

A Bethe ansatz approach to the Thirring Quantum Cellular Automaton

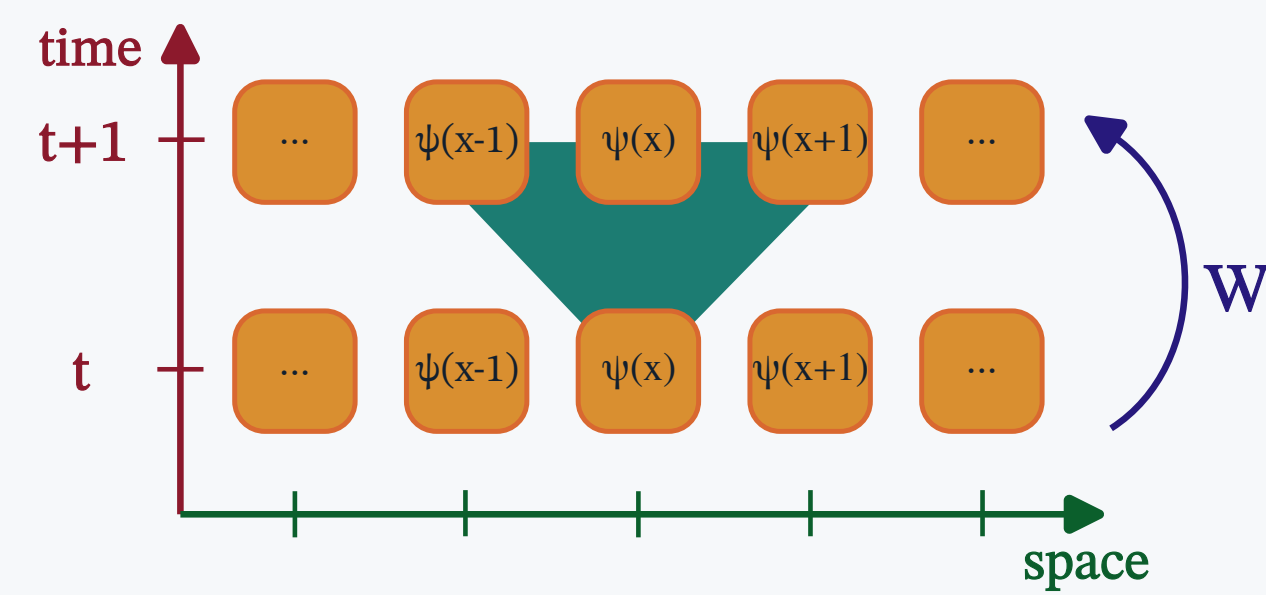
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What is a Quantum Cellular Automaton?

QCA: most general

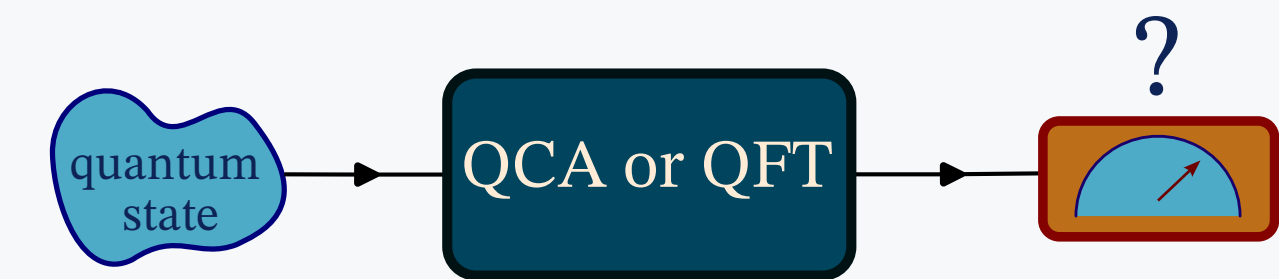
- unitary
- local
- time-discrete

dynamics of a lattice of quantum systems



Why?

- Algorithms for simulating Quantum Physics;
- Quantum Theory from principles of information processing;
- Discrete Quantum Field Theory model?



