



UNIVERSITÀ
DI PAVIA

"Repeated interactions as a thermodynamic microscope: spin rings and clustering collision models"

End of year seminars A.A 2025/2026

17/09/2026

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Supervisors: Giacomo Guarnieri
Dario Gerace

Open vs closed quantum systems

Closed quantum
systems



Open vs closed quantum systems

Closed quantum
systems



Von Neumann
equation: $\frac{d\rho_S(t)}{dt} = -i[H(t), \rho_S(t)]$



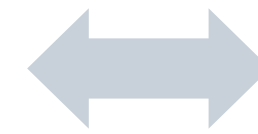
Unitary
dynamics: $\rho_S(t) = U(t)\rho_S(0)U(t)^\dagger$

Open vs closed quantum systems

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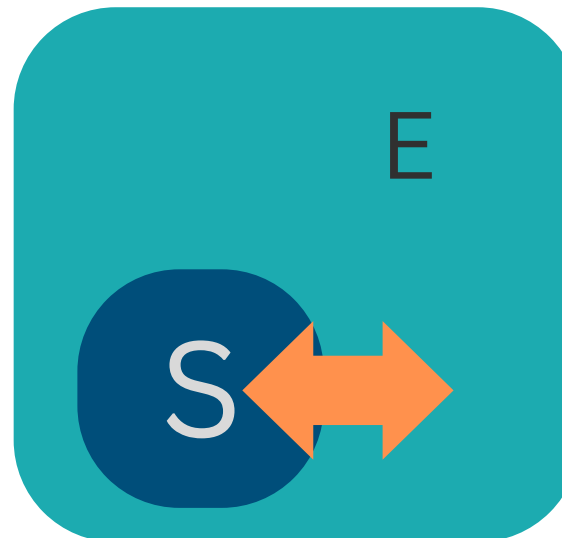
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Open quantum
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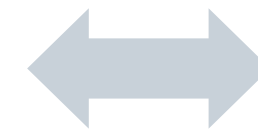


Open vs closed quantum systems

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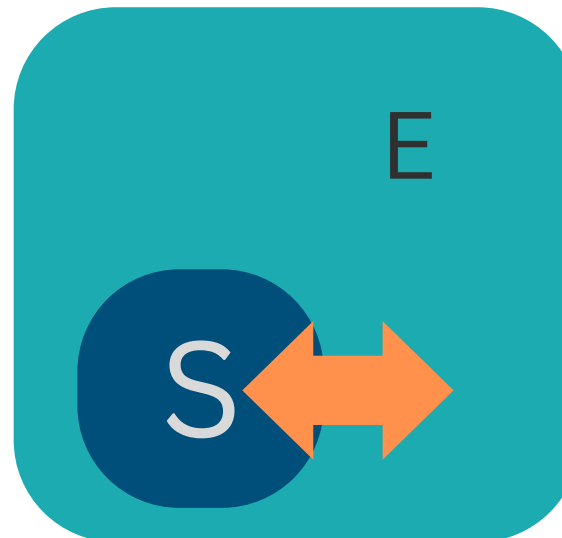
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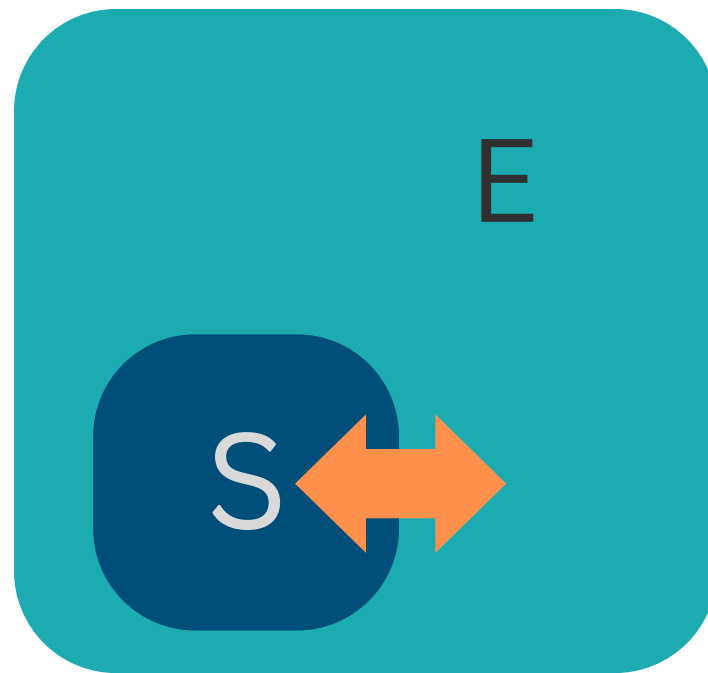
$$\rho_S(t) = U(t)\rho_S(0)U(t)^\dagger$$

Open quantum
systems



More degrees of freedom to describe
→ the formalism is not immediate

Lindblad master equations



reduced dynamics:

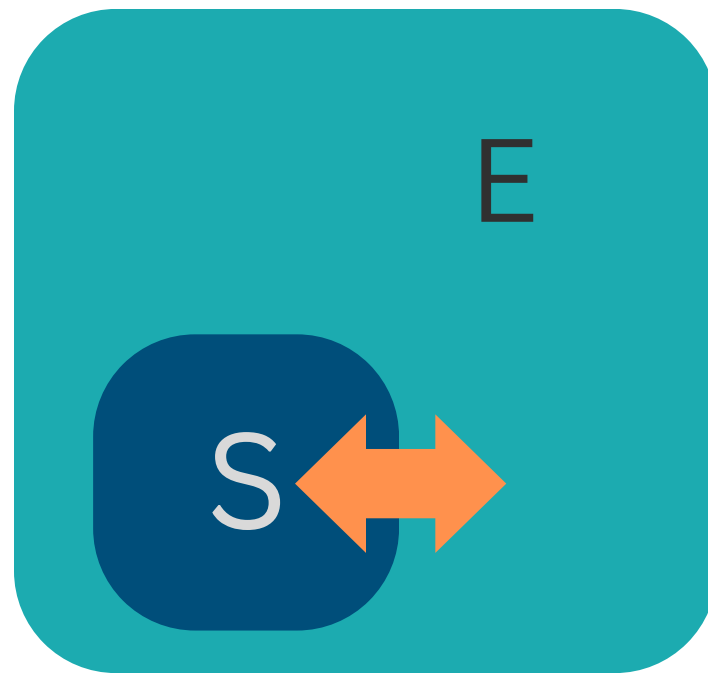
$$\frac{d\rho_S(t)}{dt} = -i[H_S(t) + H_{LS}(t), \rho_S(t)] + \mathcal{D}_t[\rho_S(t)]$$

UNITARY CONTRIBUTION

$$\mathcal{D}_t[\rho_S(t)] = \sum_k \gamma_k(t) \left(S_k(t) \rho_S(t) S_k^\dagger(t) - \frac{1}{2} \{ S_k^\dagger(t) S_k(t), \rho_S(t) \} \right) \quad \gamma_k(t) \in \mathbb{R}$$

DISSIPATOR

Lindblad master equations



reduced dynamics:

$$\frac{d\rho_S(t)}{dt} = -i[H_S(t) + H_{LS}(t), \rho_S(t)] + \mathcal{D}_t[\rho_S(t)]$$

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DISSIPATOR

Needed approximations:

Weak coupling

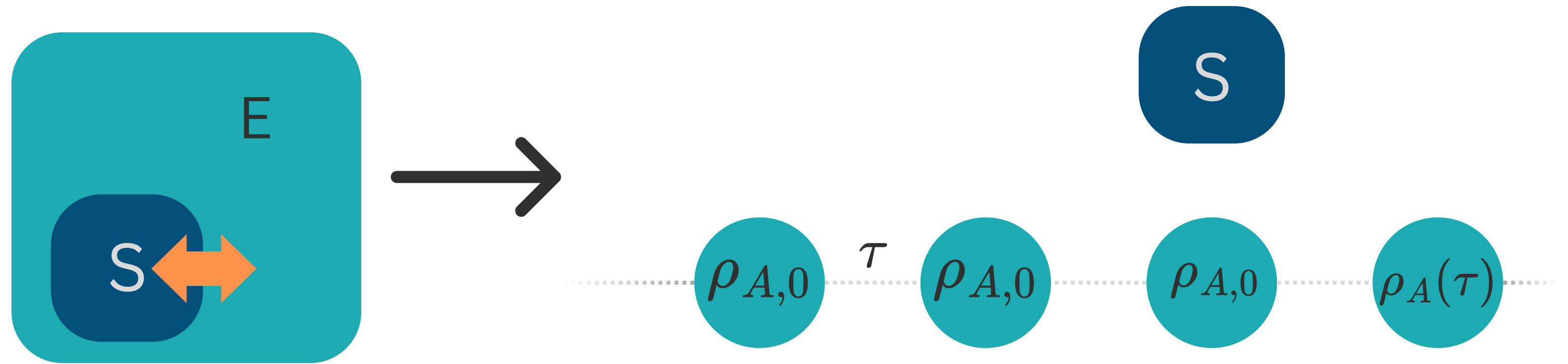
Markov approximation

Born approximation

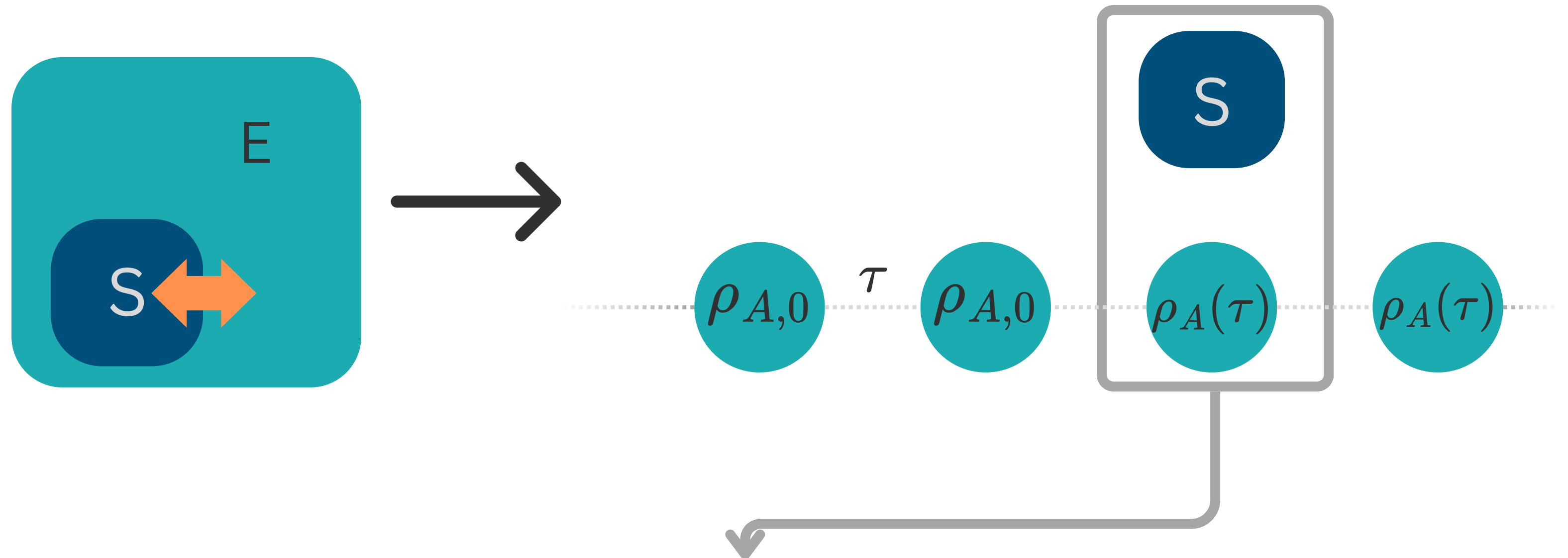
Secular approximation

3

Collision models

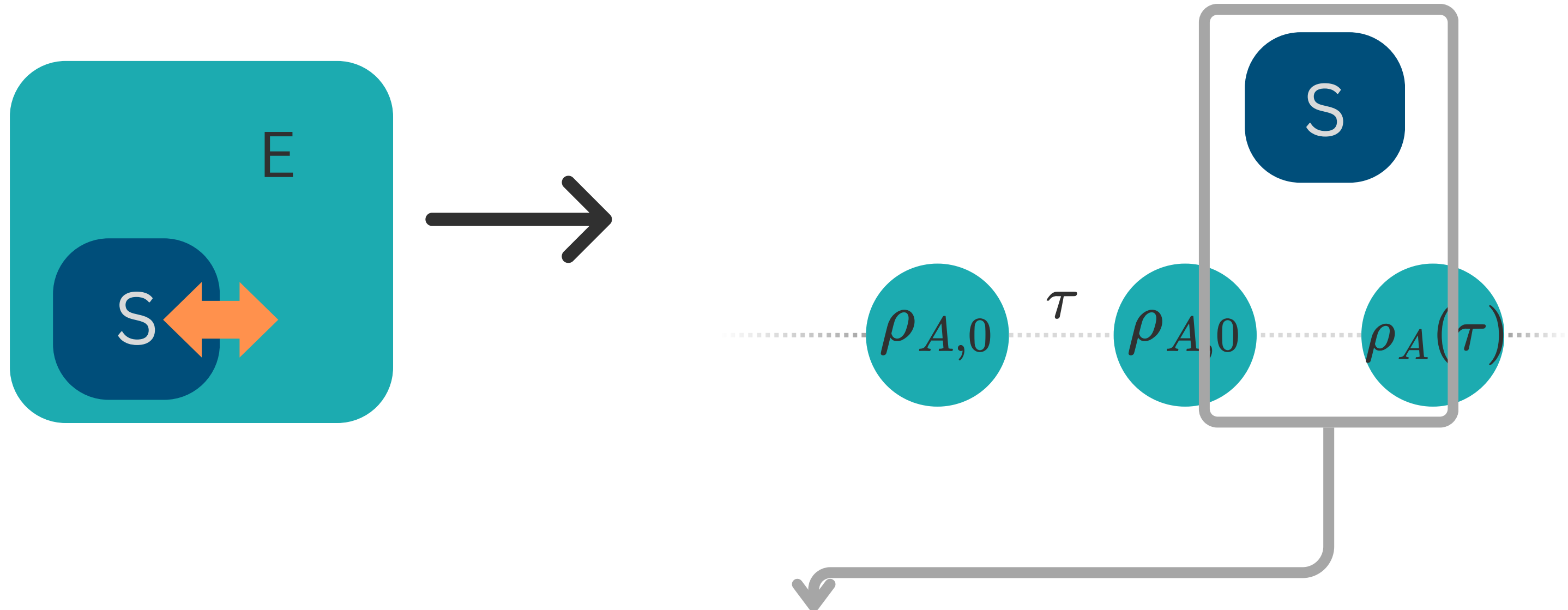


Collision models



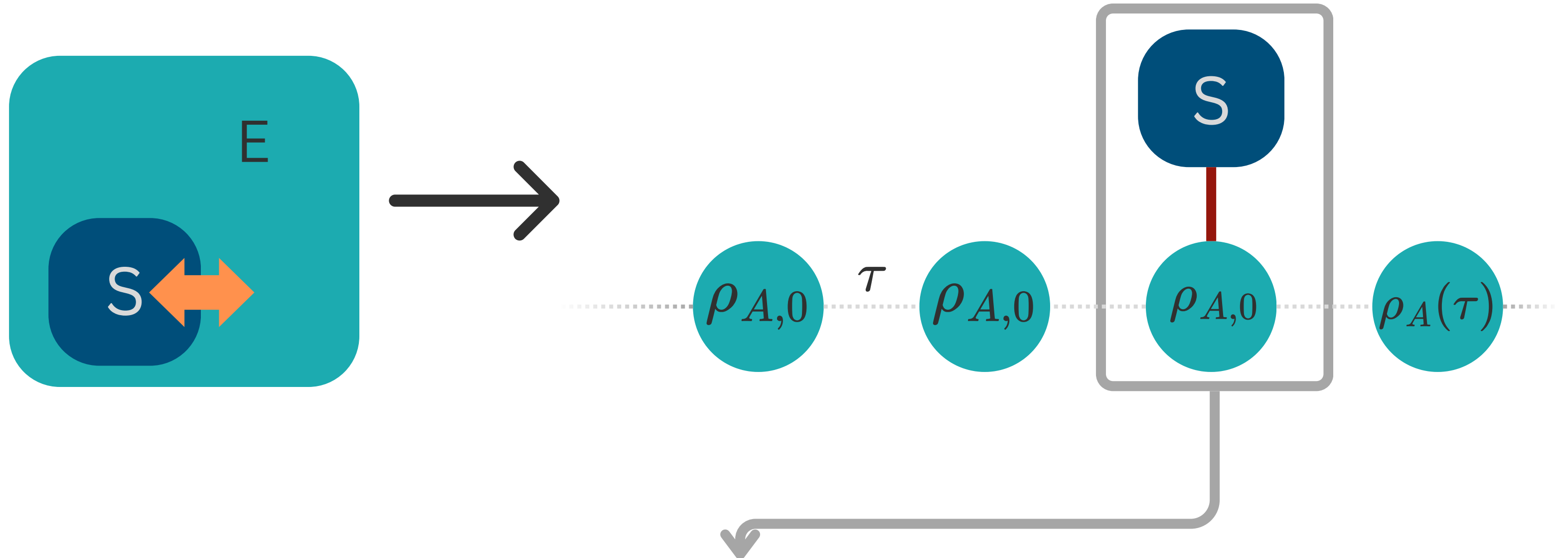
$$\rho_S(\tau) = \text{Tr}_A[U(\tau)(\rho_S(0) \otimes \rho_A(0))U(\tau)^\dagger]$$

