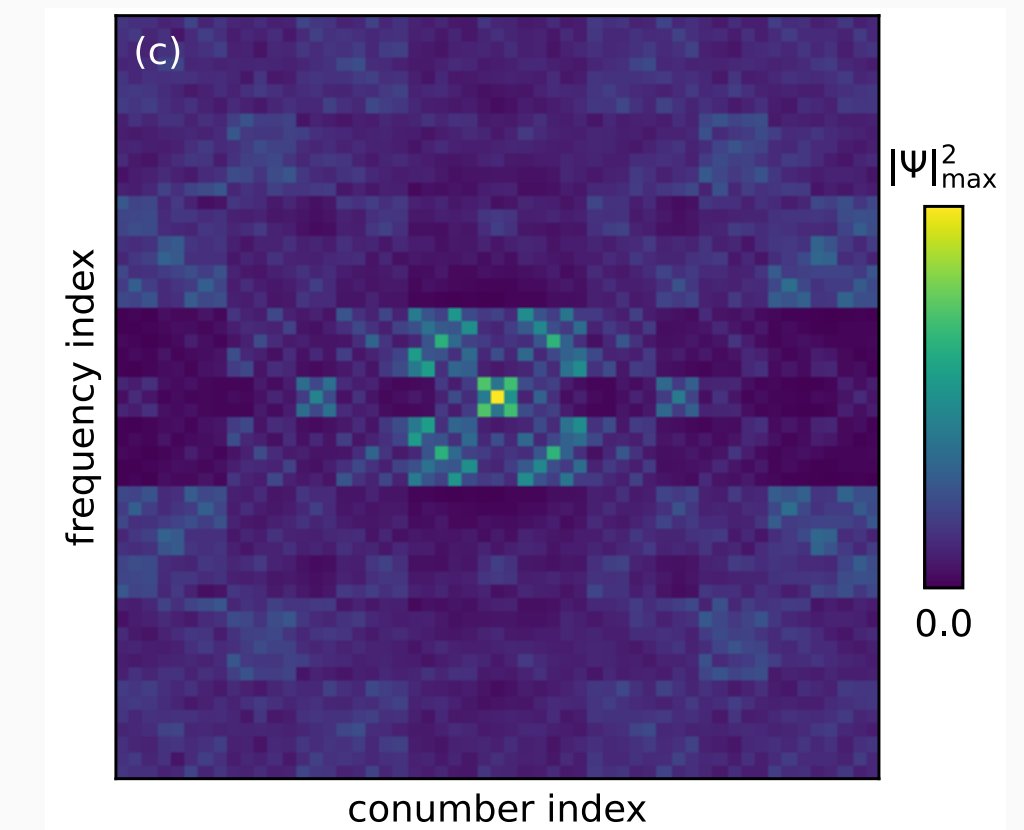
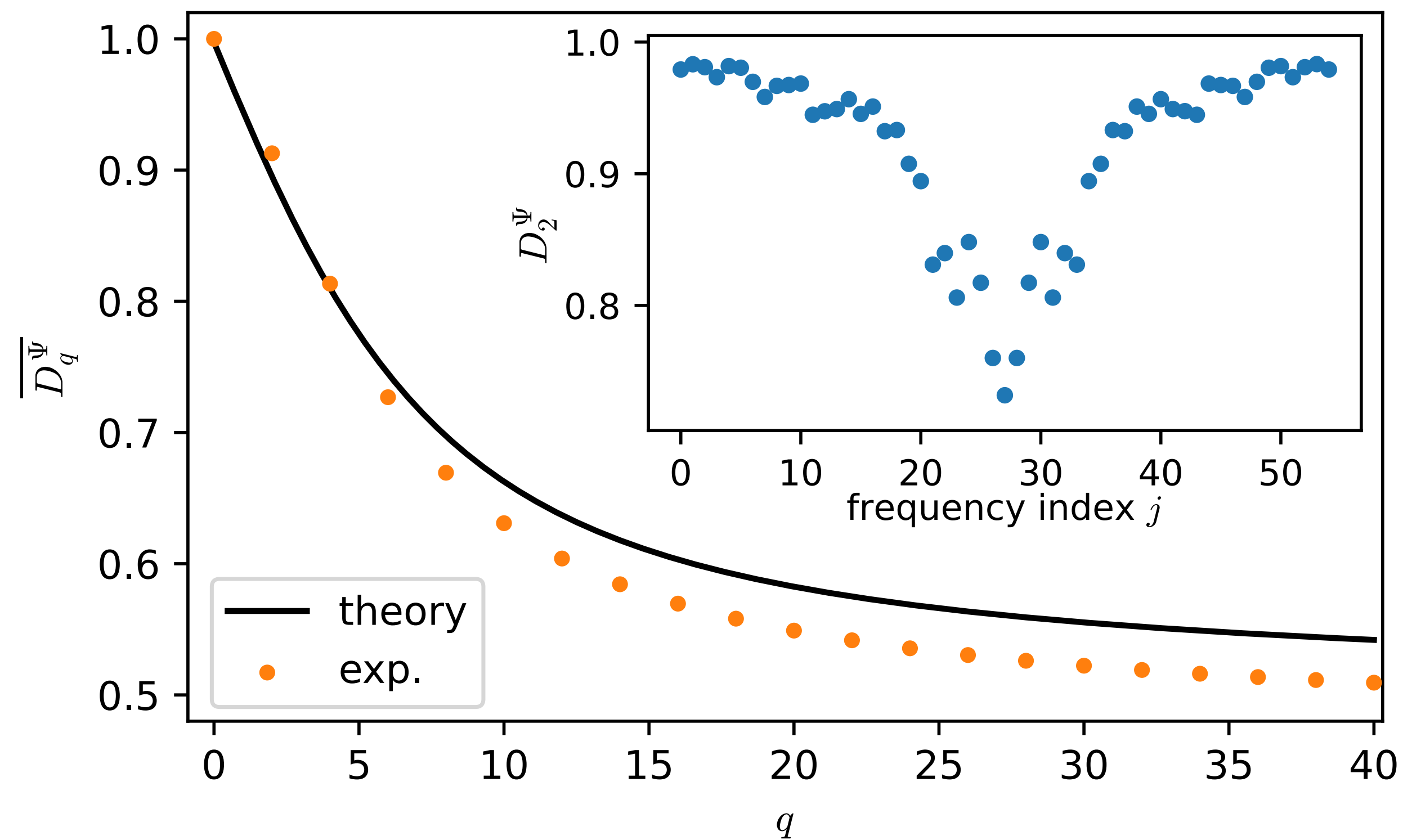
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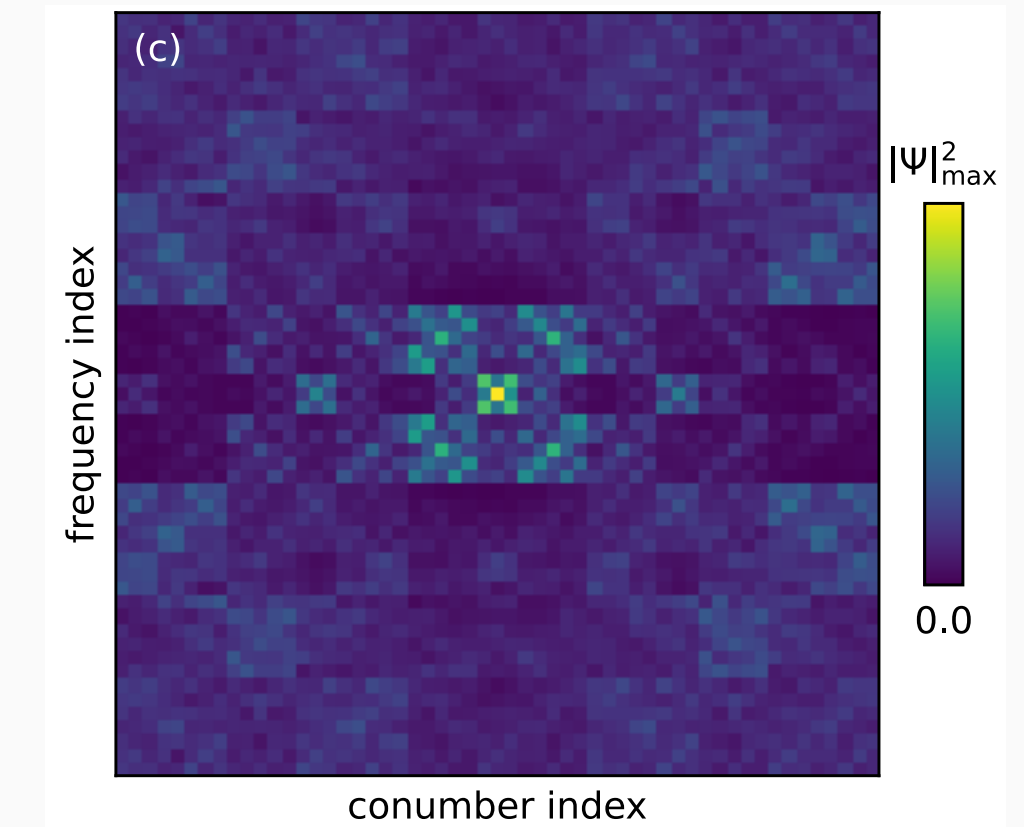
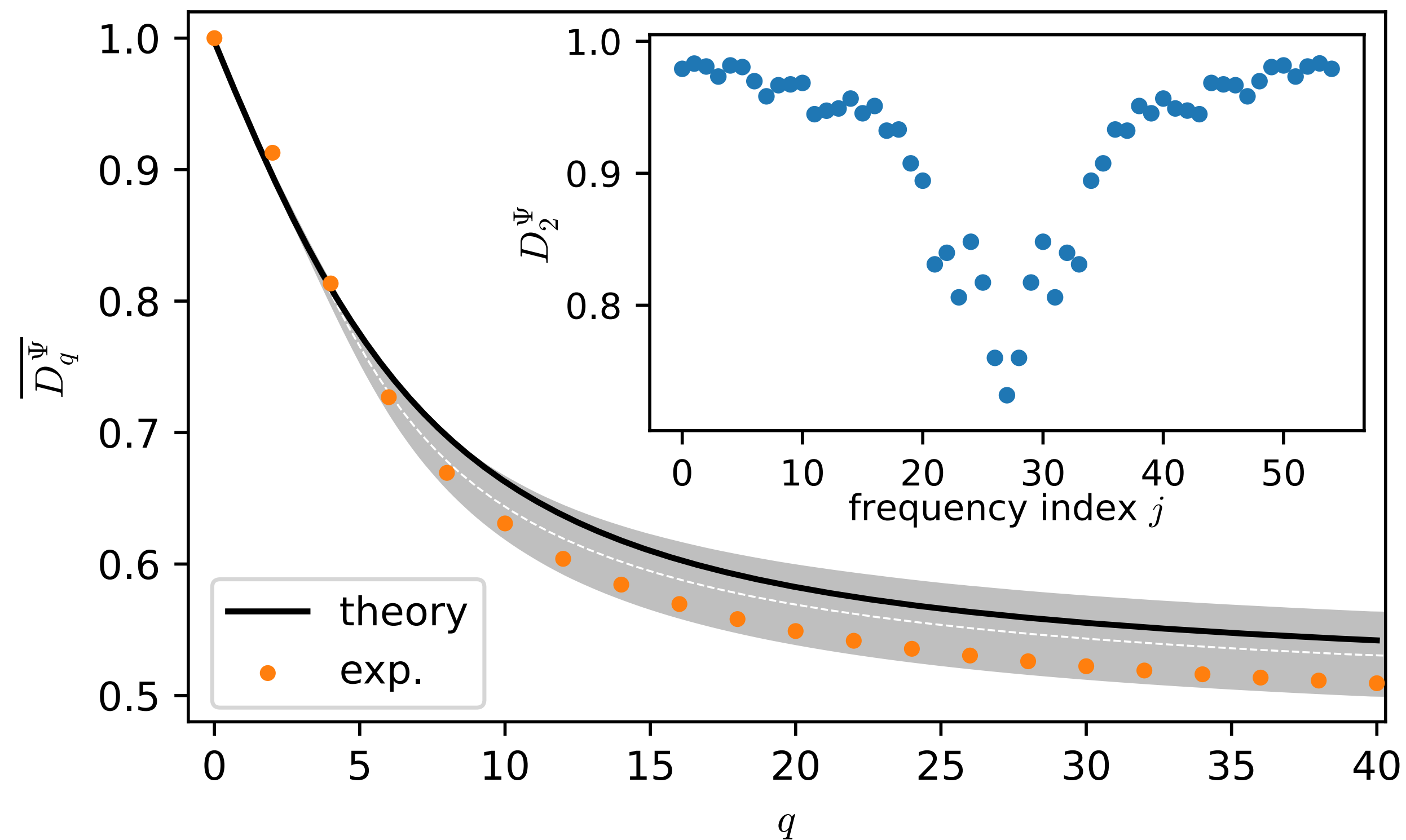
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Reasons for the offset:

- Finite system size
- Experimental fluctuations

For more details see [arXiv:2207.13755](https://arxiv.org/abs/2207.13755)

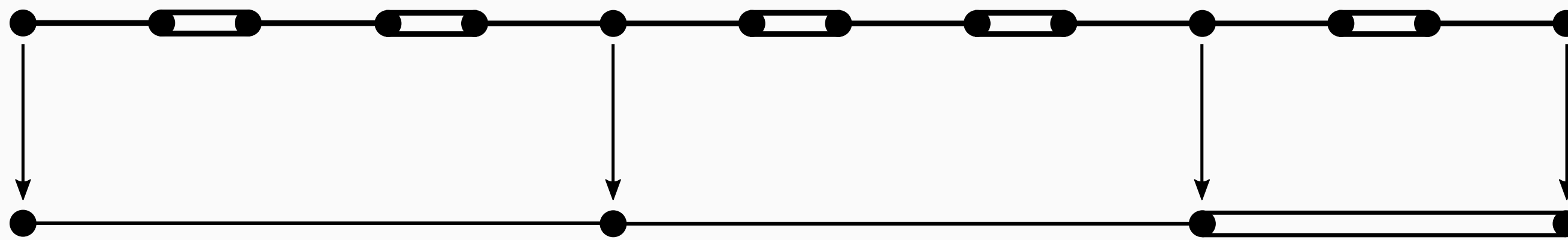


# Recursive Construction

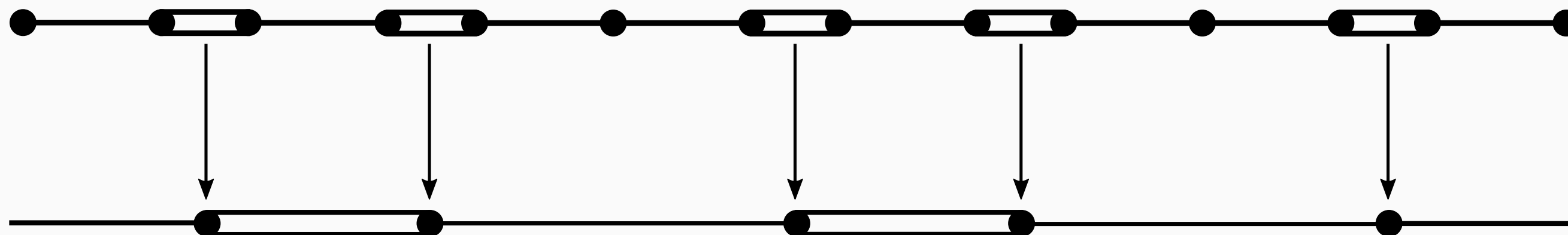
- renormalization theory to describe recursive construction
- different recursive construction for **atomic** and **molecular** sites/energies