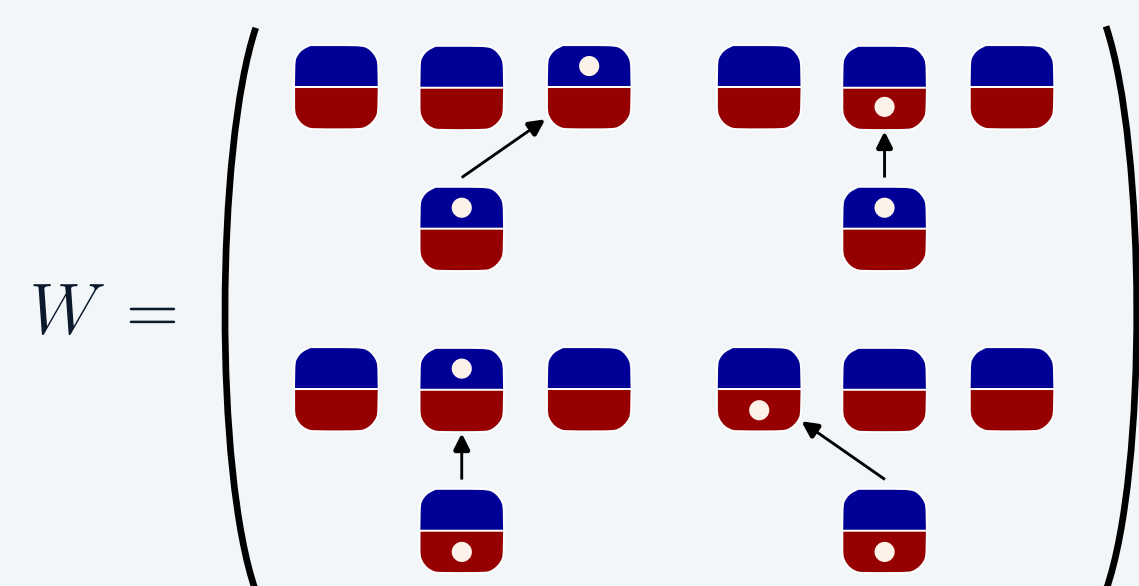
Free Theory: Dirac QCA^[1]

Assumptions:

- isotropy
- homogeneity
- minimal internal dimension of the cells

$$\text{cell} = \mathbb{C}^2 \simeq \mathbb{C}^2$$

Update rule: either flip or shift



$$W = \begin{pmatrix} \nu T & -i\mu \\ -i\mu & \nu T^\dagger \end{pmatrix} \in \text{Lin}[\mathbb{C}^2 \otimes l^2(\mathbb{Z})] \xrightarrow{\text{Fourier transform}} \tilde{W}(k) = \begin{pmatrix} \nu e^{ik} & -i\mu \\ -i\mu & \nu e^{-ik} \end{pmatrix}$$

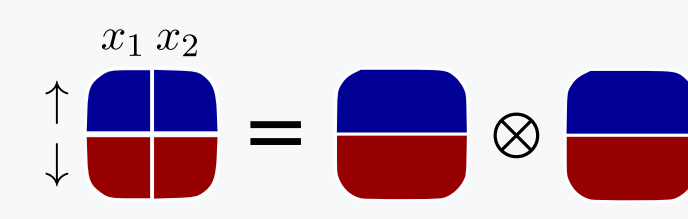
In the low-mass (μ) and low-momentum (k) regime, it recovers the dynamics described by the Dirac equation of Quantum Field Theory.

Introducing interactions: Thirring QCA^[2,3]