Collision models



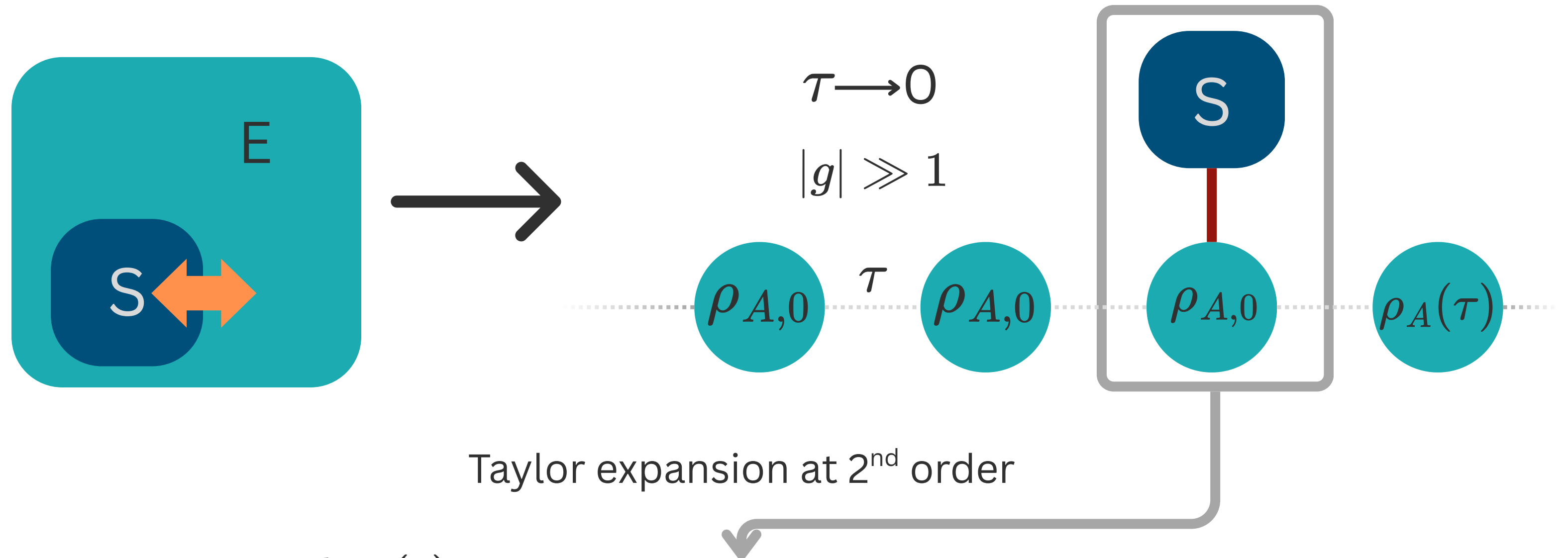
$$\rho_S(\tau) = \text{Tr}_A[U(\tau)(\rho_S(0) \otimes \rho_A(0))U(\tau)^\dagger]$$

Collision models



$$\rho_{SA}(2\tau) = U(\tau)(\rho_S(0) \otimes \rho_A(0))U(\tau)^\dagger$$

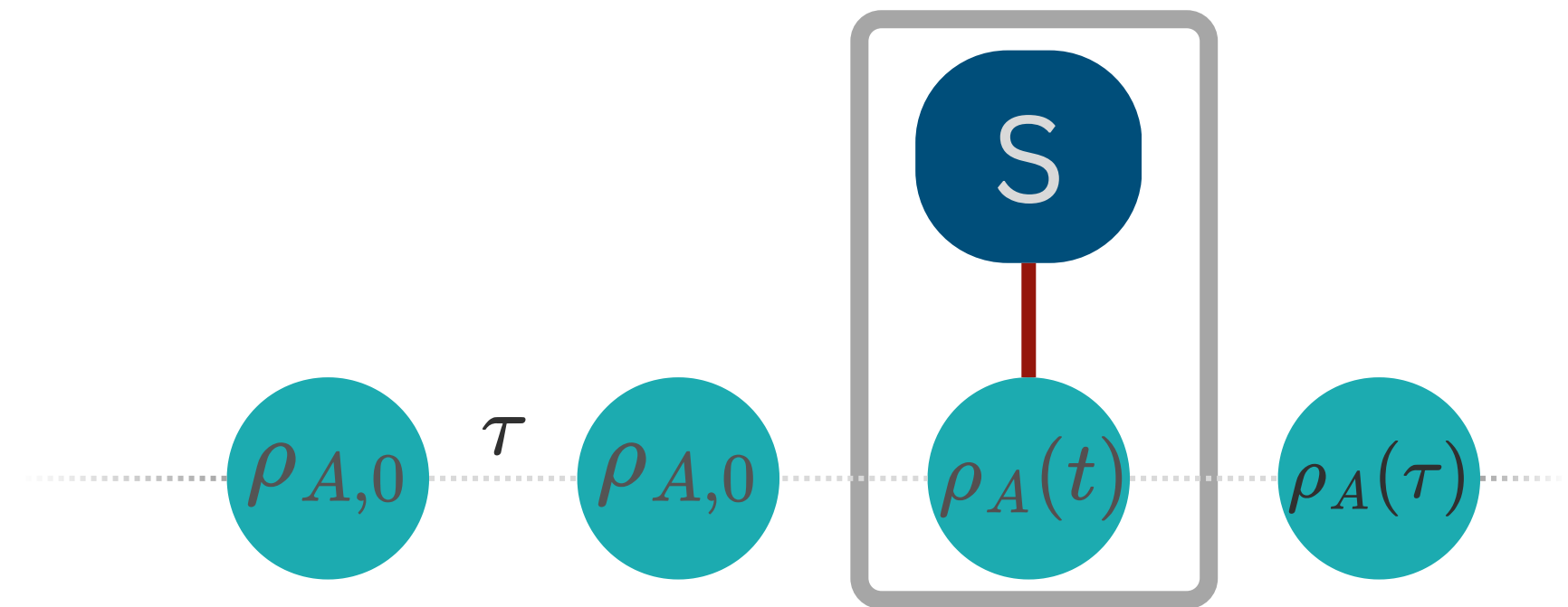
Collision models



$$\frac{d\rho_S(t)}{dt} = -i[H_S, \rho_S(t)] + \mathcal{D}[\rho_S(t)]$$

Collision models

- 1) Simple structure
- 2) No weak coupling, secular approximation



Ciccarello, Francesco, et al. Physics Reports 954 (2022): 1-70.

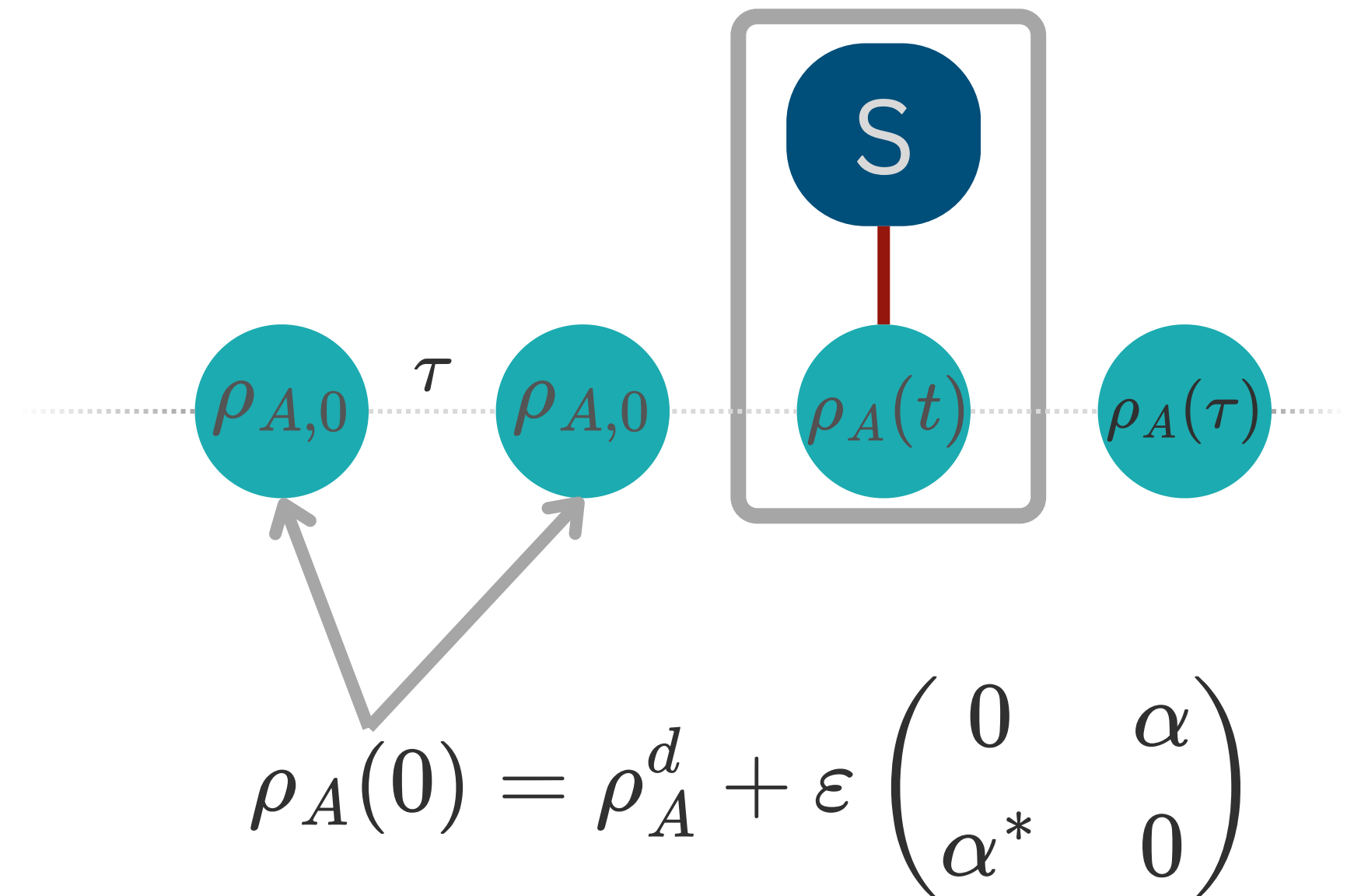
Collision models

1) Simple structure

2) No weak coupling, secular approximation

3) Versatility:

- adding coherence



Ciccarello, Francesco, et al. Physics Reports 954 (2022): 1-70.

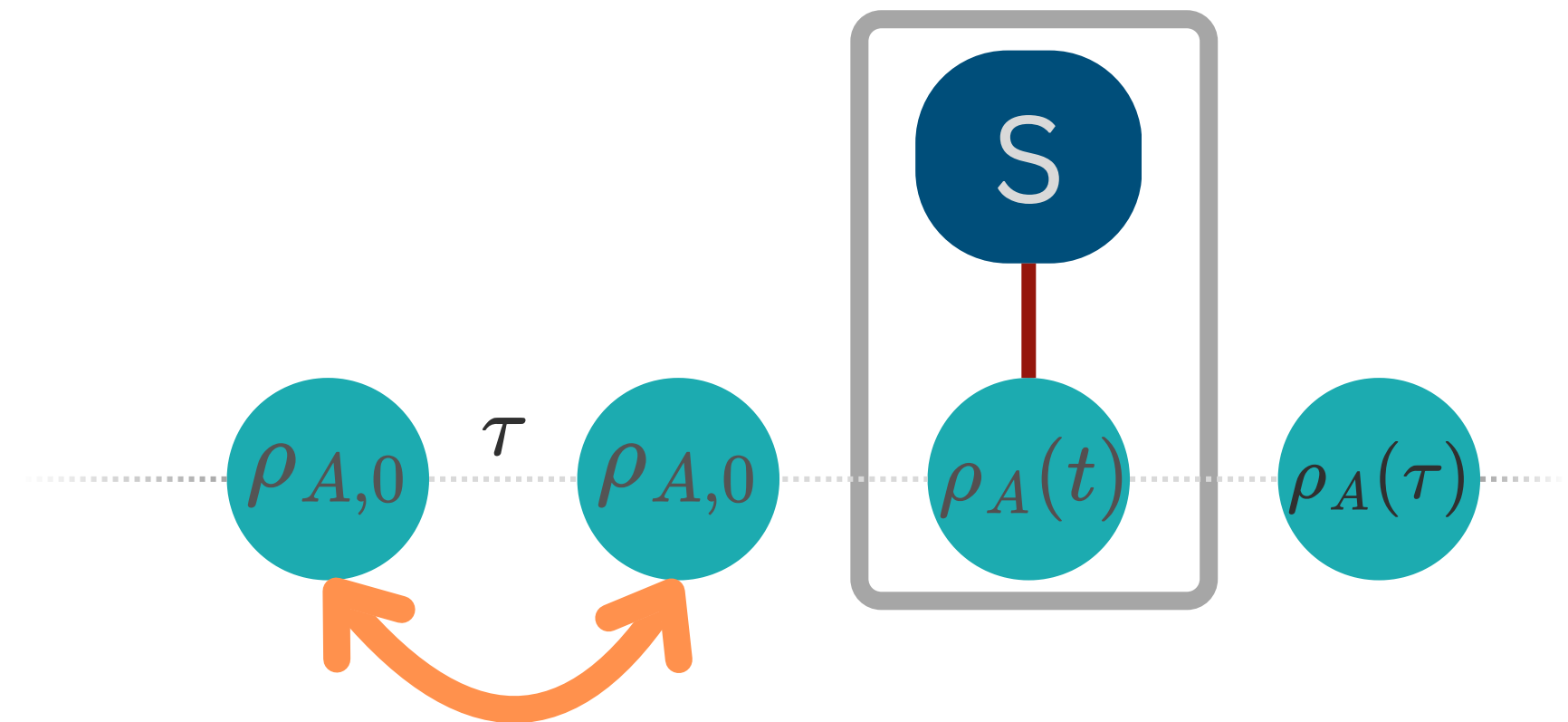
Collision models

1) Simple structure

2) No weak coupling, secular approximation

3) Versatility:

- adding coherence
- memory effects



Ciccarello, Francesco, et al. Physics Reports 954 (2022): 1-70.

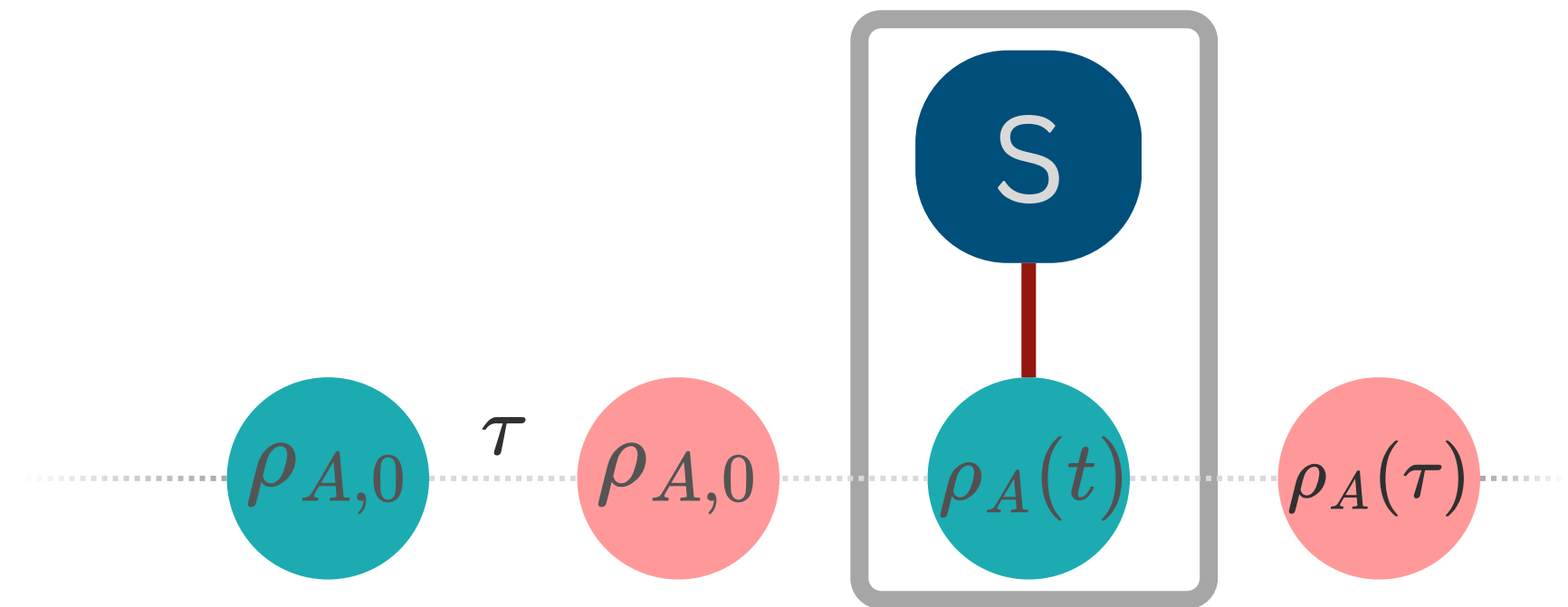
Collision models

1) Simple structure

2) No weak coupling, secular approximation

3) Versatility:

- adding coherence
- memory effects
- more than one bath



Ciccarello, Francesco, et al. Physics Reports 954 (2022): 1-70.