$F_{10}=55$

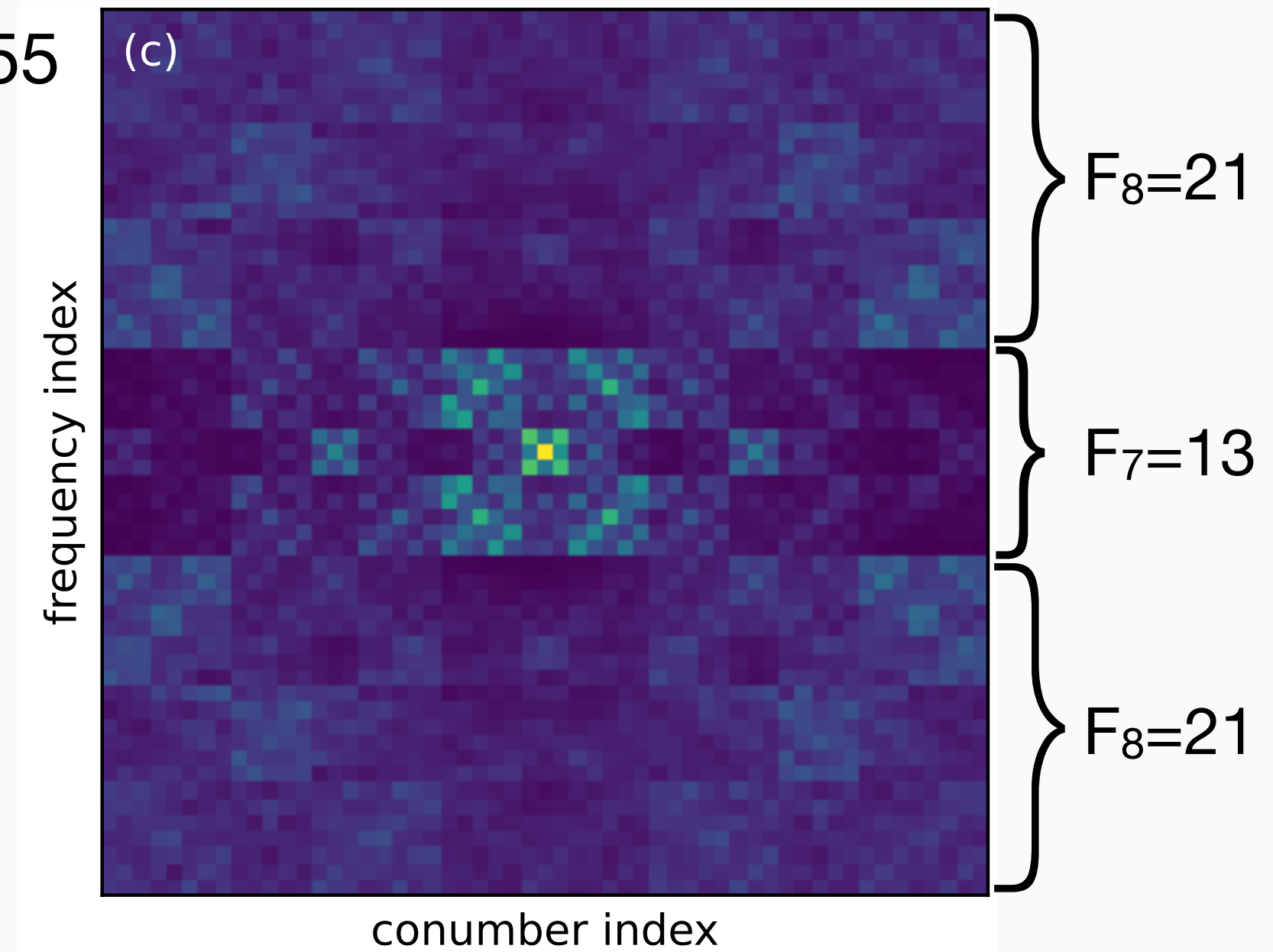
atomic



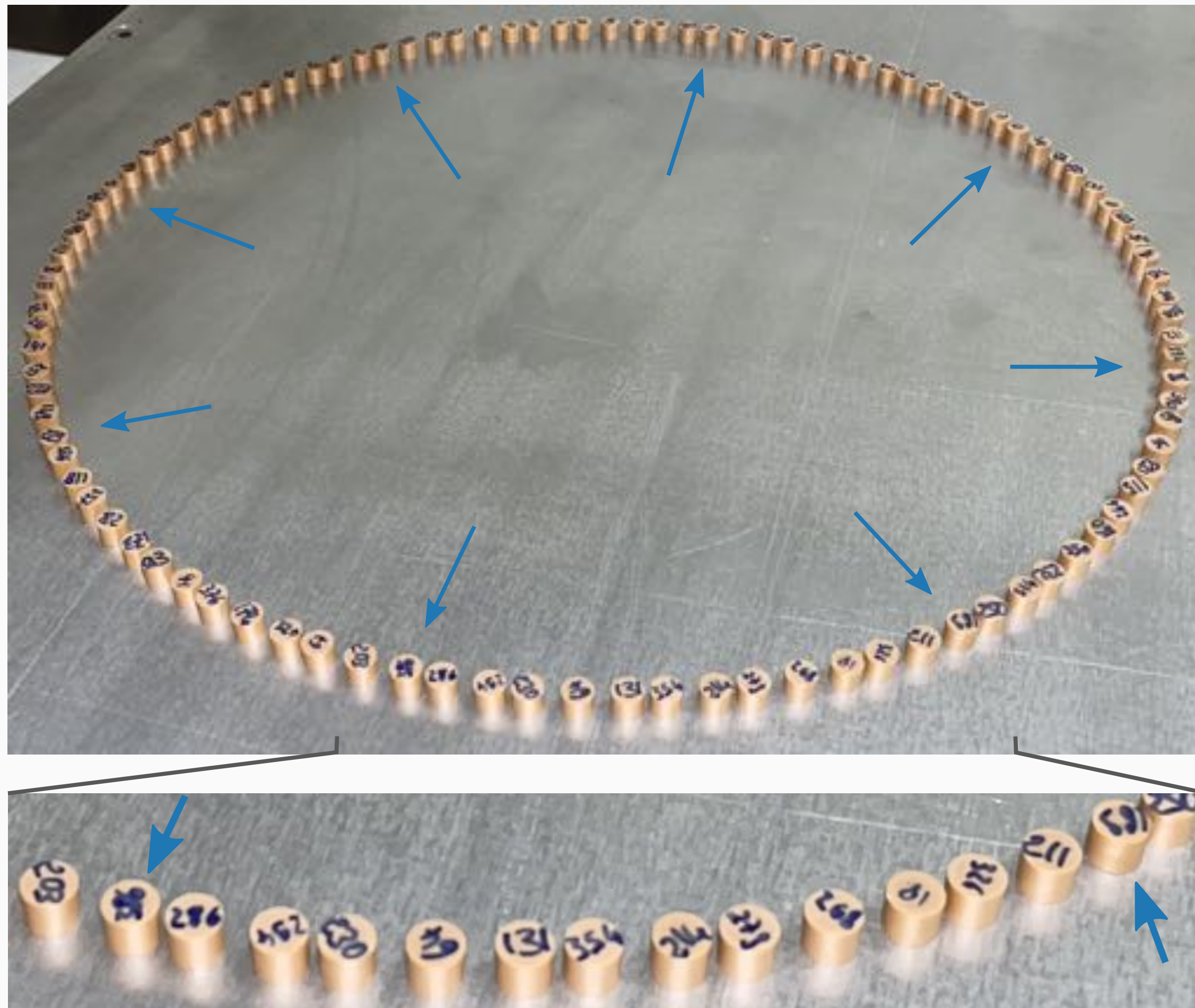
molecular



- for **atomic** sites/energies:  $\psi^{(n)} = \tilde{\lambda} \cdot \psi^{(n-3)}$
- for **molecular** sites/energies:  $\psi^{(n)} = \lambda \cdot \psi^{(n-2)}$