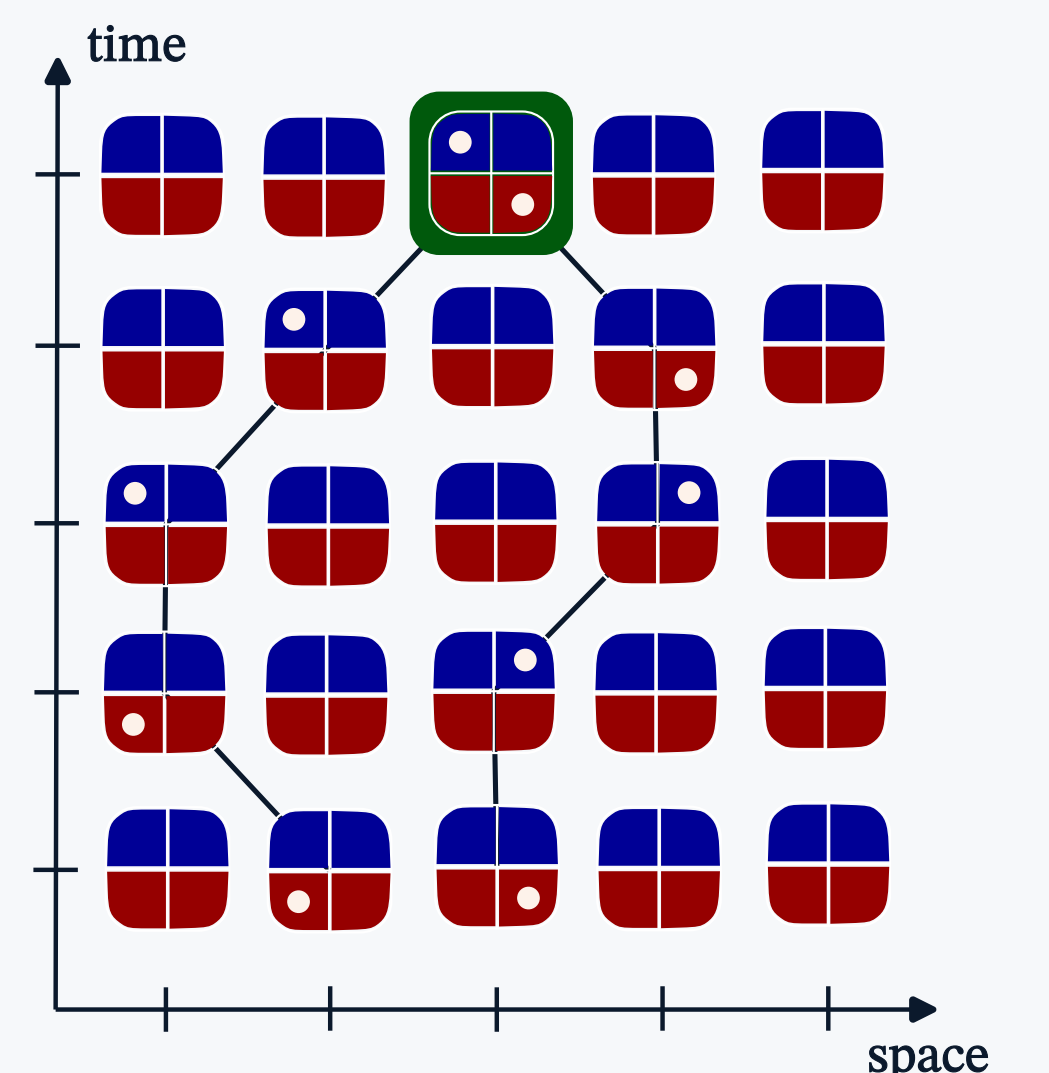
- QCA counterpart of Thirring model;
- most general on-site number preserving interaction.

$$U = J(\chi)W$$

$$J(\chi) = \begin{cases} e^{i\chi} & \text{if } x_1 = x_2, a_1 \neq a_2 \\ 1 & \text{else} \end{cases}$$



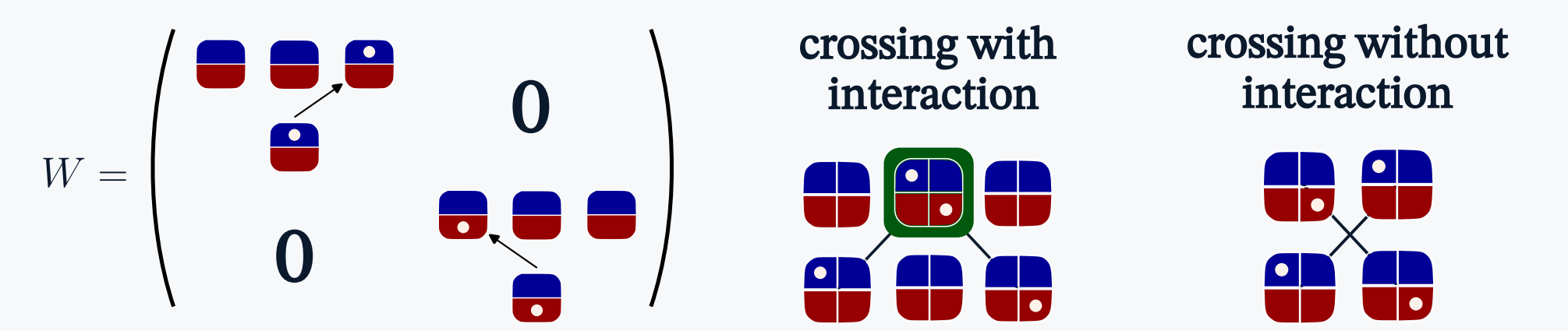
- Whenever two particles occupy
- the same lattice site;
- with opposite internal degrees of freedom; they interact.