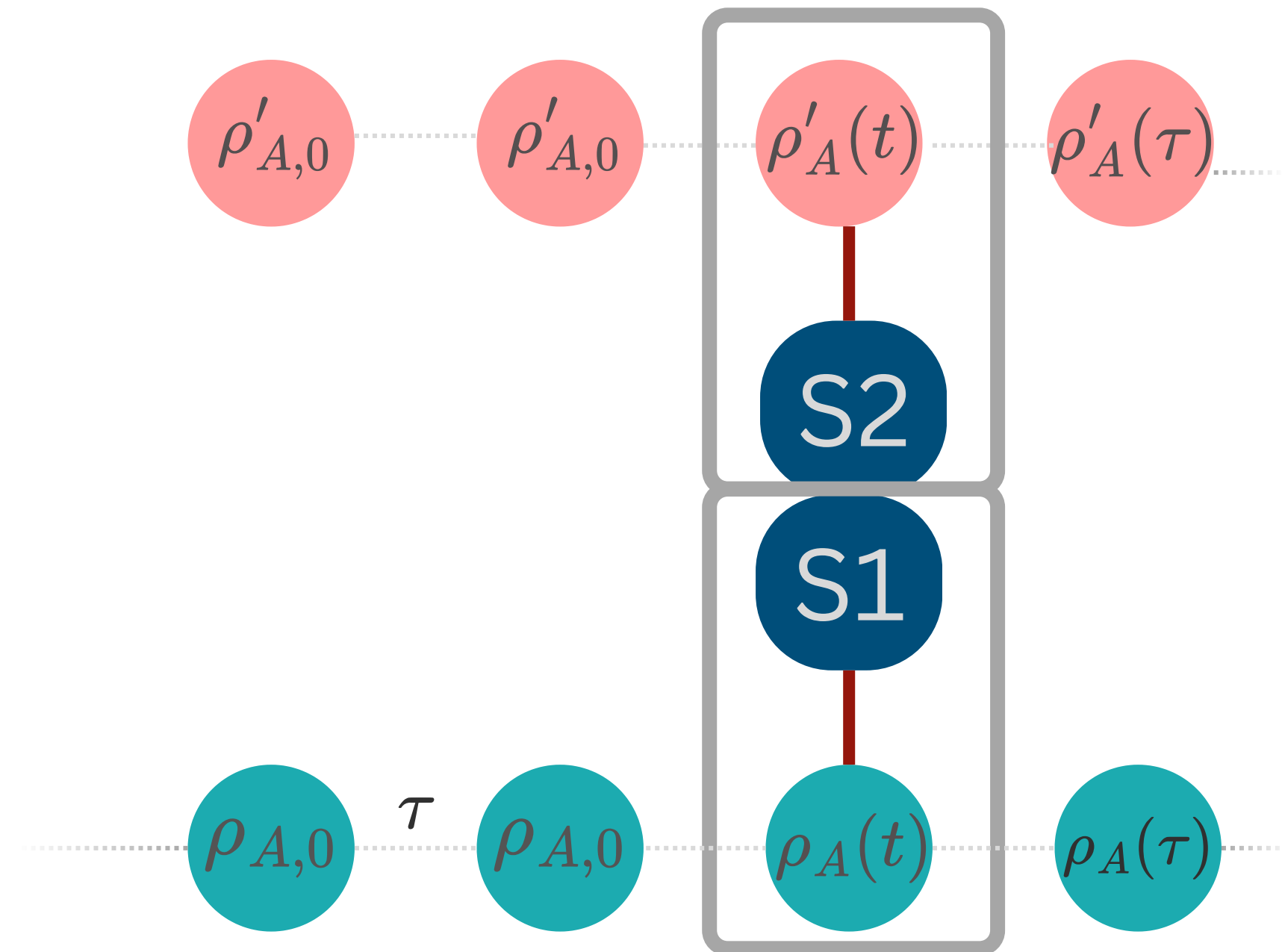
Collision models

1) Simple structure

2) No weak coupling, secular approximation

3) Versatility:

- adding coherence
- memory effects
- more than one bath
- multi-partite system

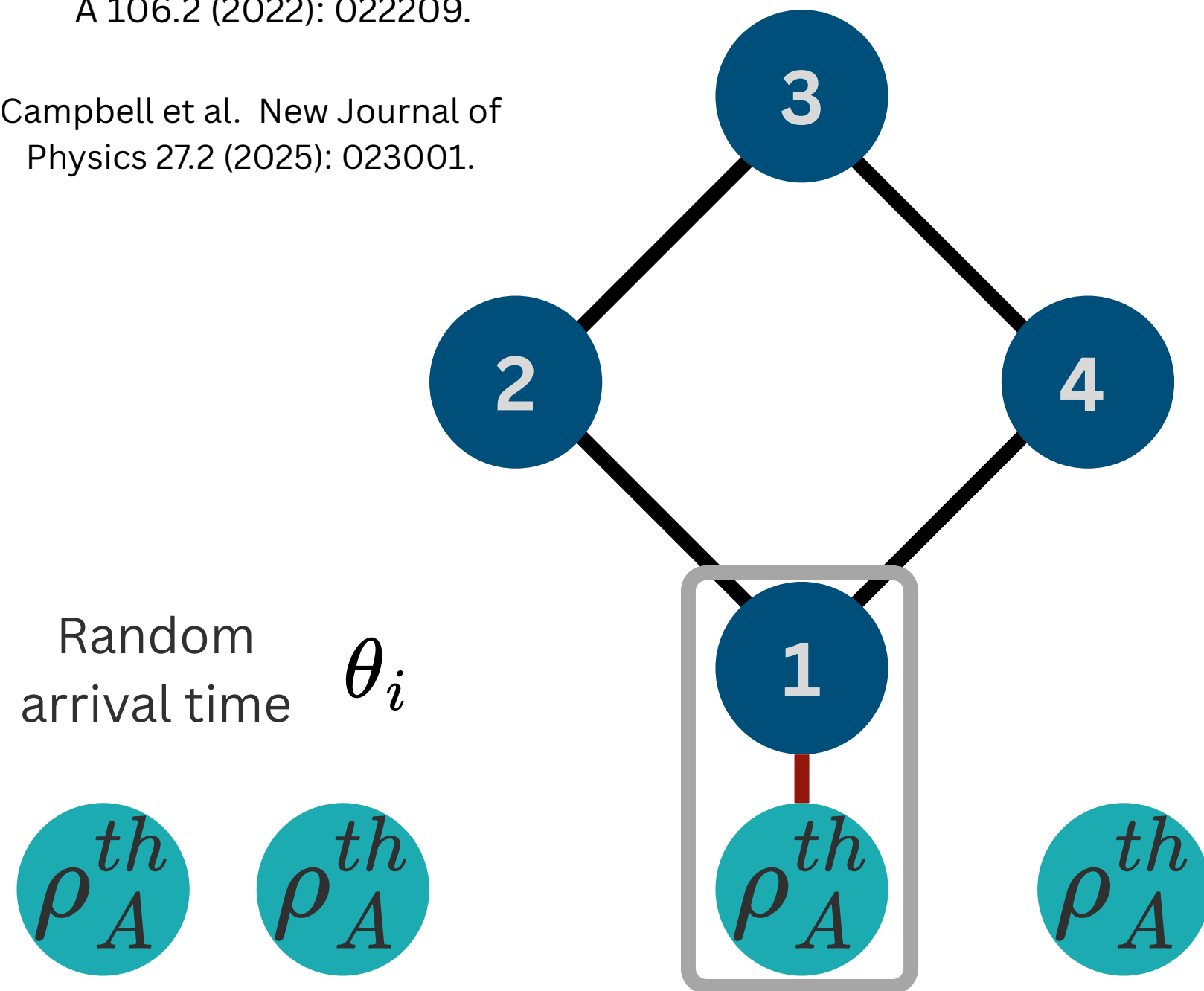


Ciccarello, Francesco, et al. Physics Reports 954 (2022): 1-70.

Heisenberg spin ring

Guarnieri, et al. Physical Review
A 106.2 (2022): 022209.

Campbell et al. New Journal of
Physics 27.2 (2025): 023001.



Hamiltonian:

$$H_{tot} = H_S + H_{SA}(t) + H_A$$

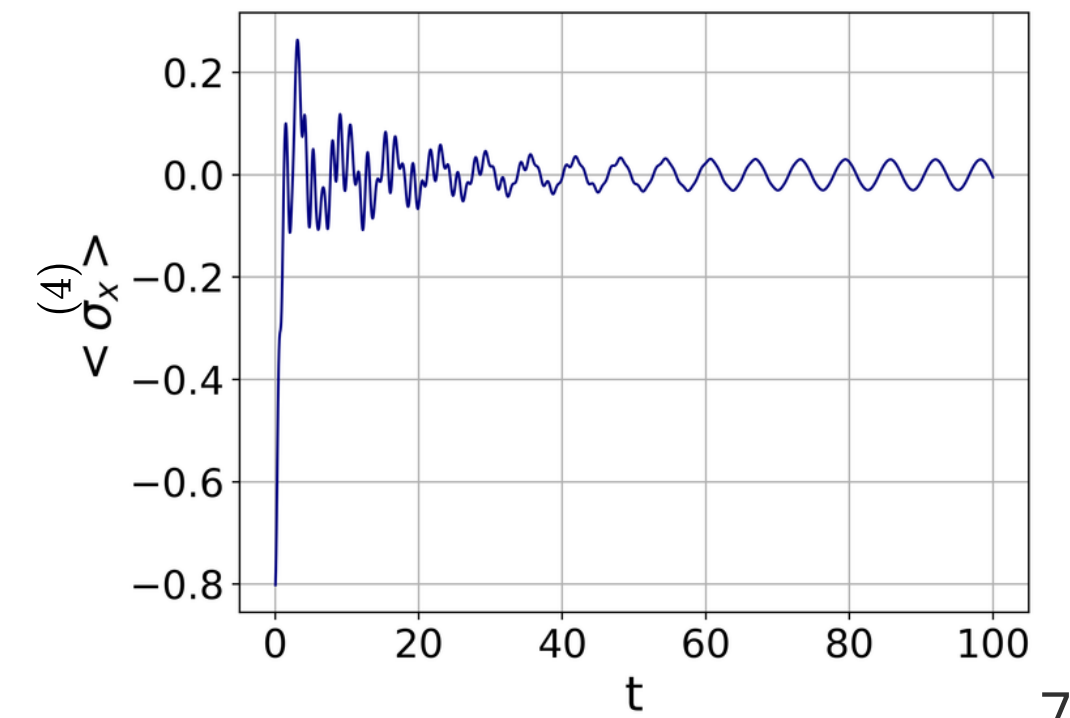
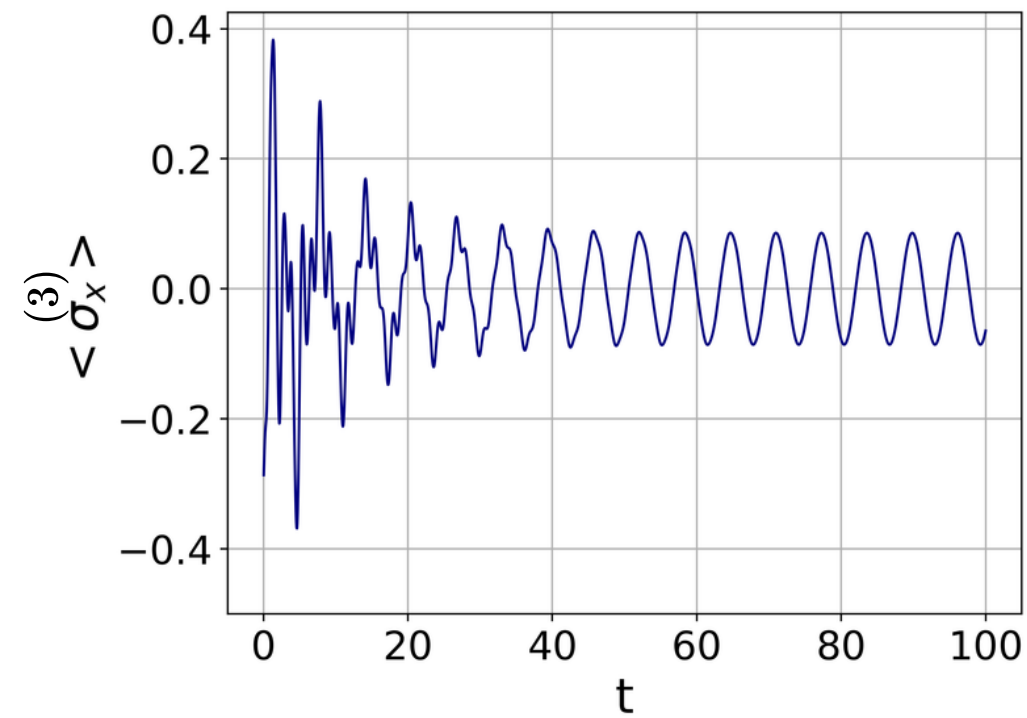
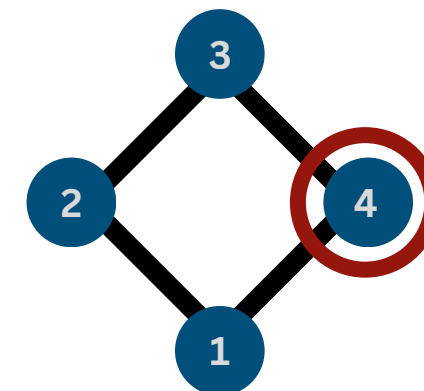
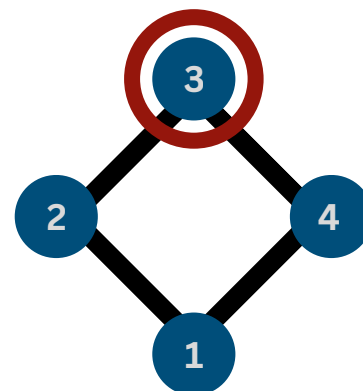
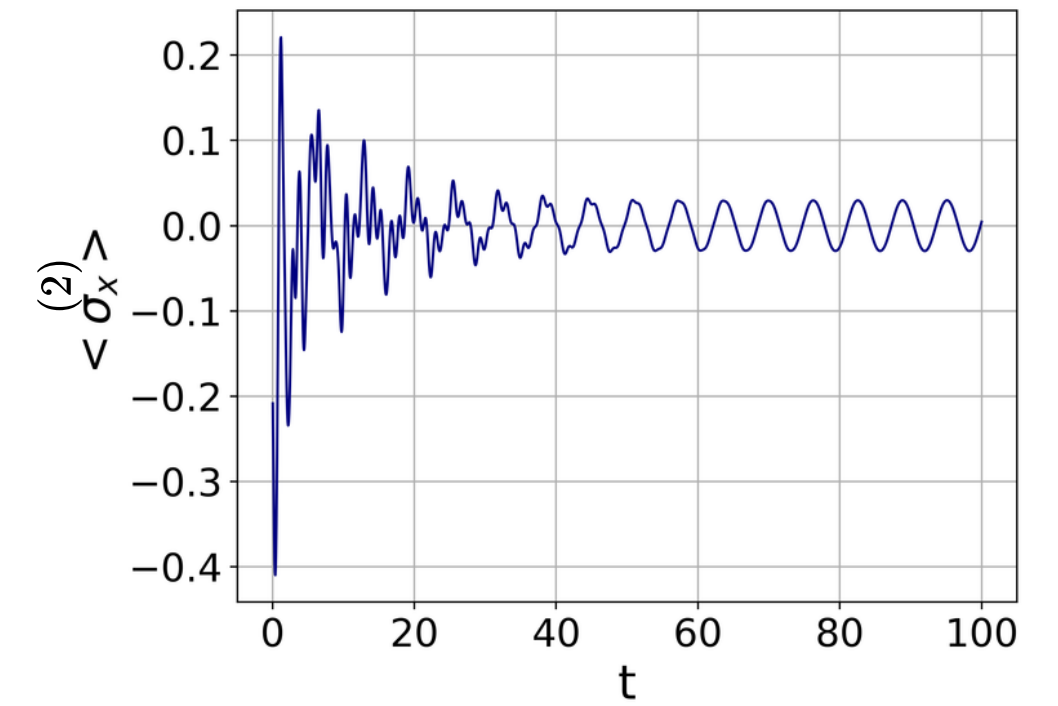
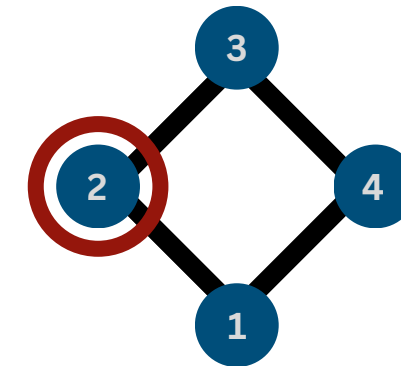
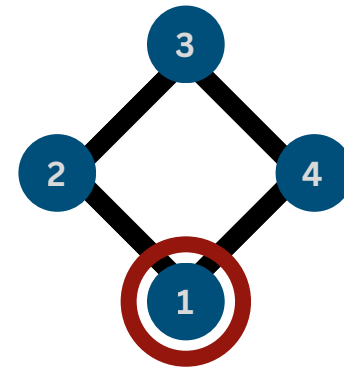
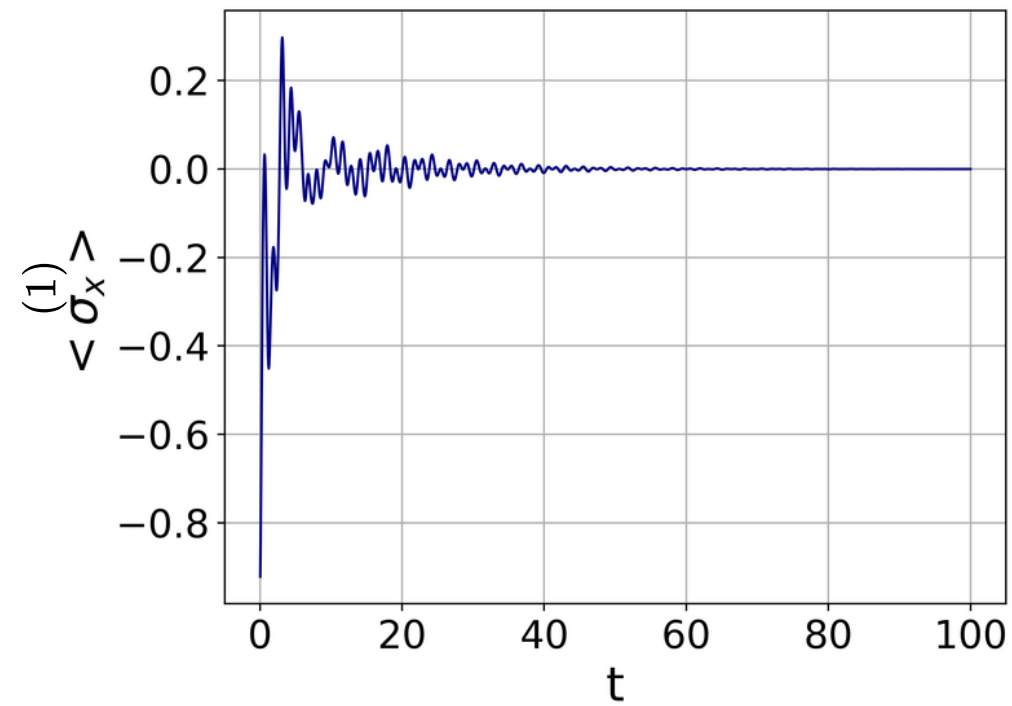
On-site energy
+ interaction between sites