# Study of the Recursive Construction with Circular Chains

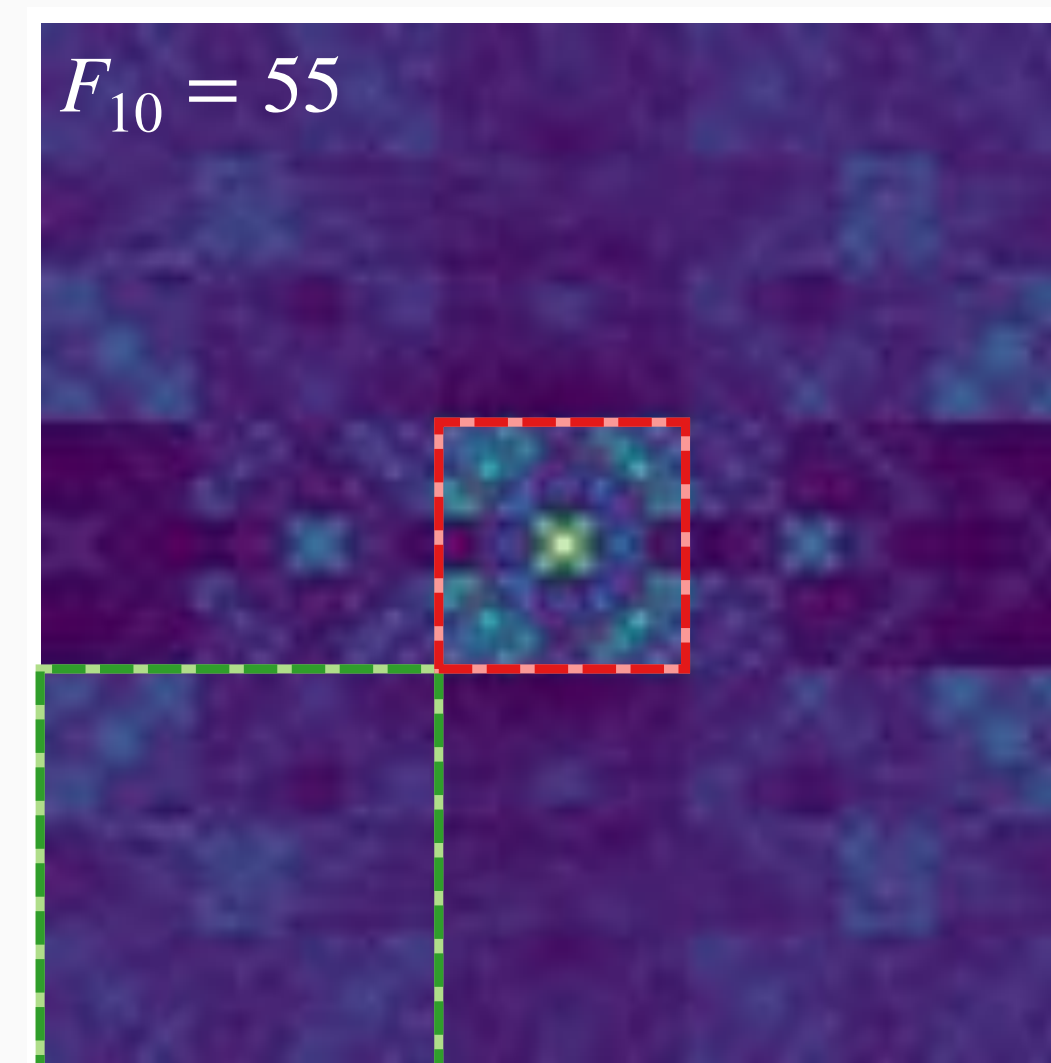