Bethe Ansatz solution

Aim: Find eigenstates of the interacting eigenvalue equation for the massless Thirring QCA.

$$J(\chi)W |\psi(\mathbf{k}, \mathbf{a})\rangle = e^{-i\omega(\mathbf{k}, \mathbf{a})} |\psi(\mathbf{k}, \mathbf{a})\rangle$$

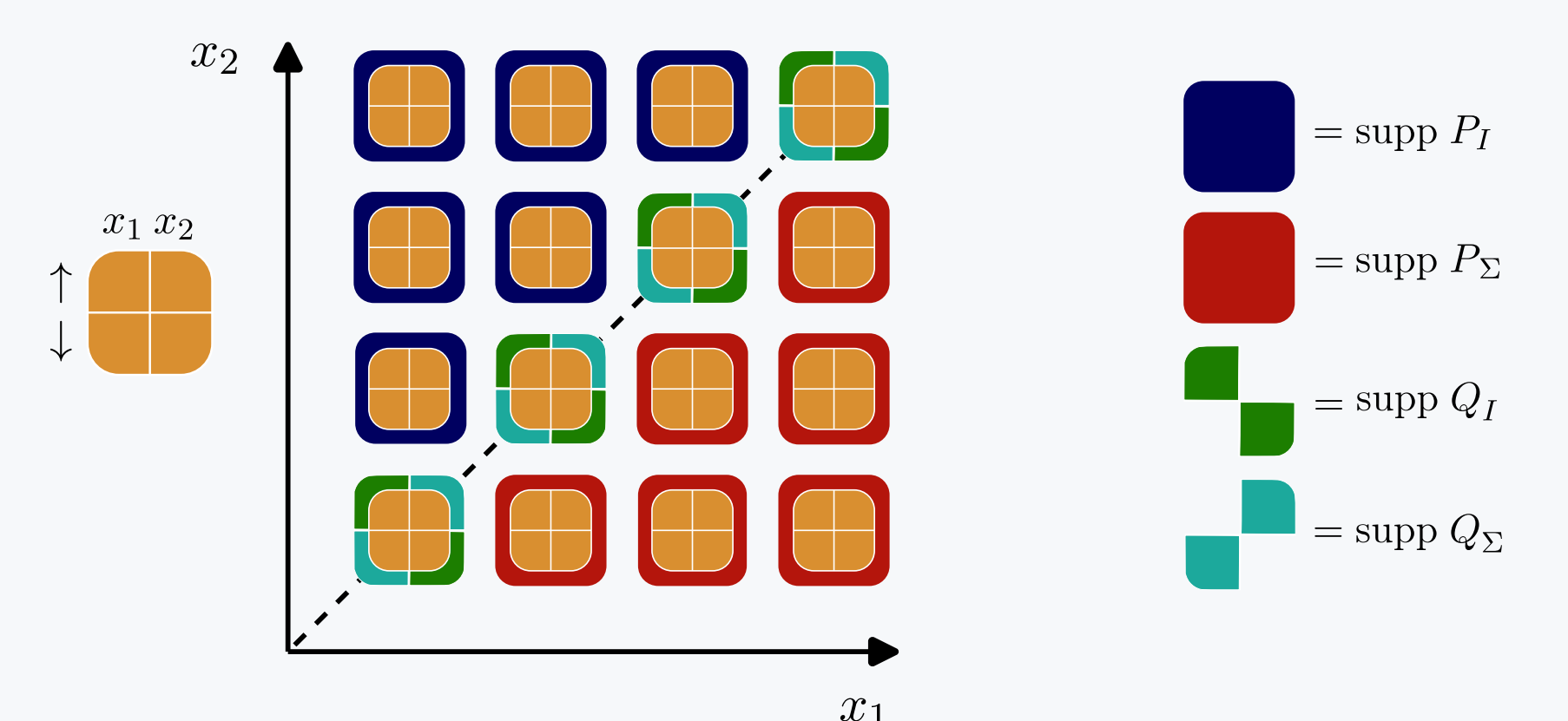


Ansatz: Anti-symmetric superposition of same-energy free eigenstates (i.e. plane waves).