On-site energy
(environment)

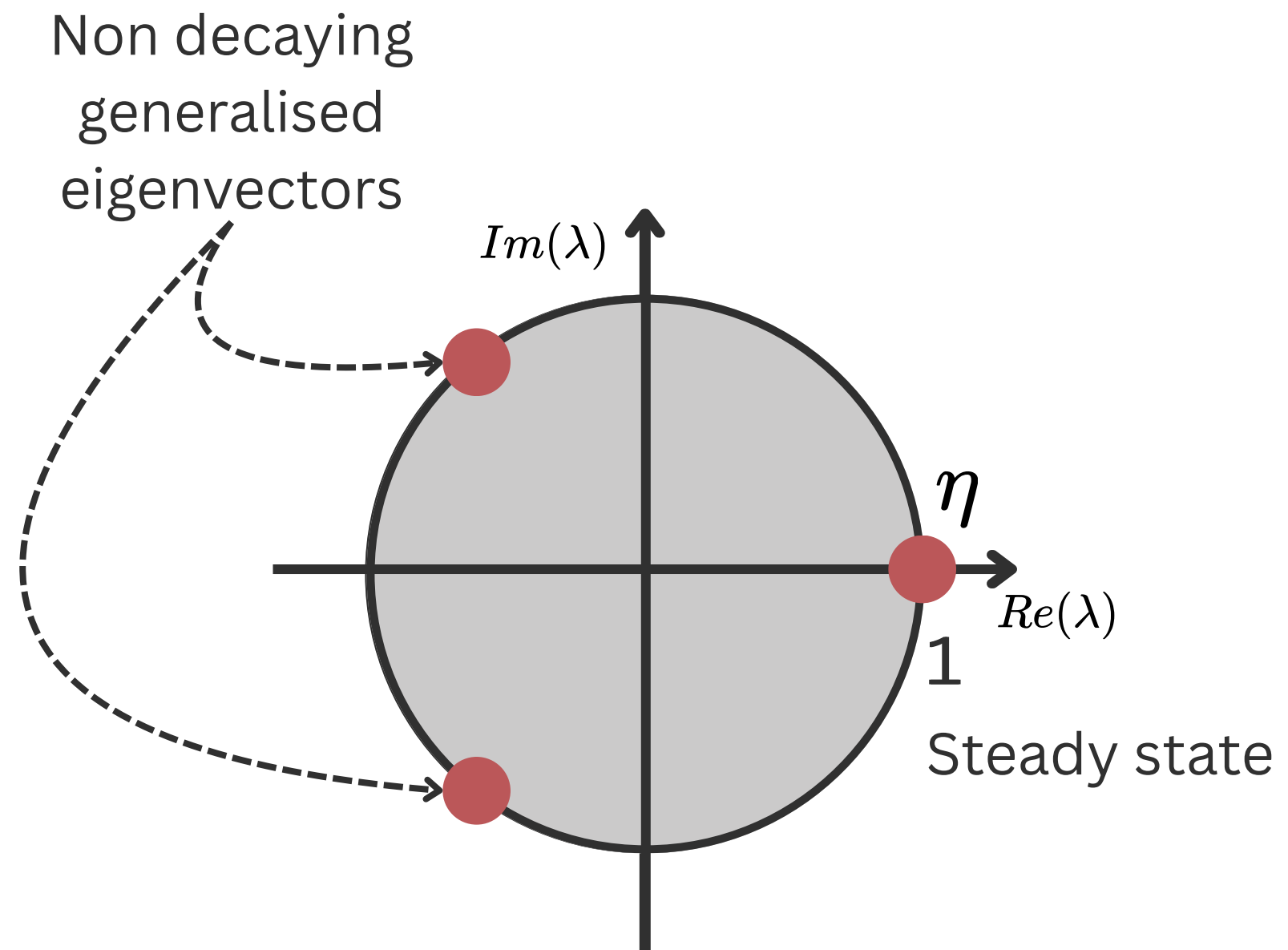
Interaction Hamiltonian

$$H_{SA}(t) = \sum_n g(\sigma_x^{(1)} \sigma_x^{(A)} + \sigma_y^{(1)} \sigma_y^{(A)}) \Theta(t - \theta_n) \Theta(\theta_n + \tau - t)$$

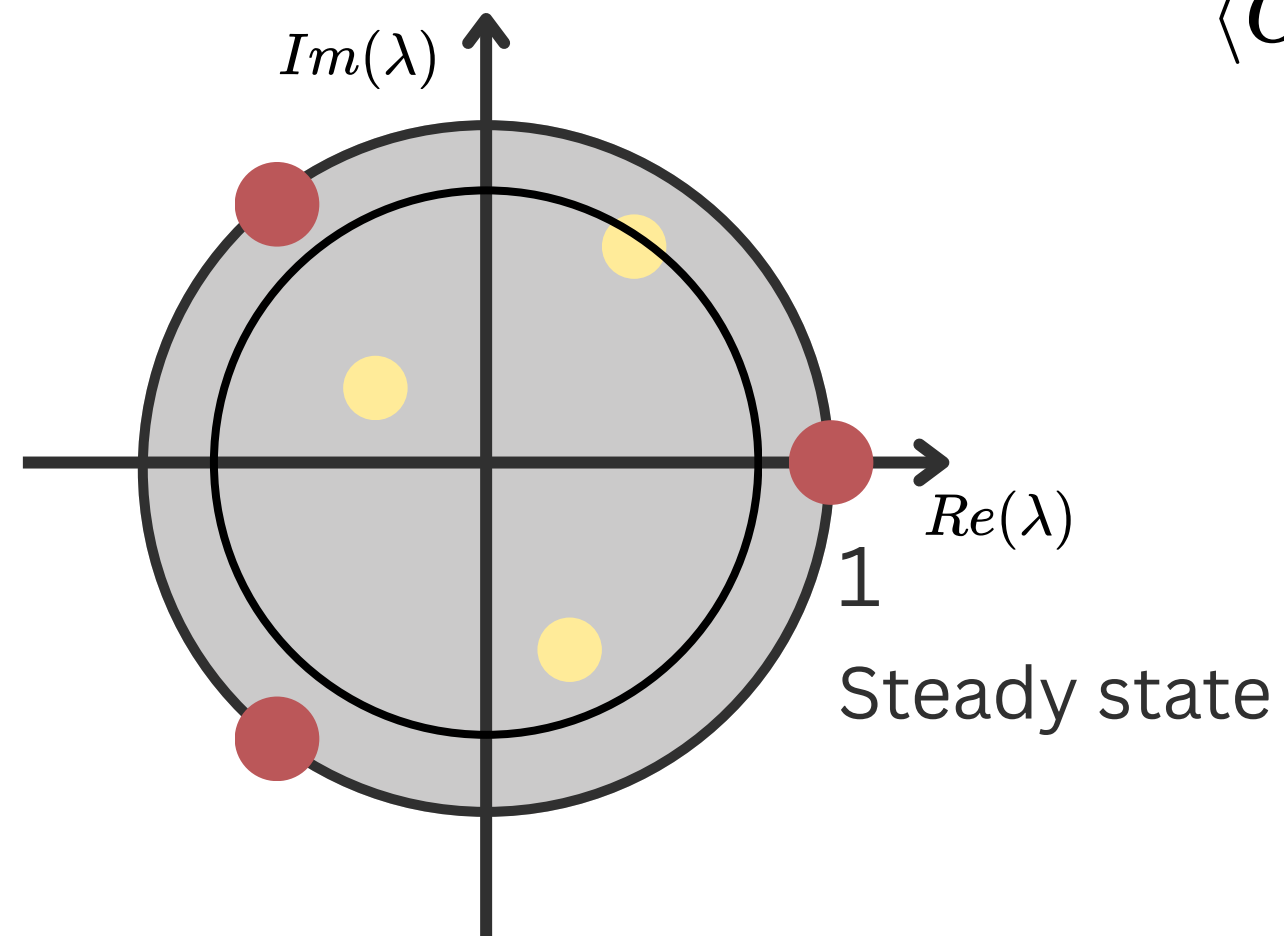
Dynamics of the N=4 ring



Dynamical symmetries



Dynamical symmetries

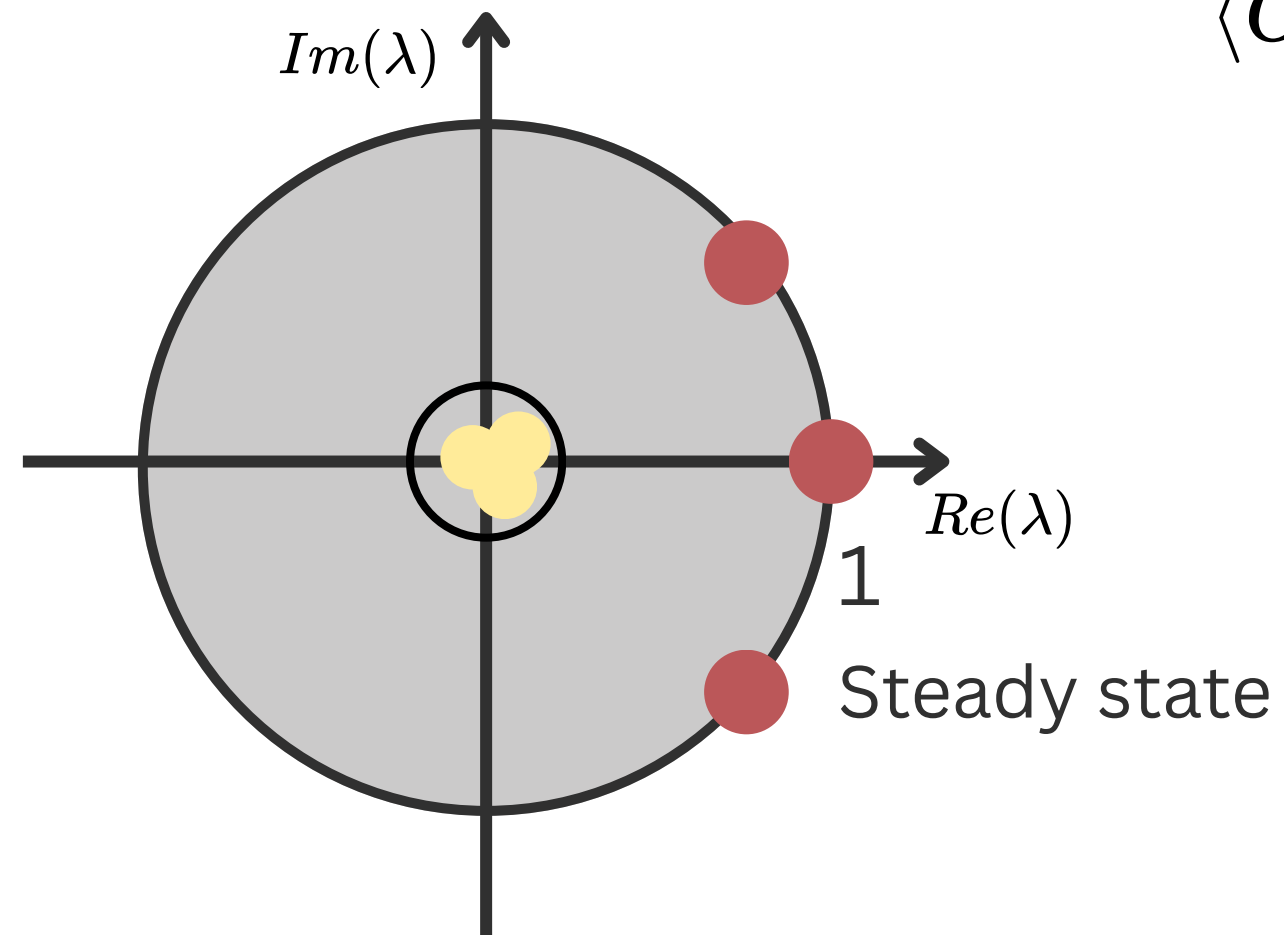


$$\langle O(nt) \rangle \xrightarrow{n \rightarrow \infty} r_0 Tr[\eta O] + \sum_m e^{i\omega_\alpha t} r_\alpha Tr[O \Xi_\alpha \eta] + h.c.$$

$$\rho_S(t) = \Lambda_t[\rho_S(0)]$$

Dynamical symmetries

Guarnieri, et al. Physical Review
A 106.2 (2022): 022209.



$$\langle O(nt) \rangle \xrightarrow{n \rightarrow \infty} r_0 Tr[\eta O] + \sum_m e^{i\omega_\alpha t} r_\alpha Tr[O \Xi_\alpha \eta] + h.c.$$