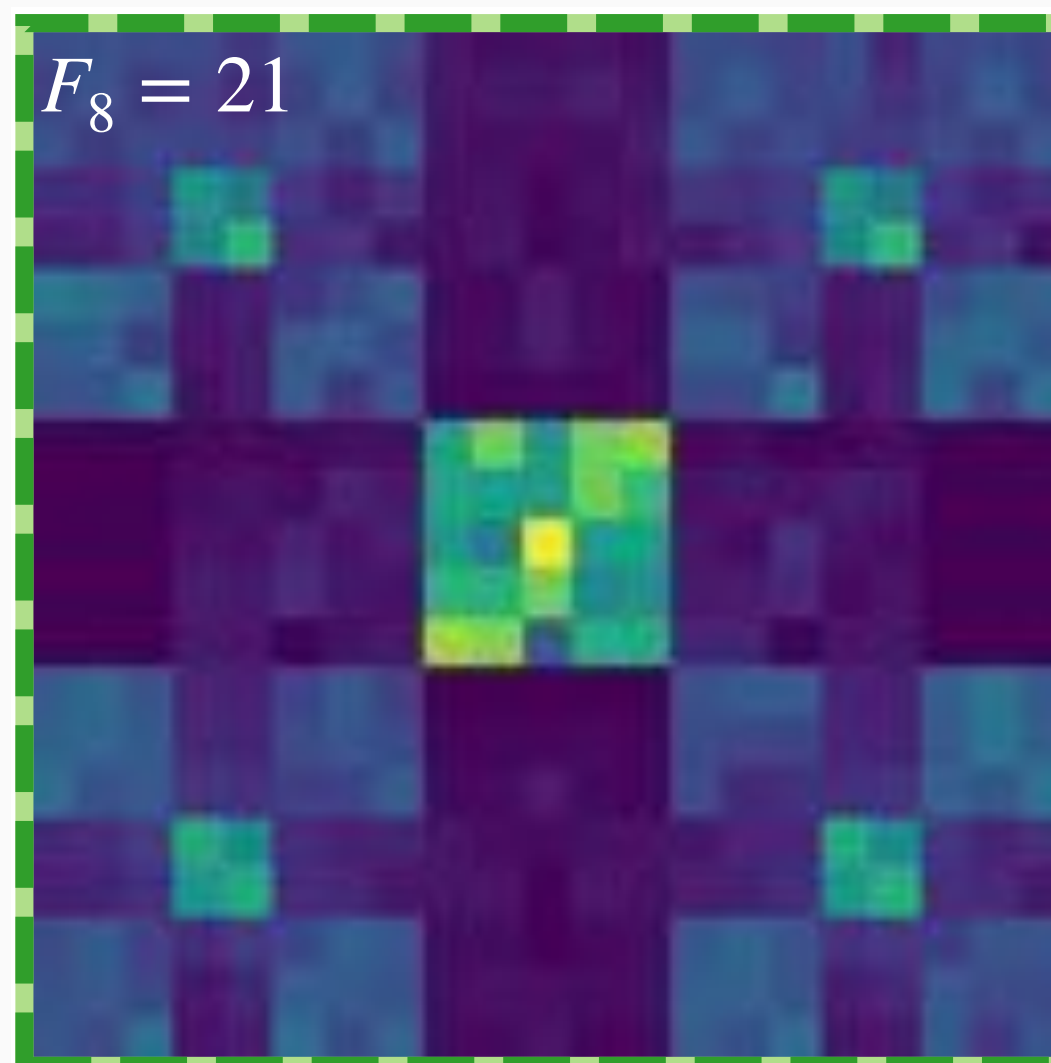
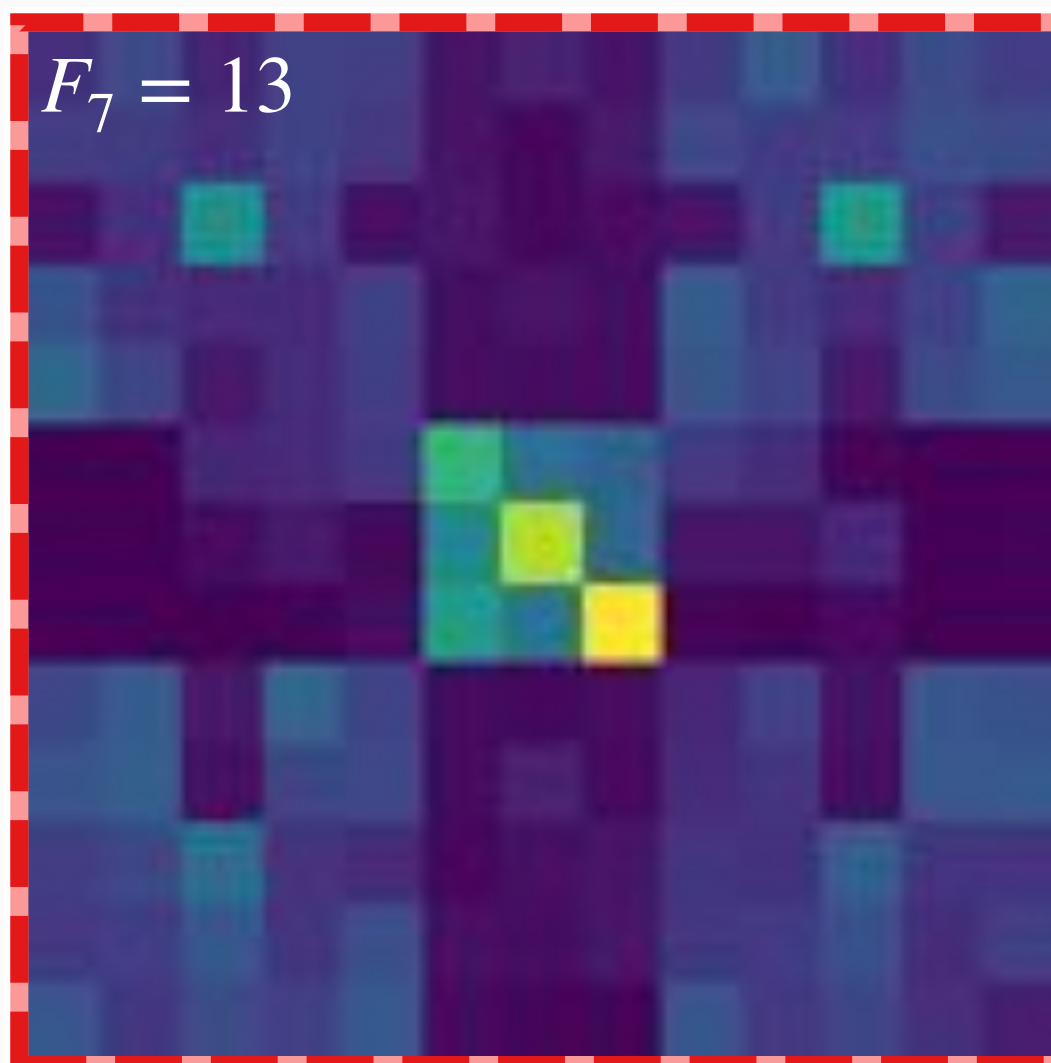
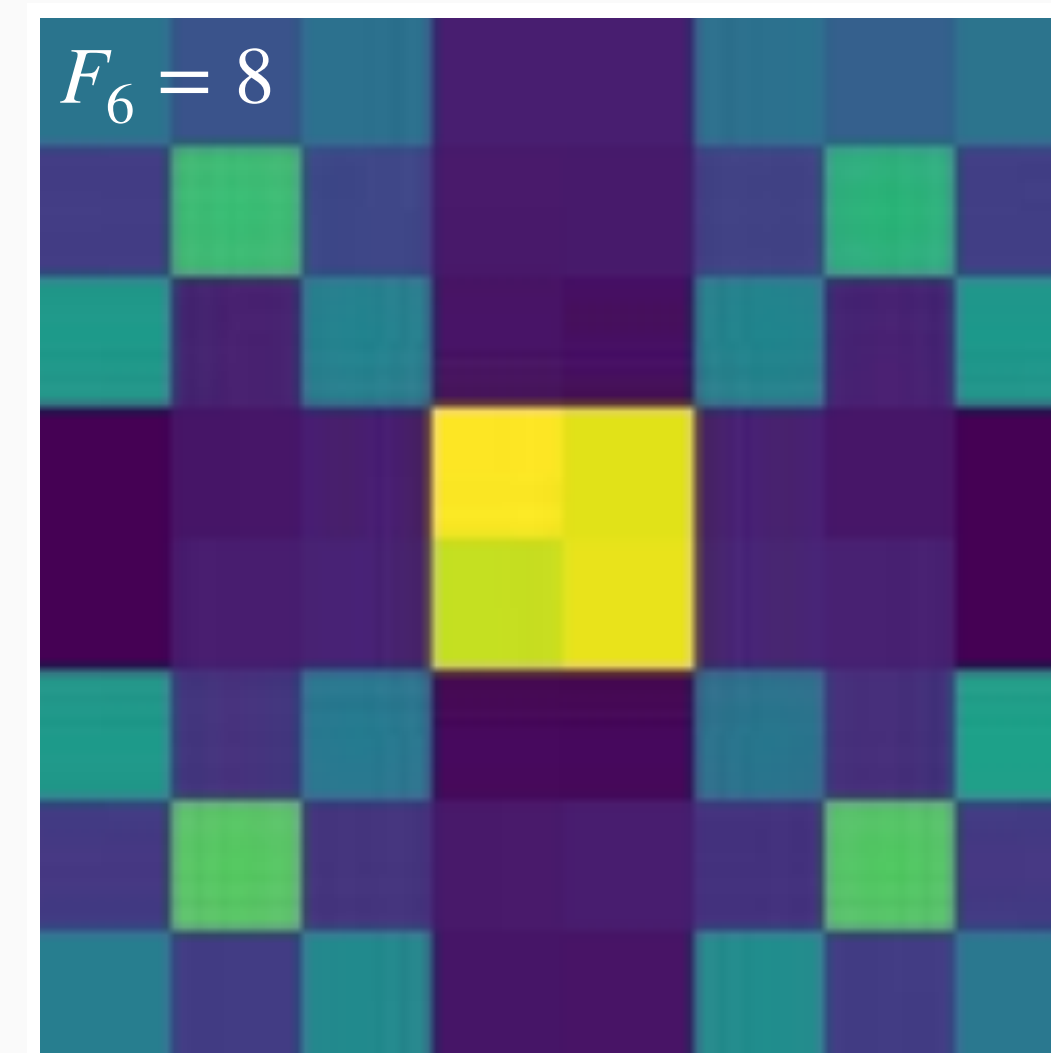