$$|\psi(\mathbf{k}, \mathbf{a})\rangle = \sum_{\sigma, \tau \in S_N} \left[P_{\sigma, \tau} + \sum_{\mathbf{v} \in B^N} Q_{\sigma, \tau}(\mathbf{v}) \right] |\sigma \mathbf{k}, \sigma \mathbf{a}\rangle$$

P and Q partition the coordinate space, based on the relative ordering of the particles.

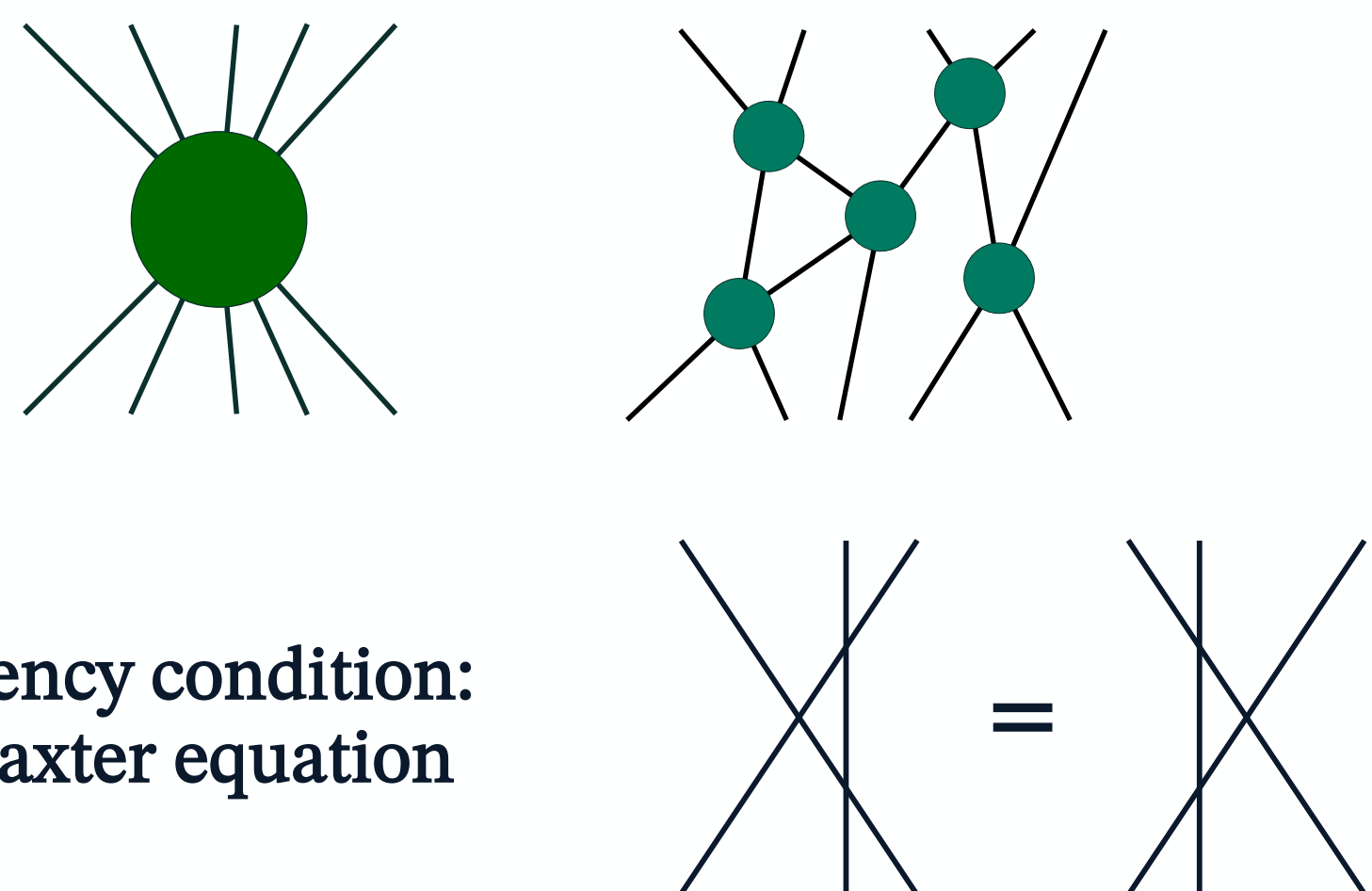
$$|\psi(\mathbf{k}, \mathbf{a})\rangle = |\psi\rangle_I + |\psi\rangle_\Sigma + |\psi\rangle_{\partial I} + |\psi\rangle_{\partial \Sigma}$$