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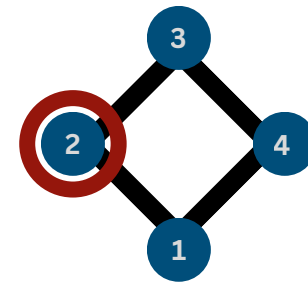
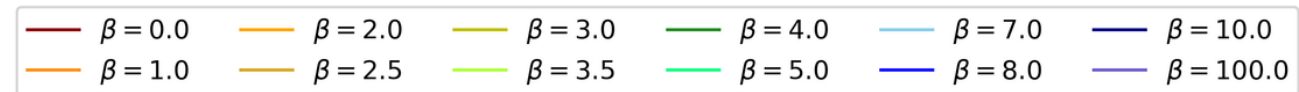
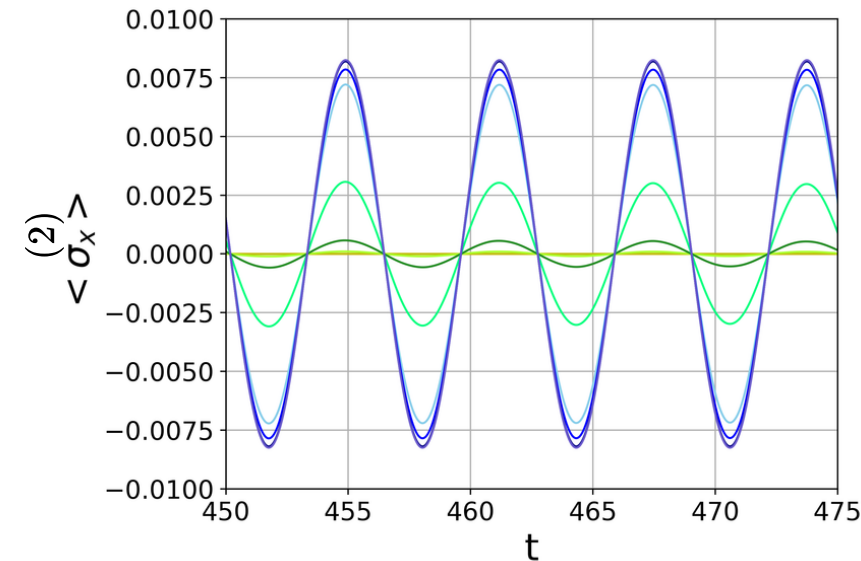
↓ n repetitions

$$\rho_S(nt) = [\Lambda_t]^n[\rho_S(0)]$$

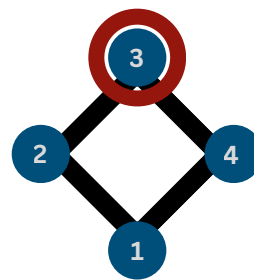
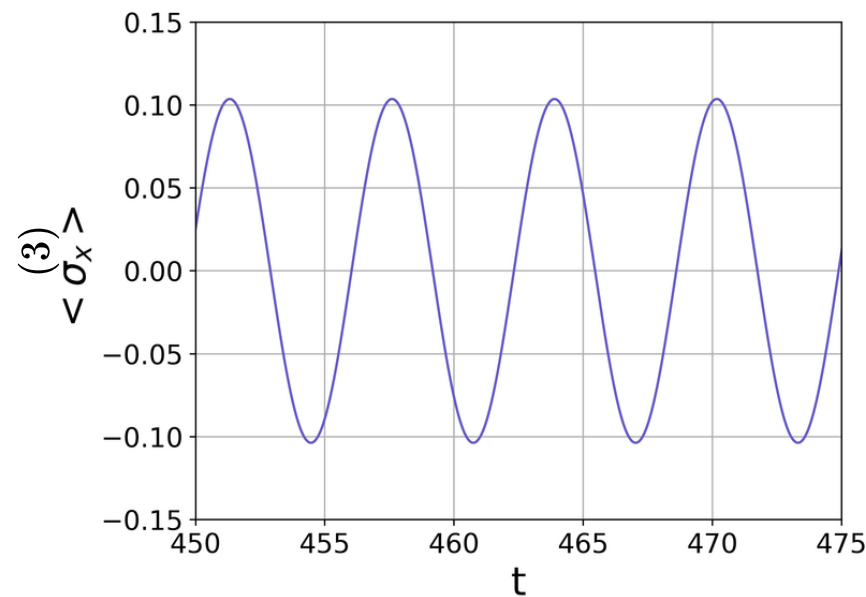
Only the generalised eigenvectors in the peripheral spectrum survive

Summary of preliminary results

1) Robustness to temperature



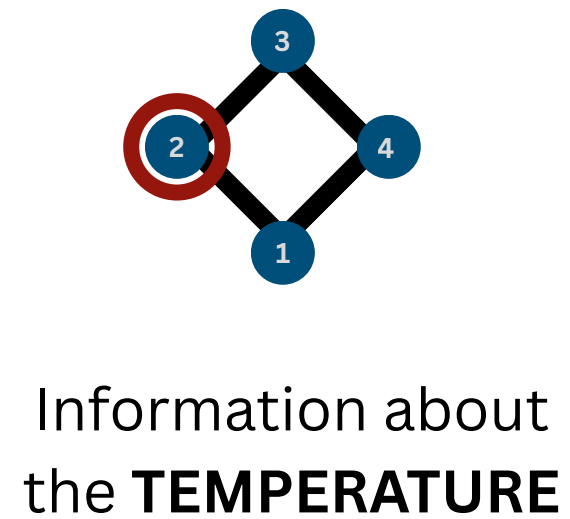
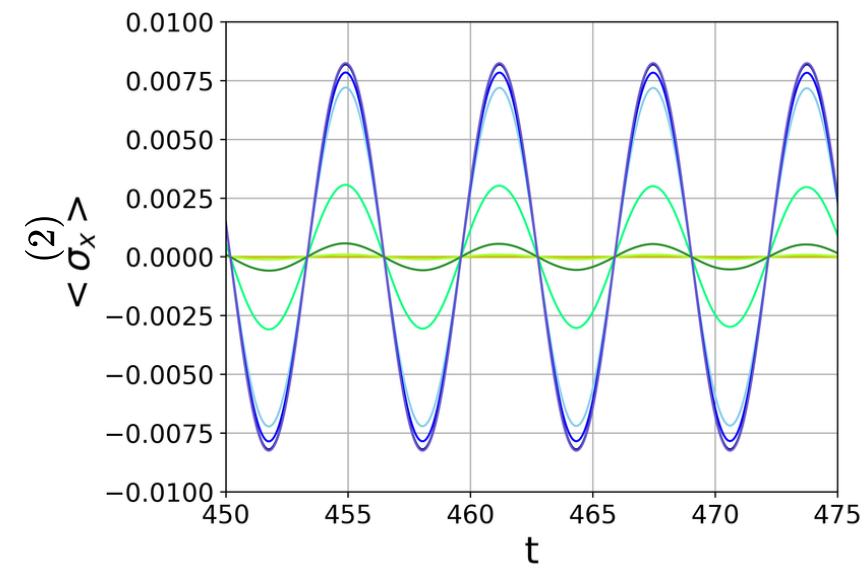
Information about
the **TEMPERATURE**



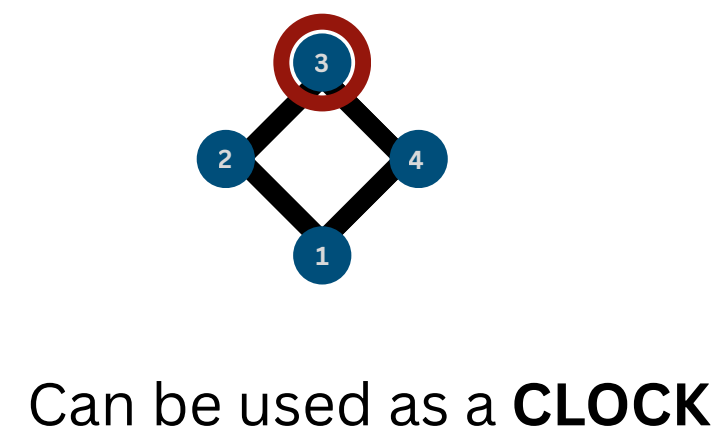
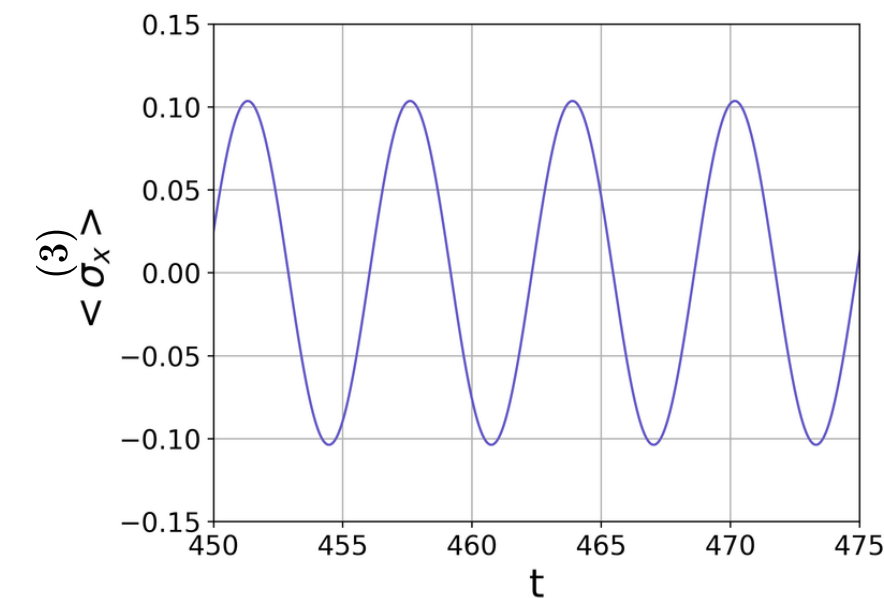
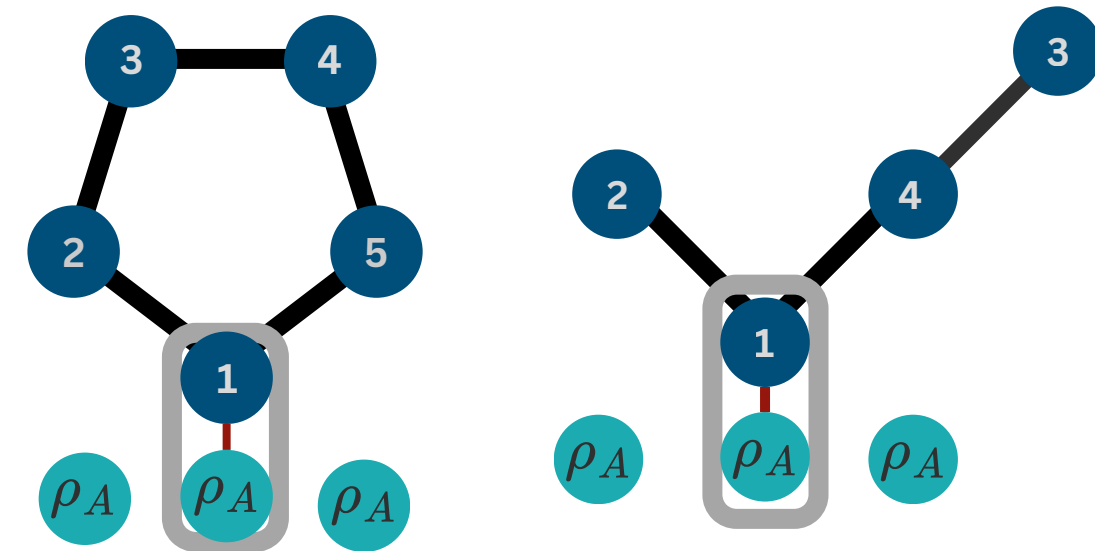
Can be used as a **CLOCK**

Summary of preliminary results

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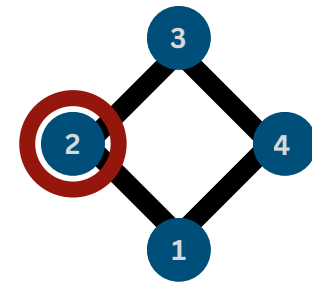
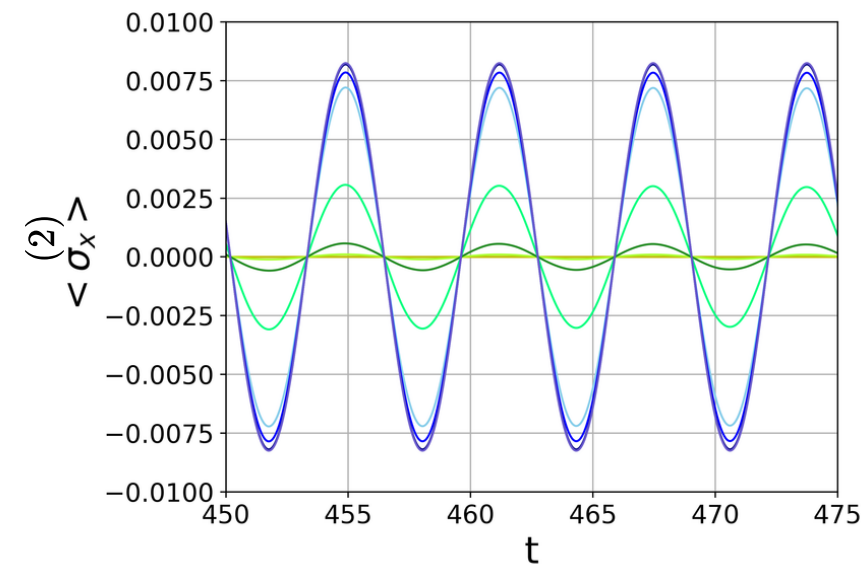


2) Different dynamical configurations

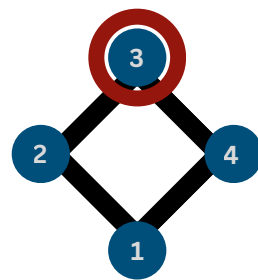
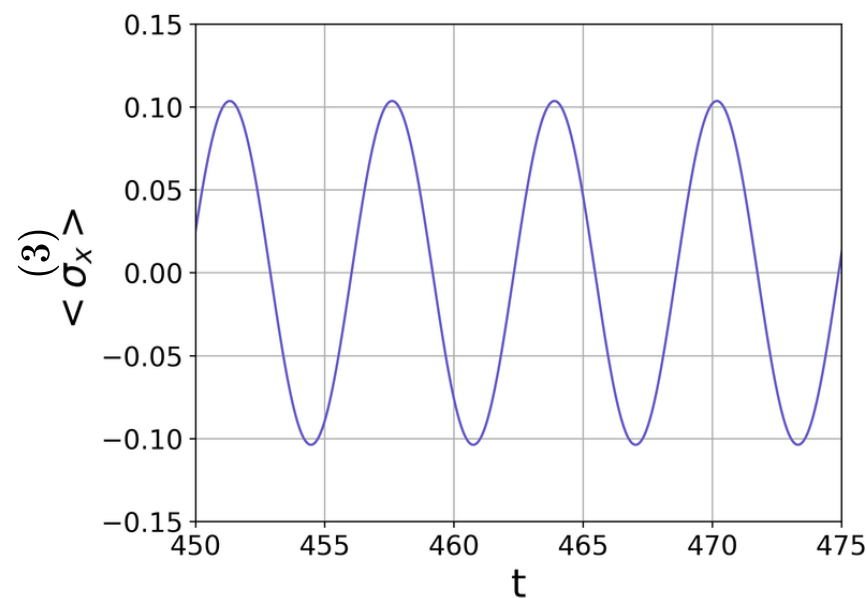


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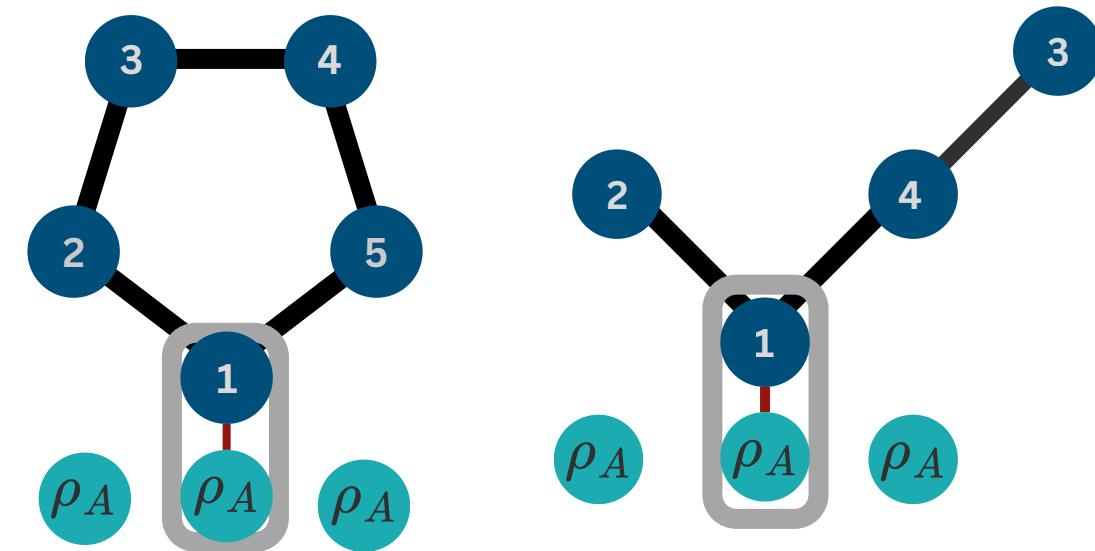