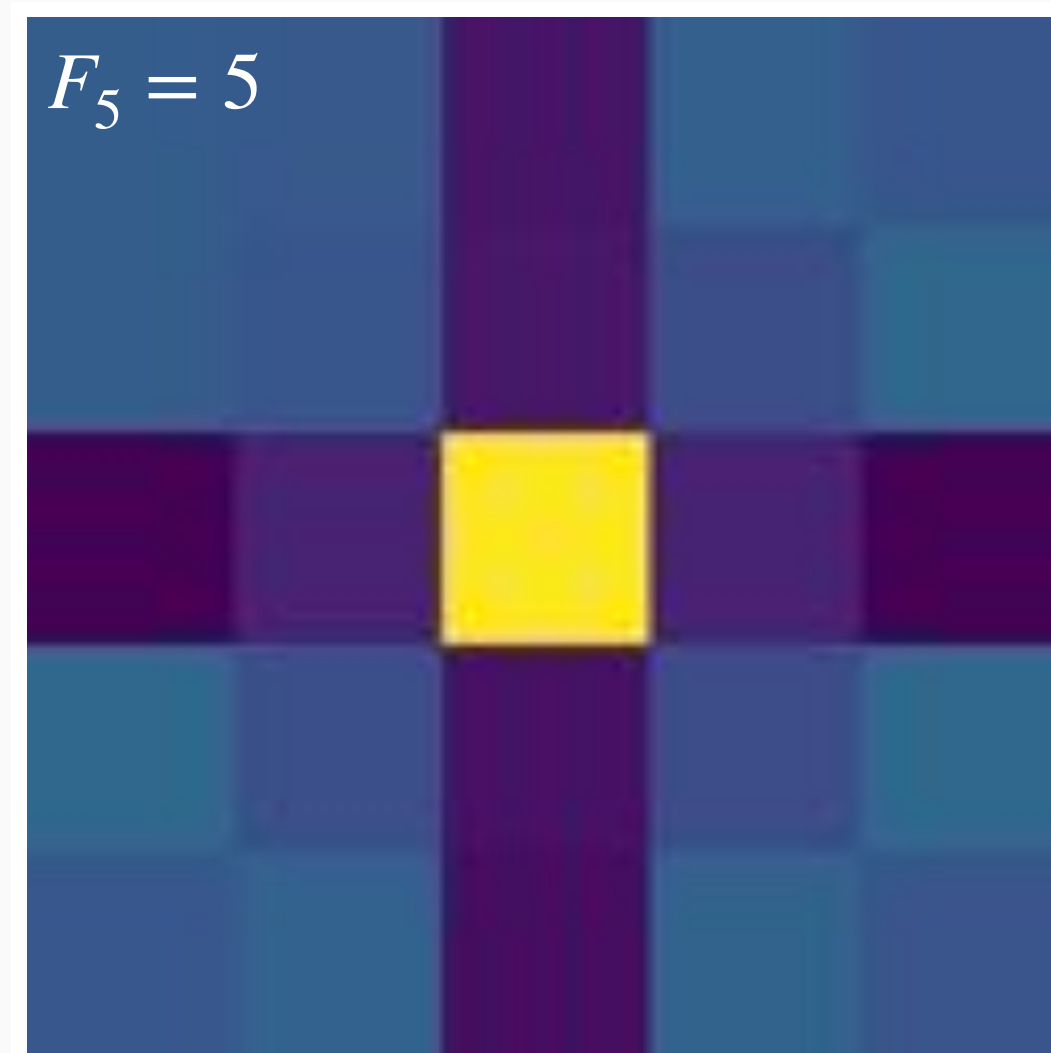
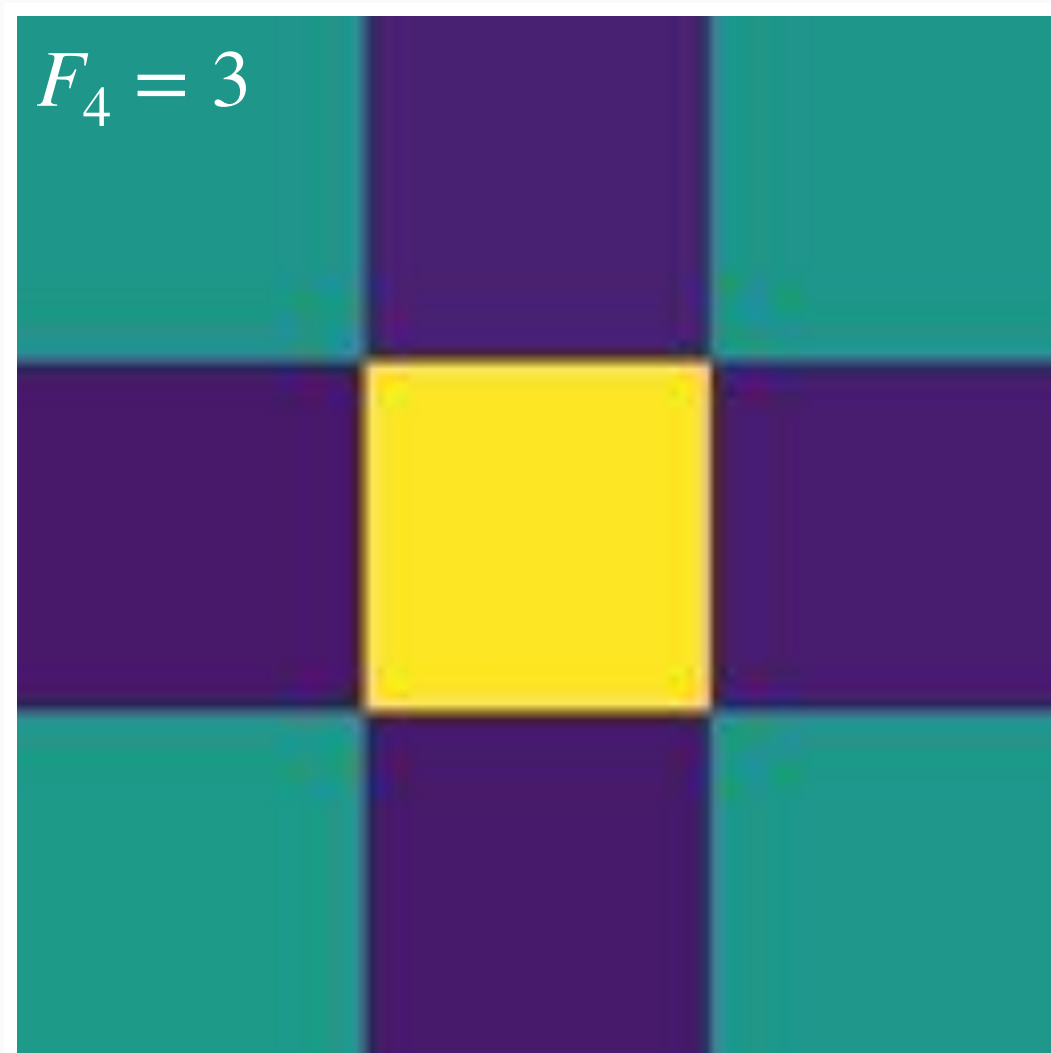