What is quantum integrability?

No formal definition, loosely speaking, we call a quantum system integrable if:

$$\text{Many-body interactions} = \prod (\text{Pairwise interactions})$$



Consistency condition: Yang-Baxter equation

Solutions: We map the solutions on a directed graph $\Gamma_{\mathbf{a}} := \{\Gamma_{\mathbf{a}}(\sigma, \tau)\}_{\sigma, \tau \in S_N}$.

Each sub-graph is associated with a particle ordering and a permutation of spins and momenta:

$$\Gamma_{\mathbf{a}}(\sigma, \tau) = (V, E) \quad V := \{\text{particle orderings}\} \quad E := \{\text{orderings connected by a single QCA step}\}$$

Edges encode relations between bulk and boundary coefficients:

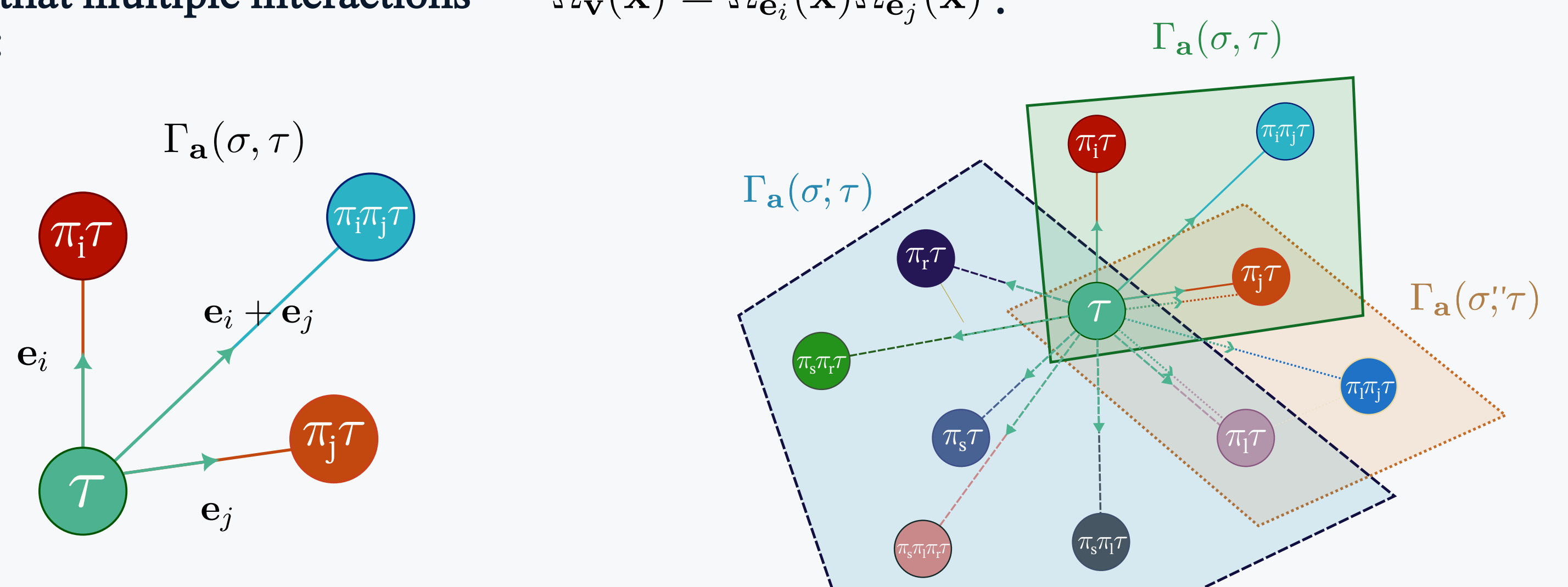
$$(\tau, \pi_{\mathbf{e}_i} \tau) \in E \iff \begin{cases} f_{\sigma, \pi_{\mathbf{e}_i} \tau}(\mathbf{x}) = \Omega_{\mathbf{e}_i}(\mathbf{x}) f_{\sigma, \tau}(\mathbf{x}), \\ g_{\sigma, \tau}^{\mathbf{e}_i}(\mathbf{x}) = e^{-i|\mathbf{e}_i|^2 \chi} f_{\sigma, \tau}(\mathbf{x}). \end{cases}$$