Information about the **TEMPERATURE**

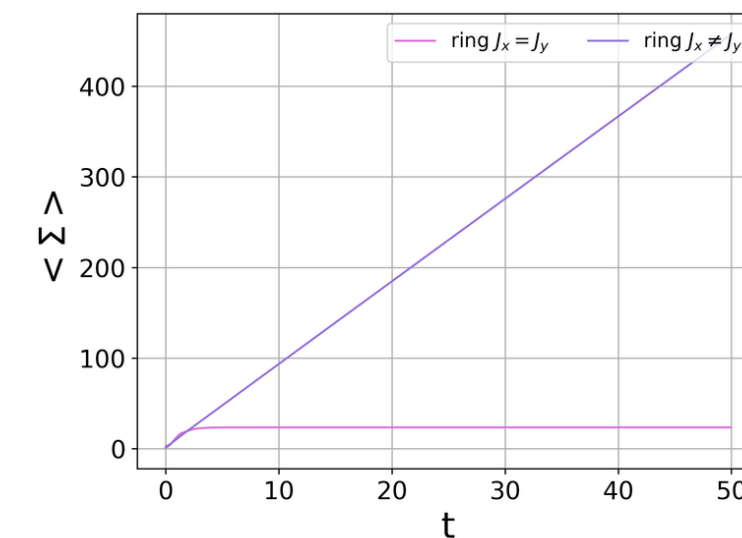


Can be used as a **CLOCK**

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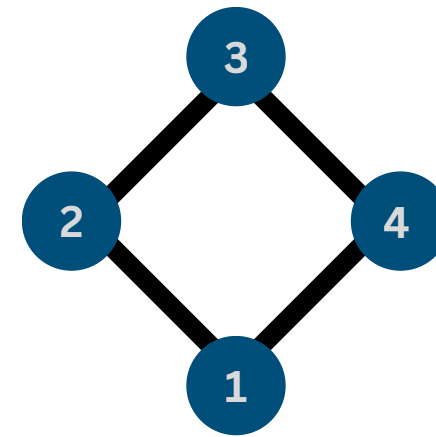


3) Characterisation of the thermodynamic cost

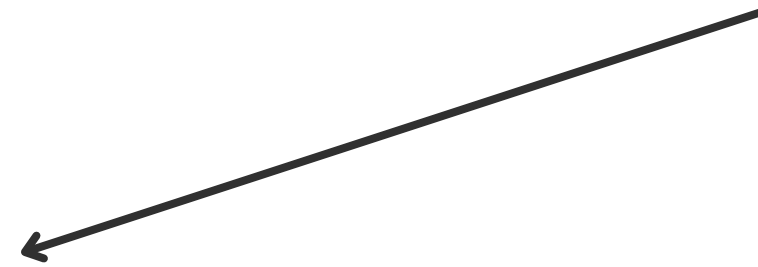
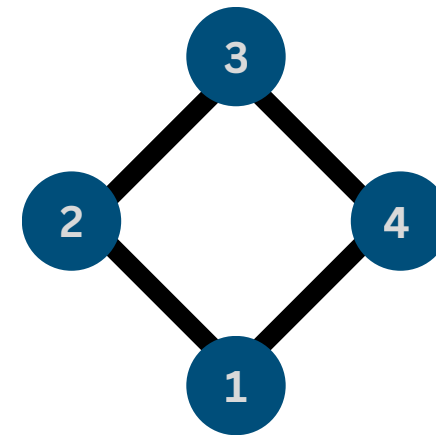


$$\Sigma = \Delta S_S + \Delta S_E \geq 0$$

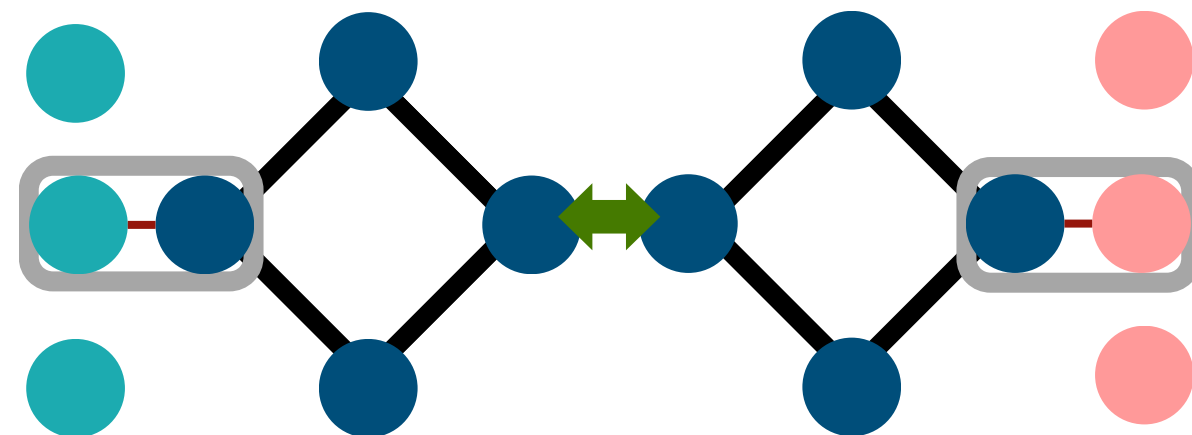
New questions



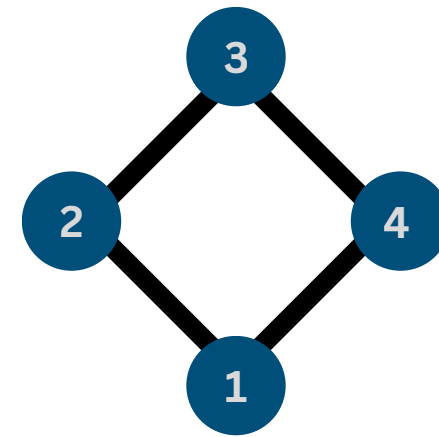
New questions



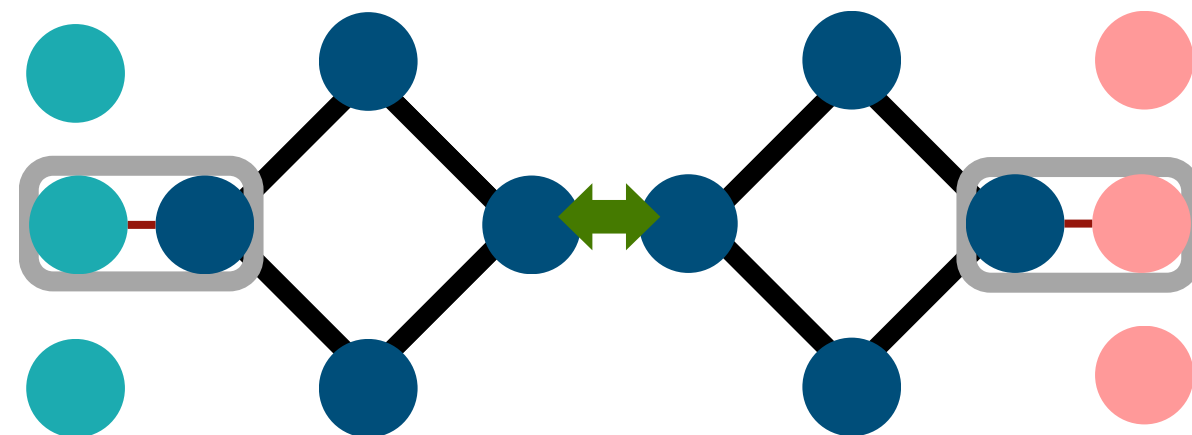
What about **TRANSPORT PROPERTIES** between two thermal baths?



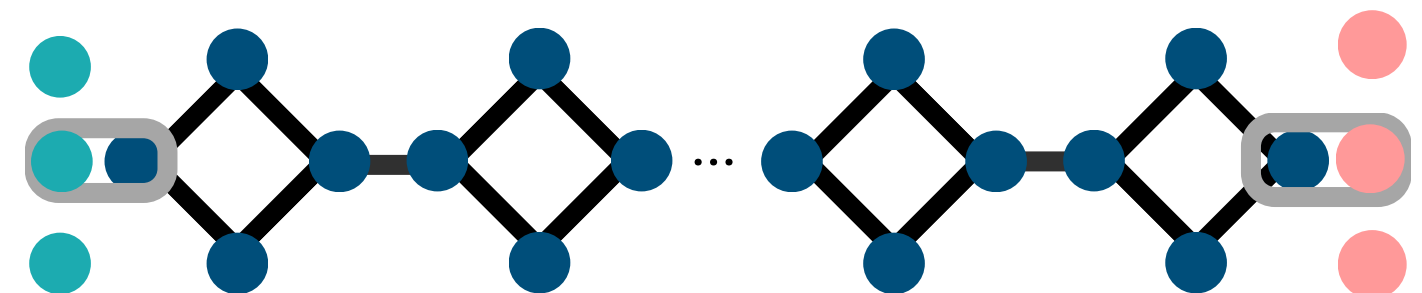
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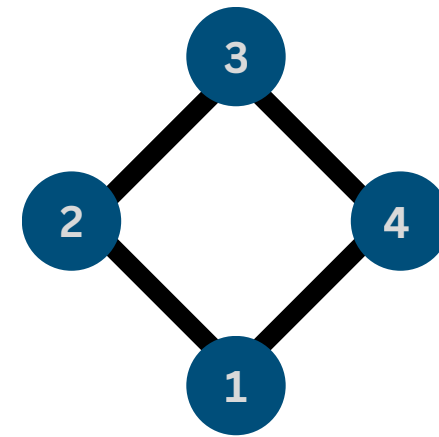
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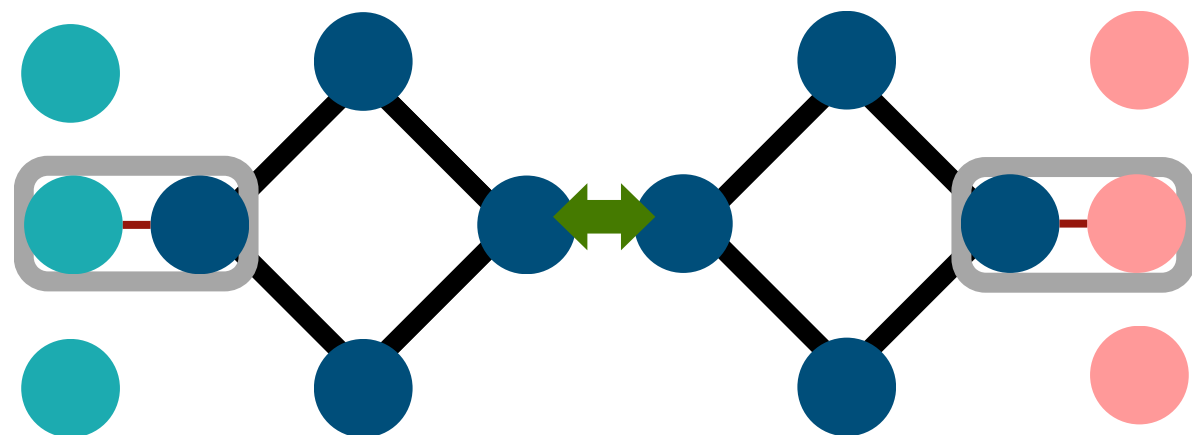
What happens when we **scale up** the system?



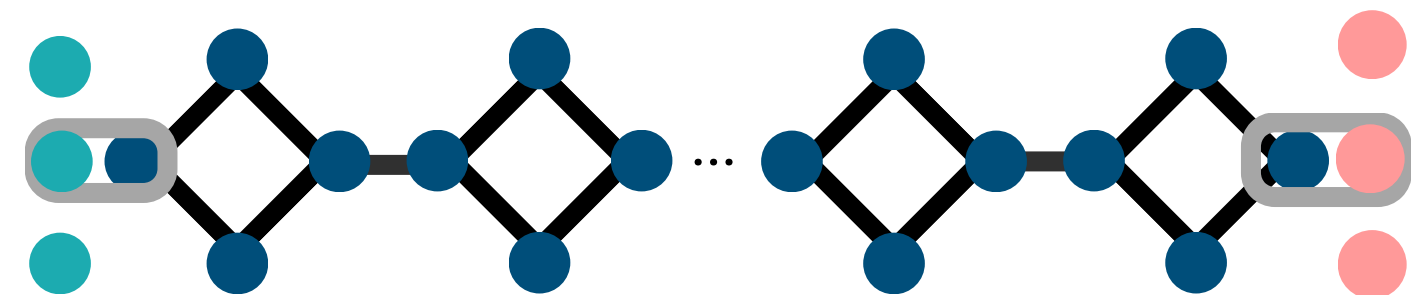
New questions



What about **TRANSPORT PROPERTIES** between two thermal baths?



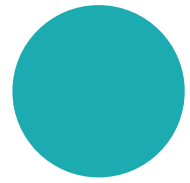
What happens when we **scale up** the system?



PROBLEM: Computational complexity is too high

Tensor networks - graphical representation

Basic elements

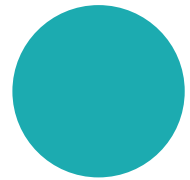


Number

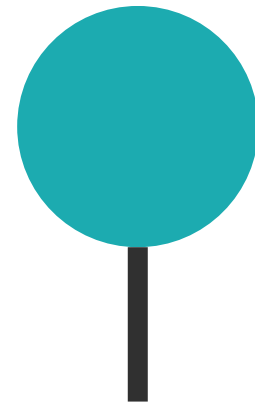
a

Tensor networks - graphical representation

Basic elements



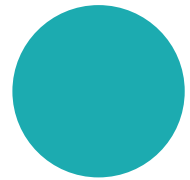
Number
 a



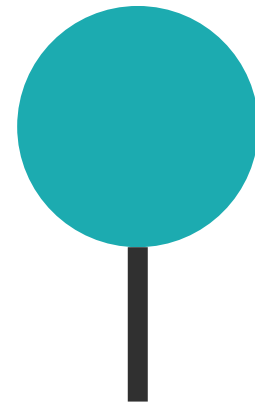
Vector
 v_i

Tensor networks - graphical representation

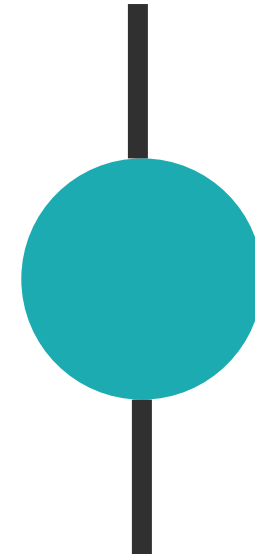
Basic elements



Number
 a



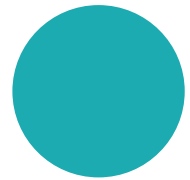
Vector
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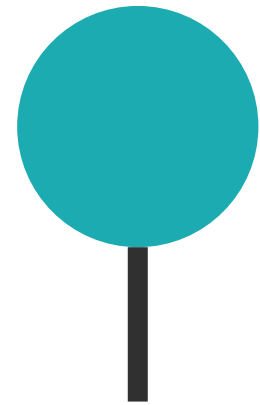
Matrix
 $M_{i,j}$

Tensor networks - graphical representation

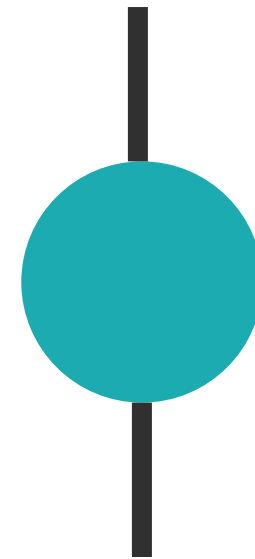
Basic elements



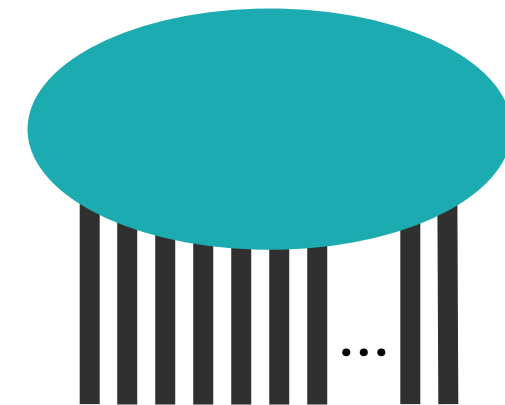
Number
 a



Vector
 v_i



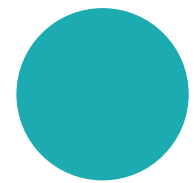
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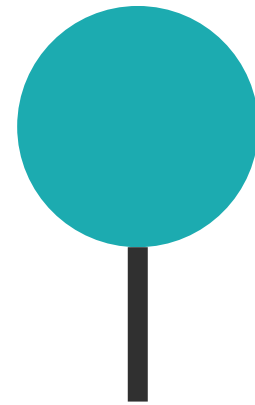
Tensor
 $T_{l,m,n...}^{i,j,k,...}$