Different method to experimentally realise the periodic approximates of infinite length

- small motif length  $\rightarrow$  few permutations
- basic motif ( $F_7 = 13$ ) repeated 8 times
- ring of  $\approx 100$  resonators
- no edge states
- frequency-bands get populated
- average over periodic sites



# Study of the Recursive Construction with Circular Chains



- **atomic** sites/energies:  
 $\psi^{(n)} = \tilde{\lambda} \cdot \psi^{(n-3)}$

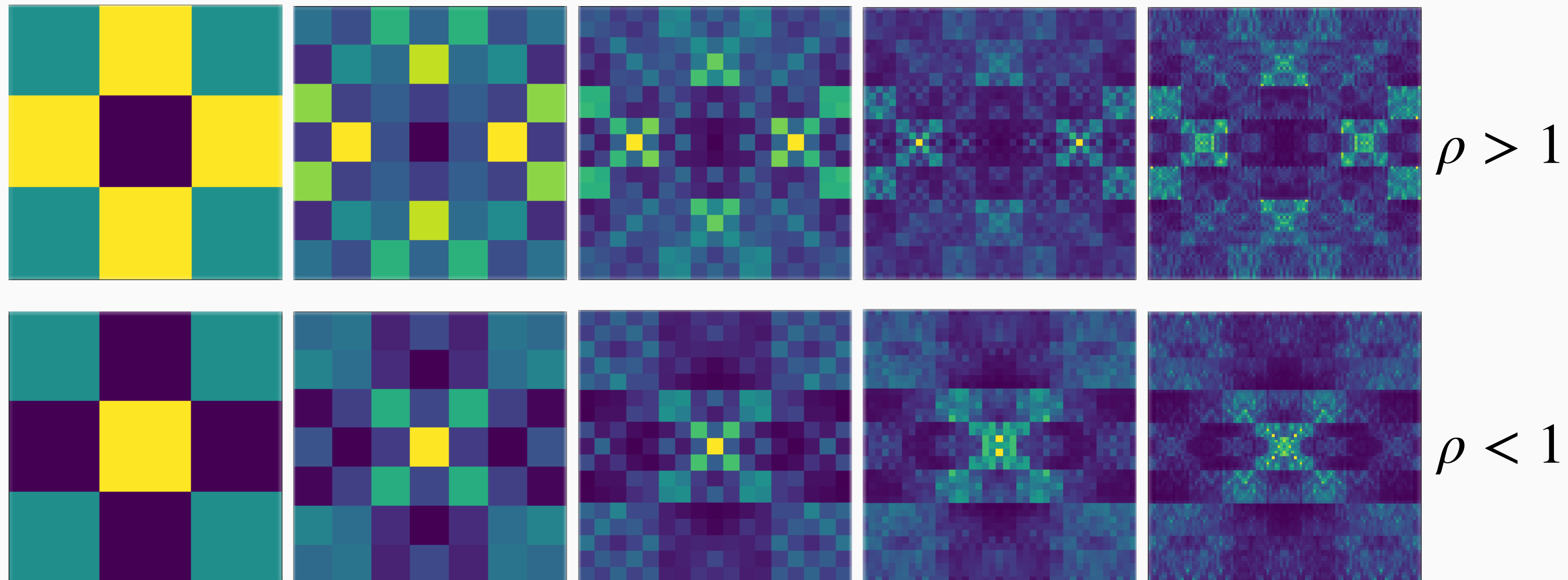
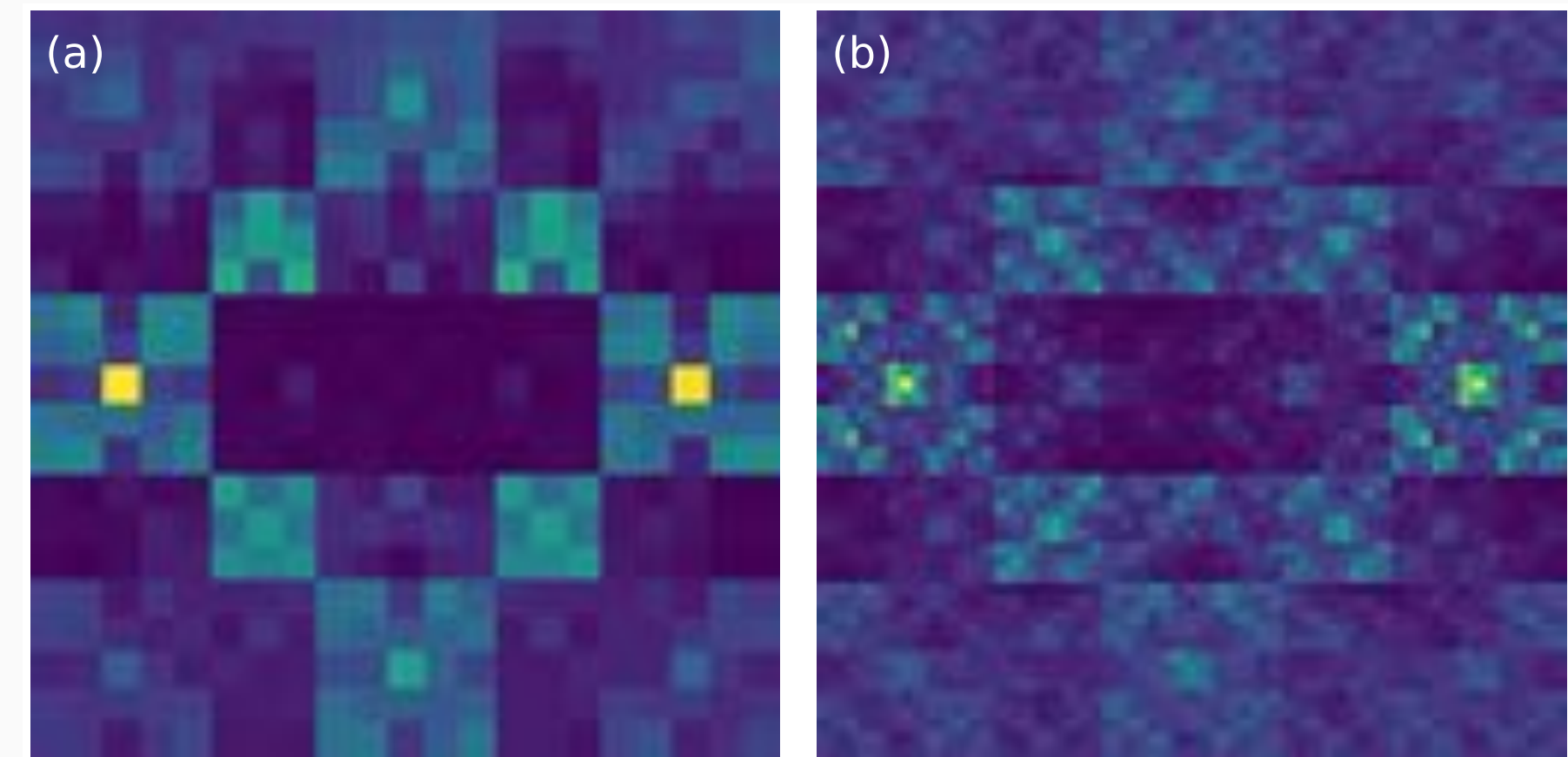
- **molecular** sites/energies:  
 $\psi^{(n)} = \lambda \cdot \psi^{(n-3)}$