The composition rule

$$\mathbf{v} = \mathbf{e}_i + \mathbf{e}_j \mapsto (\tau, \pi_{\mathbf{v}} \tau) = (\tau, \pi_{\mathbf{e}_i} \pi_{\mathbf{e}_j} \tau),$$

guarantees that multiple interactions factorise as:

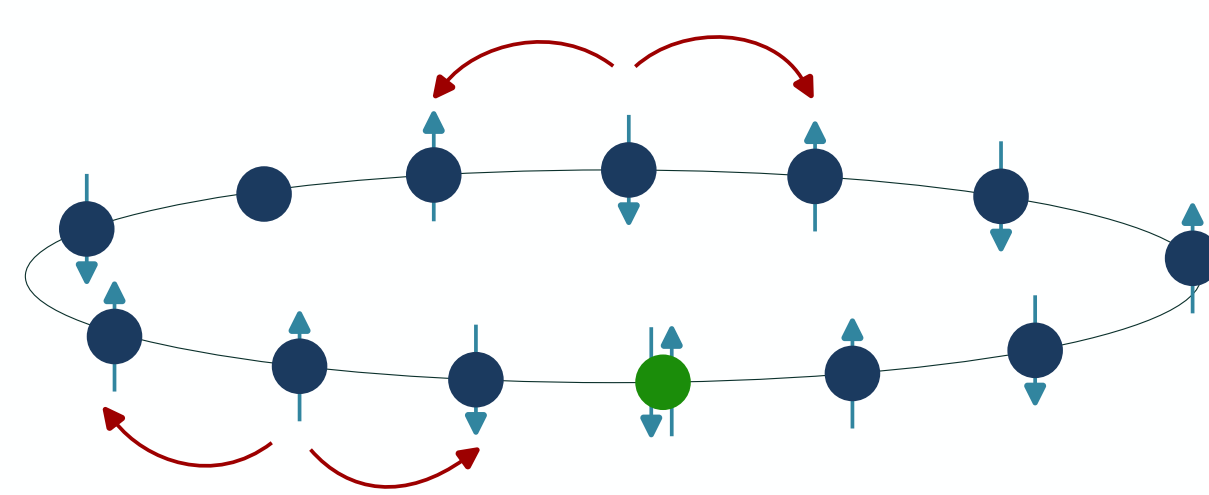
$$\Omega_{\mathbf{v}}(\mathbf{x}) = \Omega_{\mathbf{e}_i}(\mathbf{x}) \Omega_{\mathbf{e}_j}(\mathbf{x}).$$



Traditional approach (e.g. Hubbard model)

Closed spin chain with Hamiltonian

$$H = \text{Hopping term} + \text{Local interaction}$$



Approach: exploit symmetries to make an educated guess about the form of the wave function; then plug it into the Schrödinger equation to find its parameters.

Main differences of the QCA framework:

- time-discrete evolution;
- (quasi)-energy invariant mod 2π ;
- spin-dependent dynamics;
- infinite open chain.