Tensor networks - graphical representation

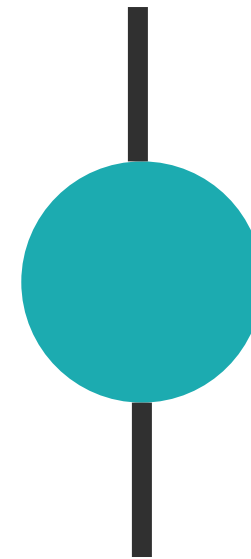
Basic elements



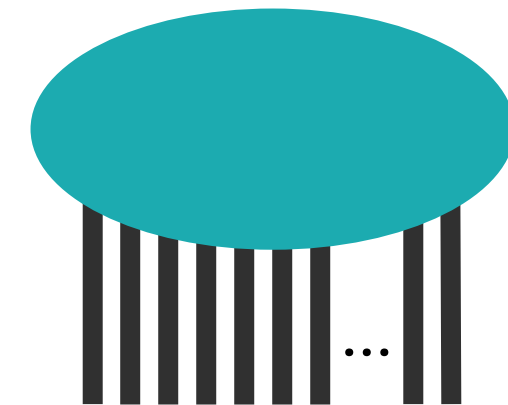
Number
 a



Vector
 v_i



Matrix
 $M_{i,j}$



Tensor
 $T_{l,m,n...}^{i,j,k,...}$

Basic operation: contraction

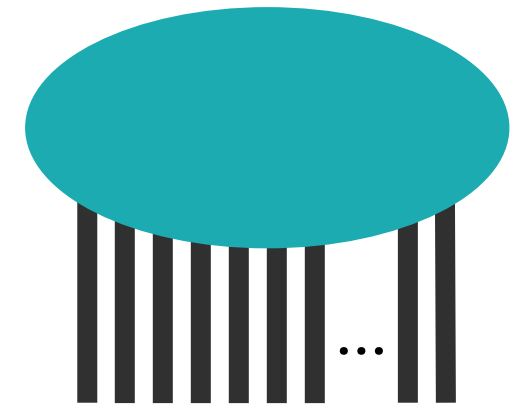
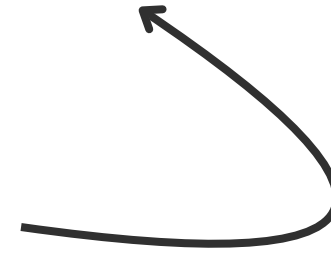


$$w_j = \sum_i M_{i,j} v_i$$

Tensor networks

Generic quantum state: $|\psi\rangle = \sum_{\alpha_1, \alpha_2, \dots, \alpha_N} \psi_{\alpha_1, \alpha_2, \dots, \alpha_N} |\alpha_1, \alpha_2, \dots, \alpha_N\rangle$

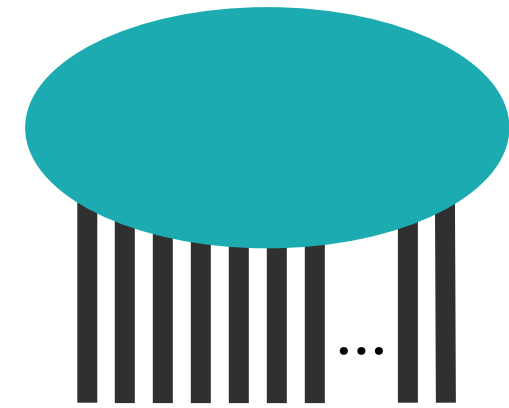
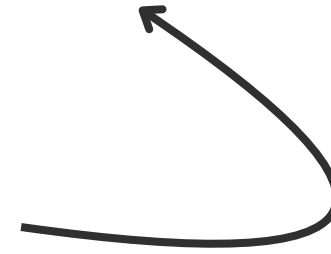
Exponential scaling



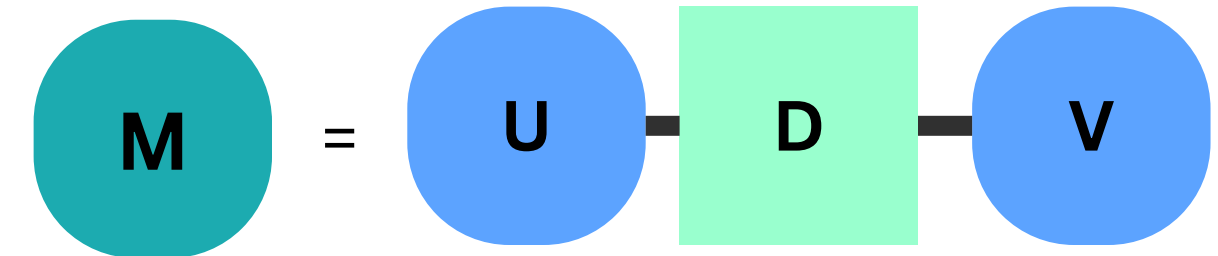
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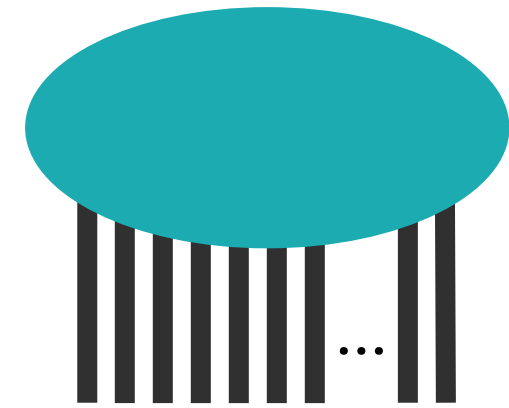
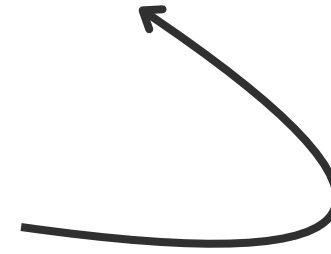
It is possible to use SVD to add an internal leg and obtain a MPS



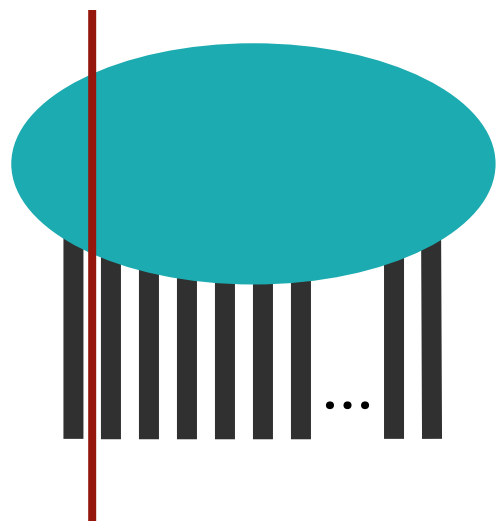
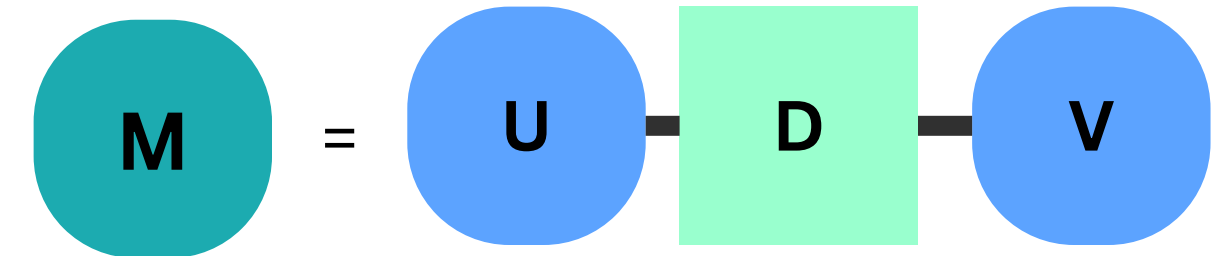
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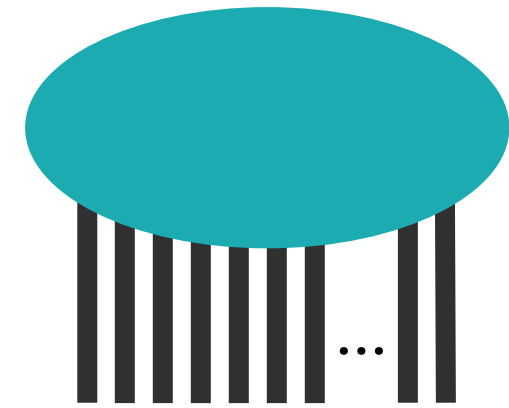
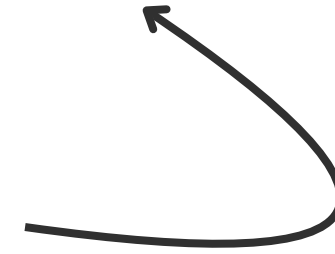
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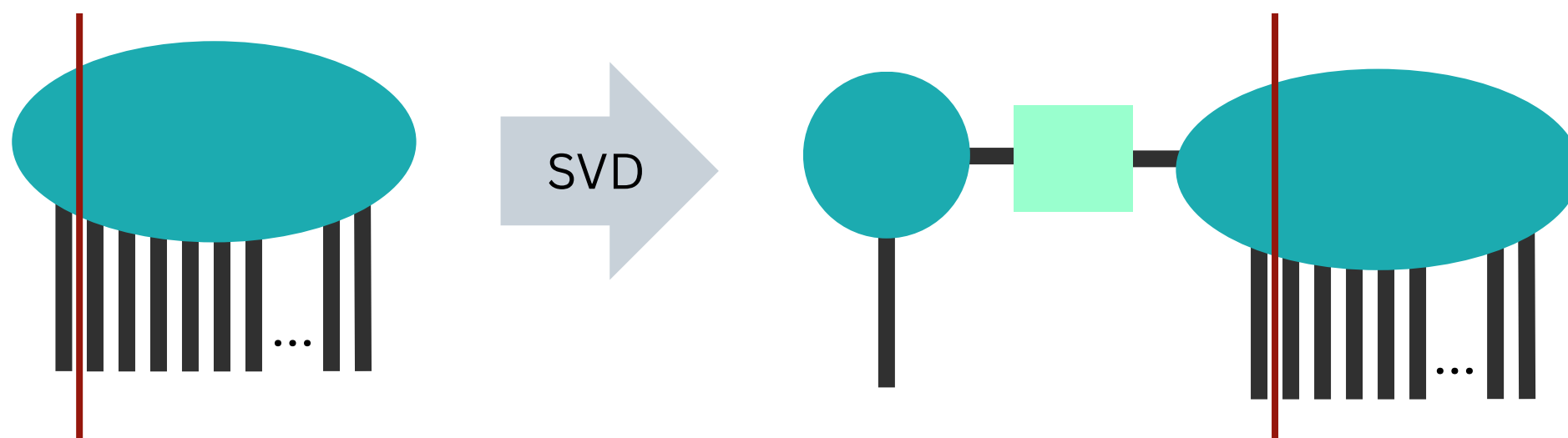
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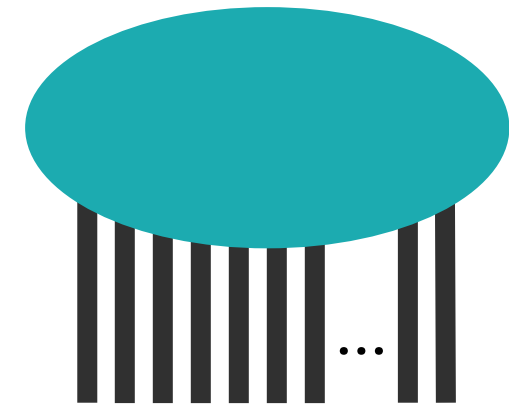
$$\mathbf{M} = \mathbf{U} \mathbf{D} \mathbf{V}$$



Tensor networks

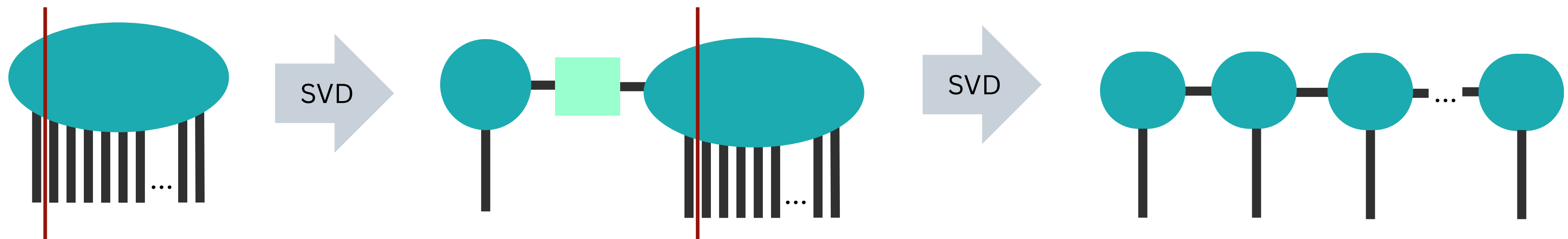
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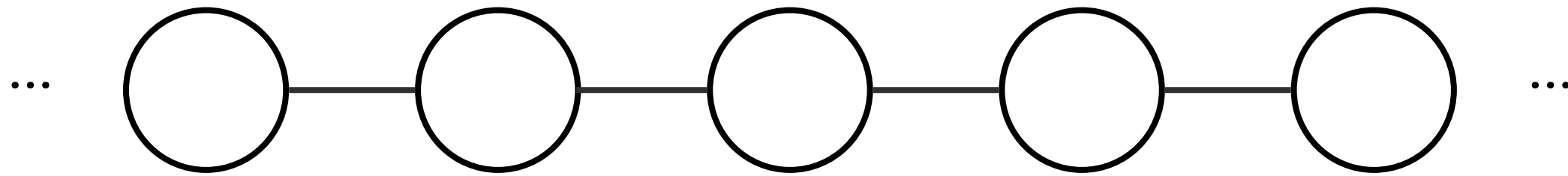
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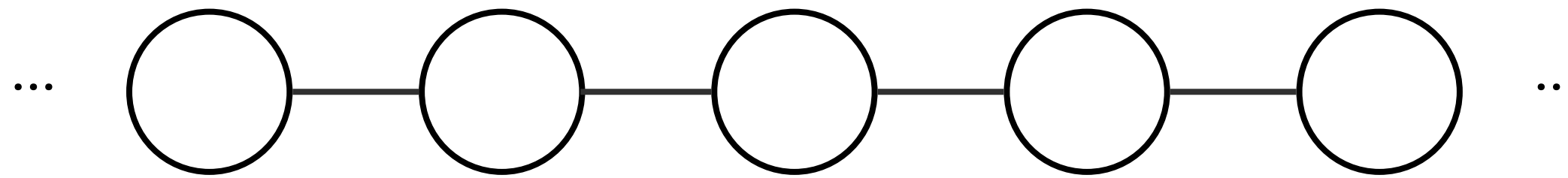
Tensor networks

The gain comes from the physics itself!



Tensor networks

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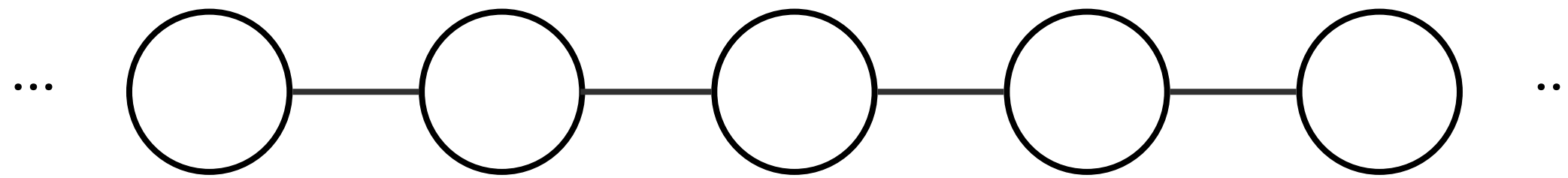
Typical Hamiltonian terms:

$$\propto h_n h_{n+1}$$

Schollwöck Annals of physics
326.1 (2011): 96-192.