# Outlook

- Interchange weak and strong couplings:  $t_A \rightleftharpoons t_B \rightarrow \rho > 1$

- other metallic means: silver mean, bronze mean, ....

$$\rightarrow F_{n+1} = m \cdot F_n + F_{n-1}$$



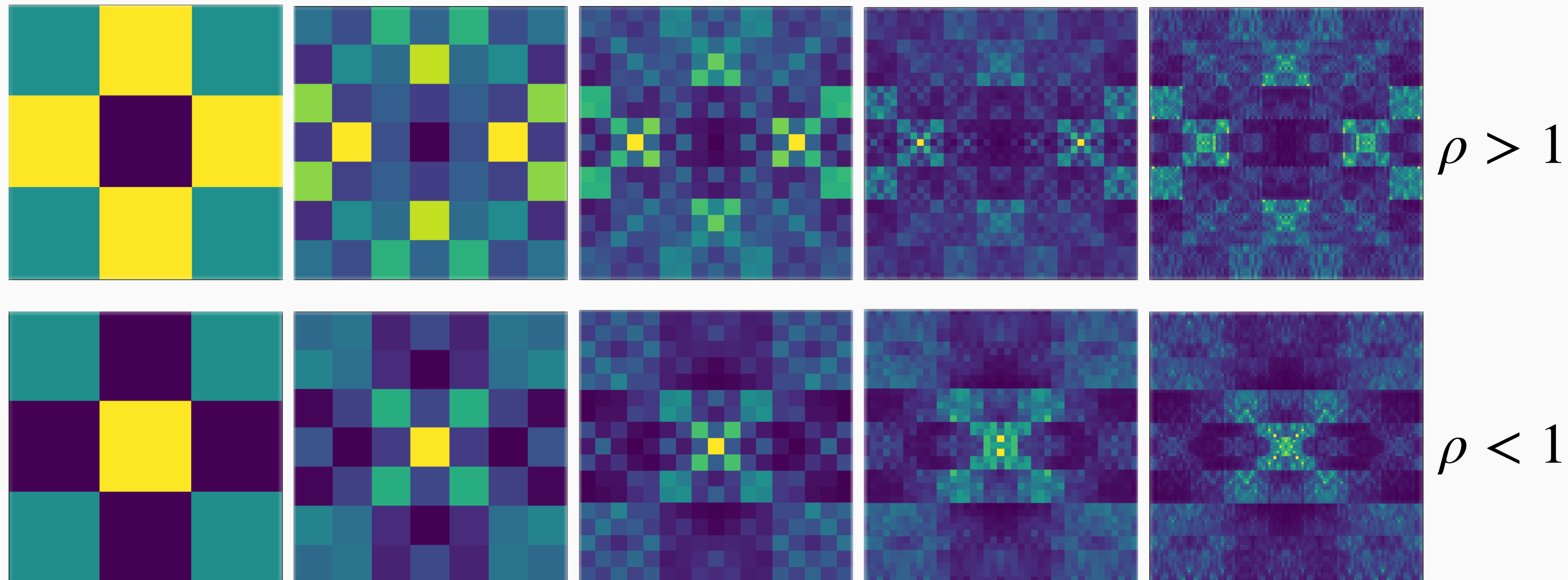
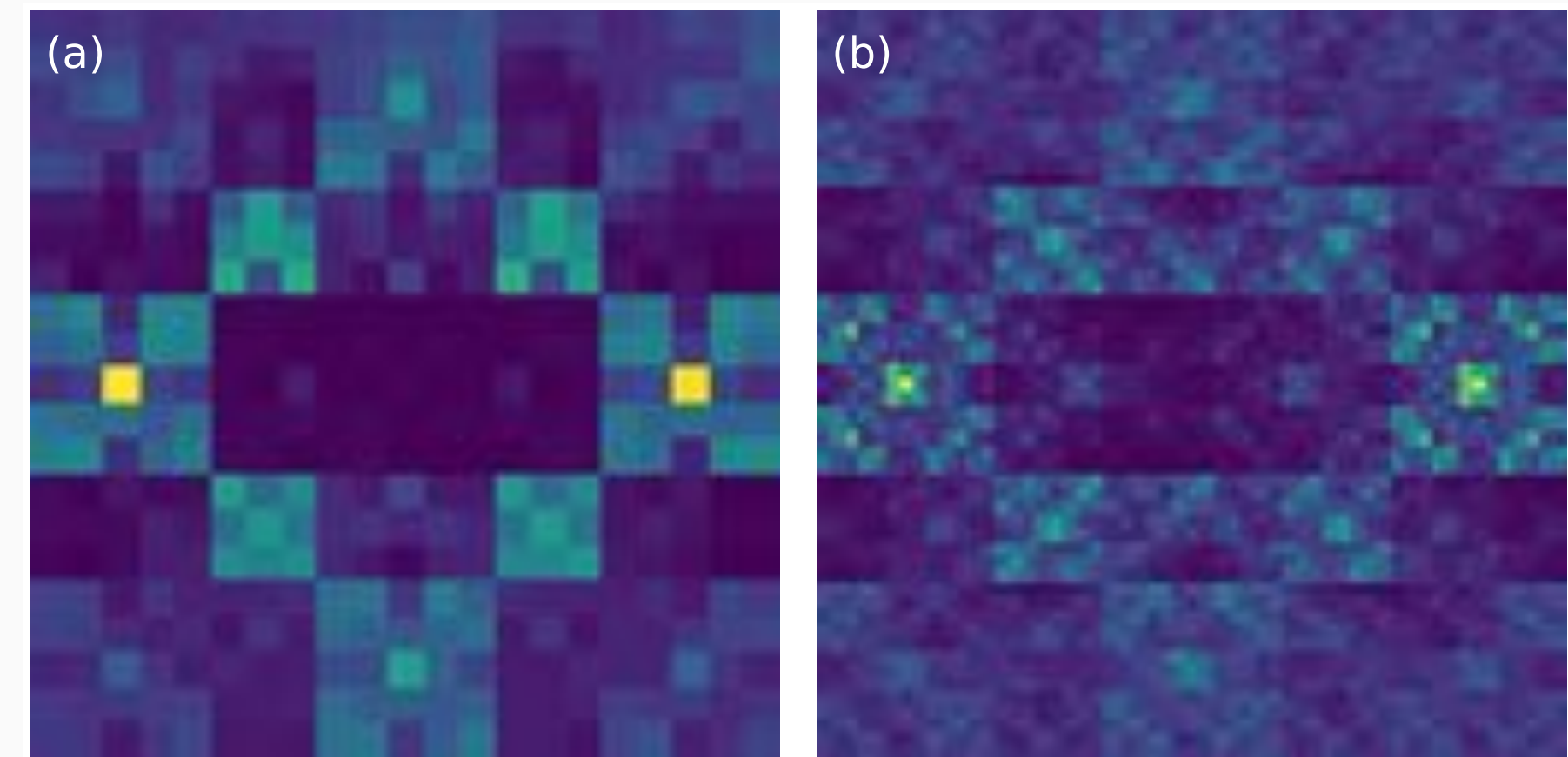


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**Thank you for  
your attention!**