Tensor networks

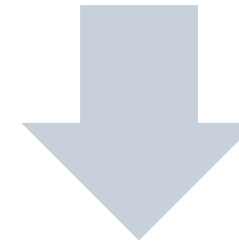
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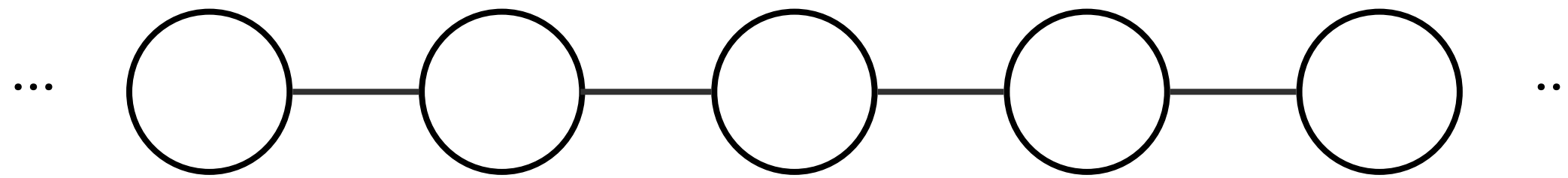
This automatically limits entanglement growth across sites in the lattice → it is possible to identify



TOTAL Hilbert space

Tensor networks

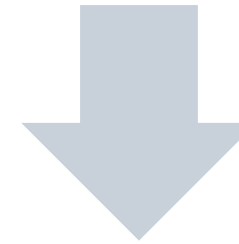
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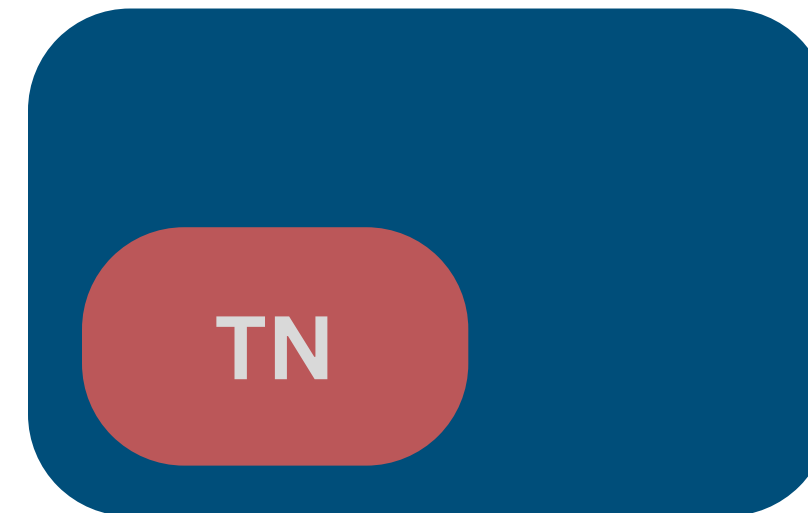
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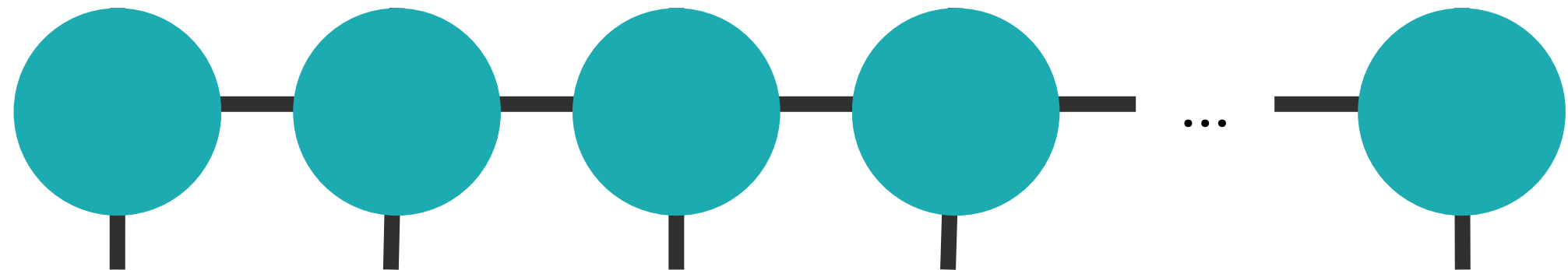
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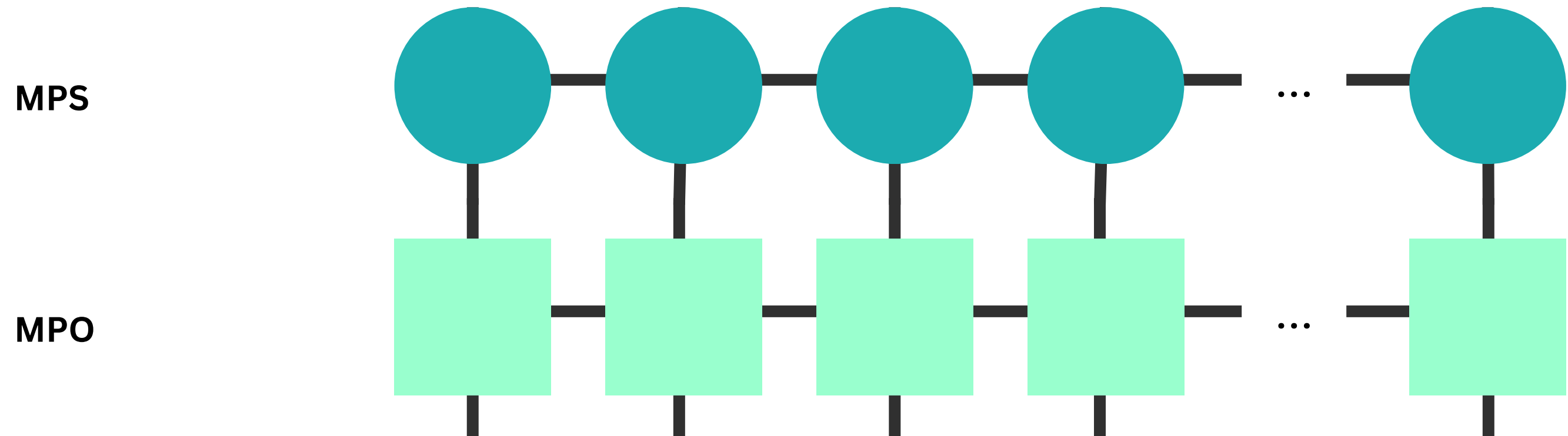
RELEVANT PORTION of the
Hilbert space

Tensor networks

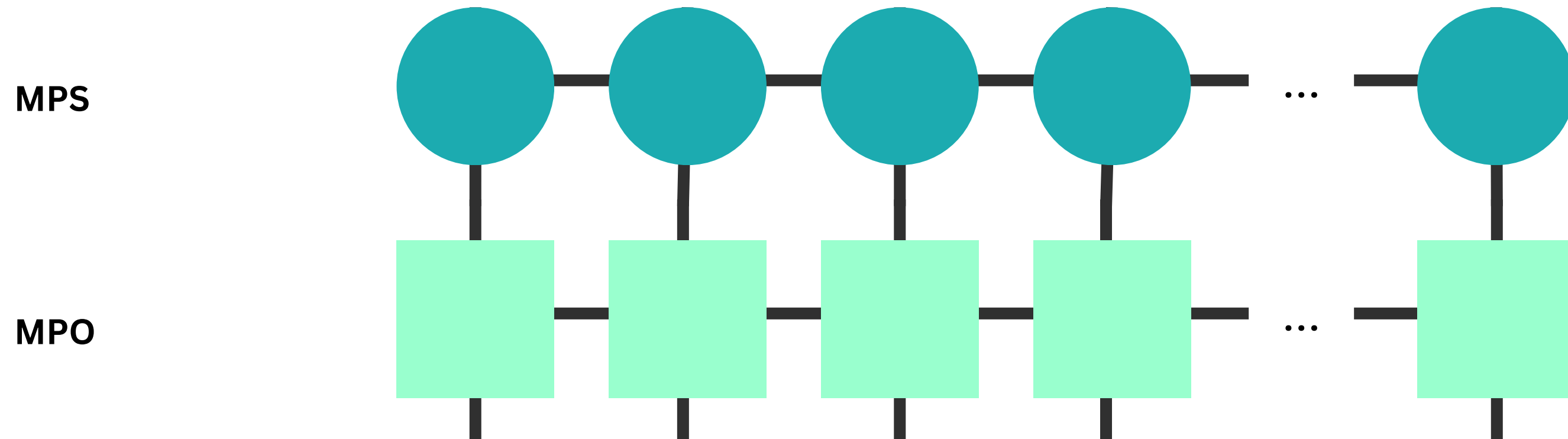
MPS



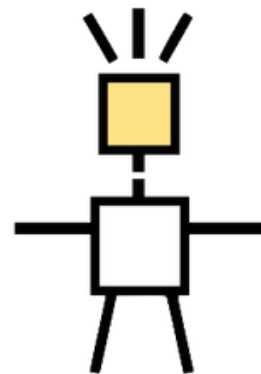
Tensor networks



Tensor networks



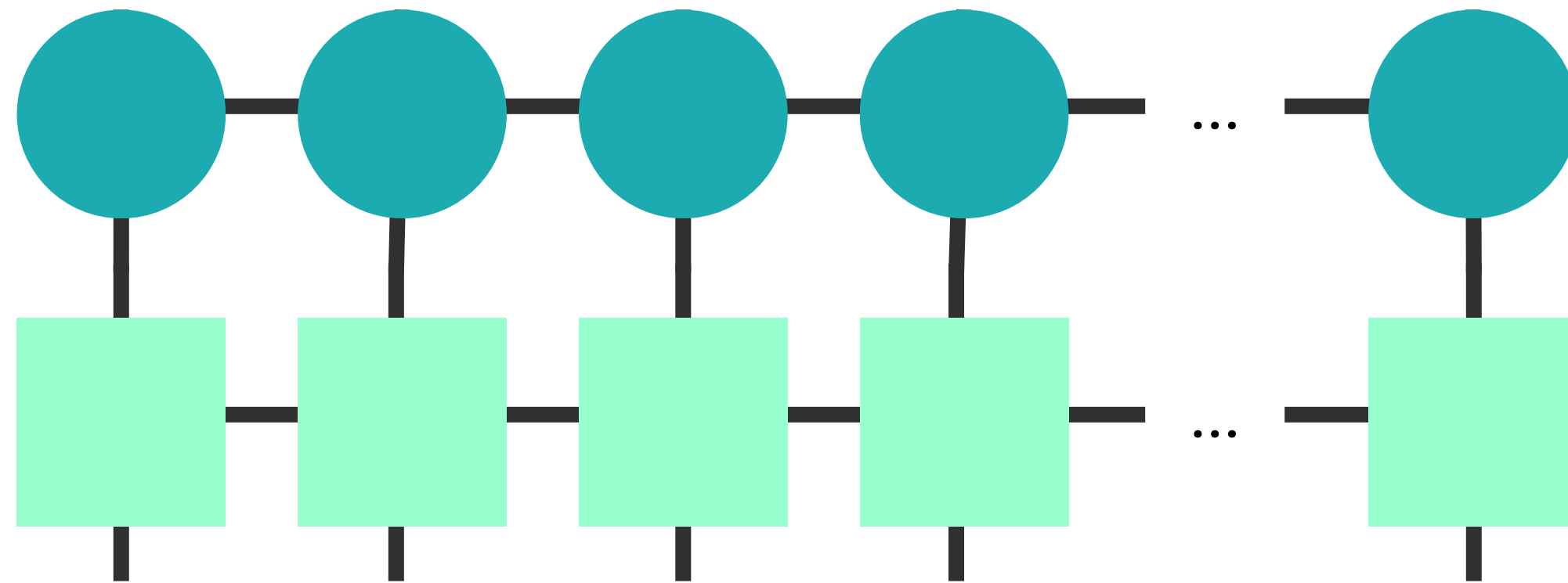
Adding some more tools
(thankfully) provided by the
ITensor library



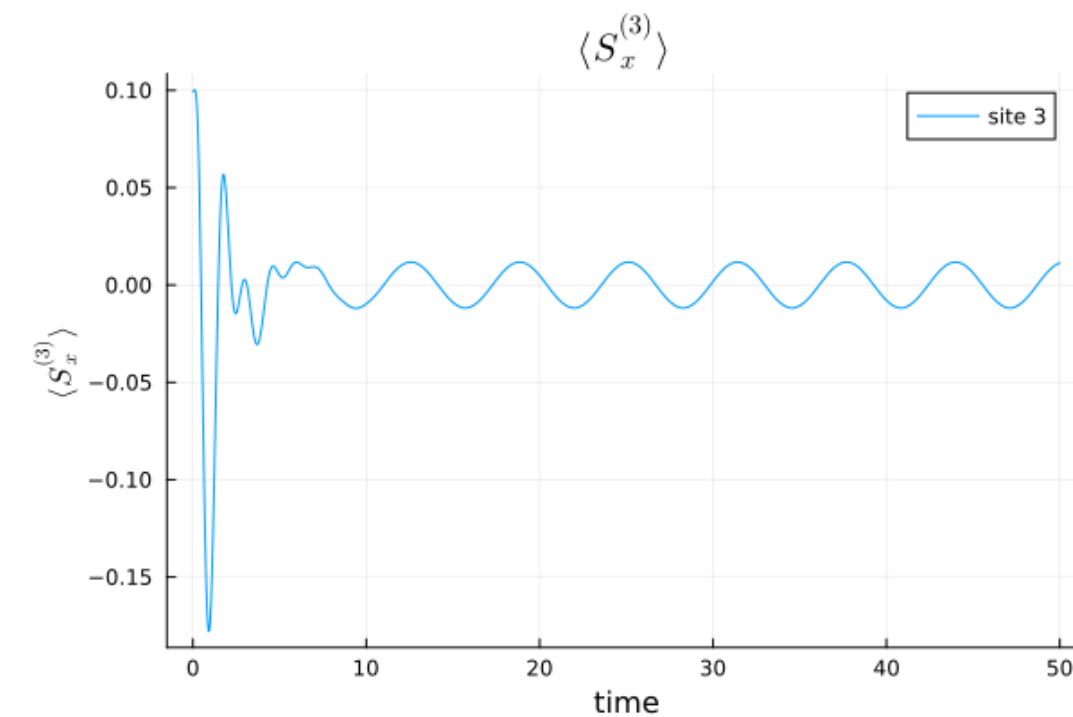
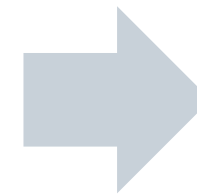
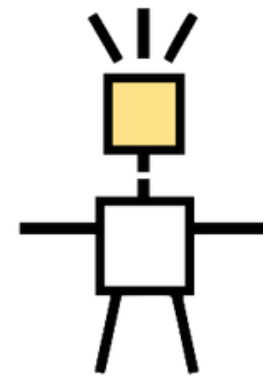
Tensor networks

MPS

MPO

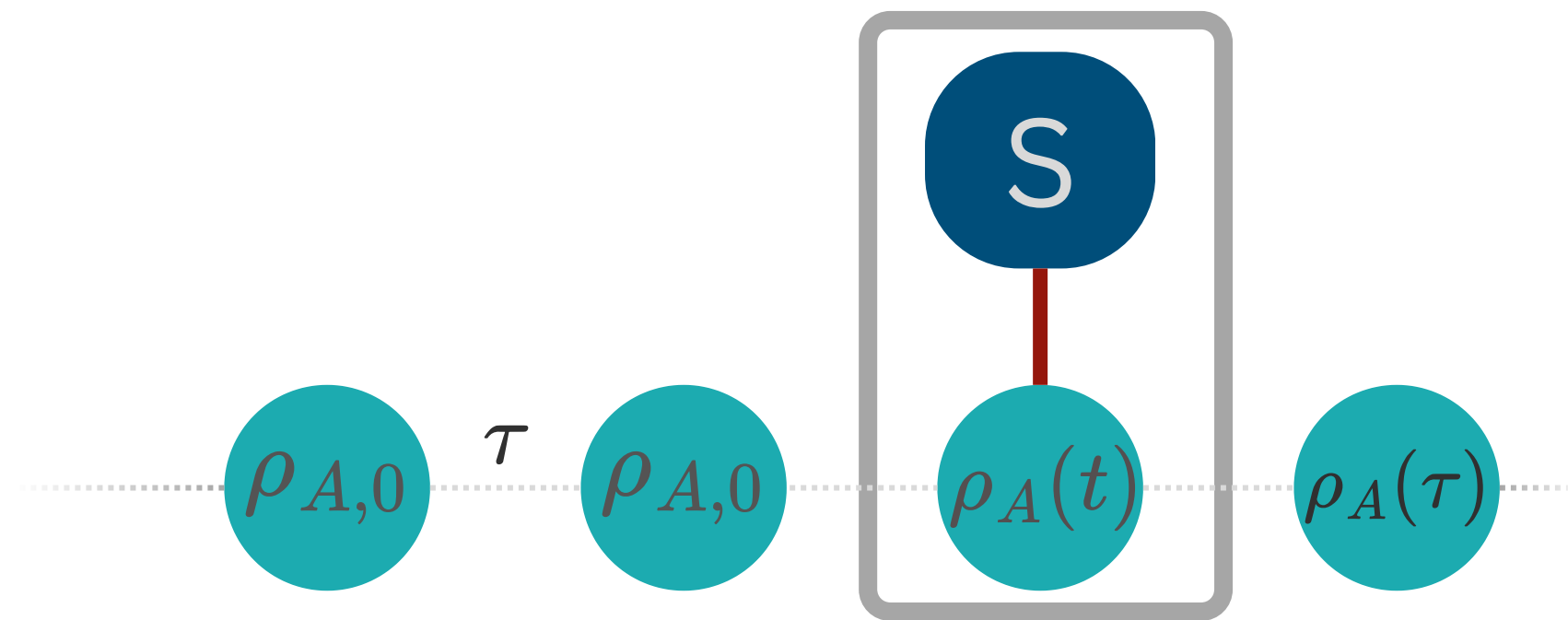


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Matthew Fishman, et al. SciPost Phys. Codebases 4 (2022)

Problem: matter exchange



Collision models:

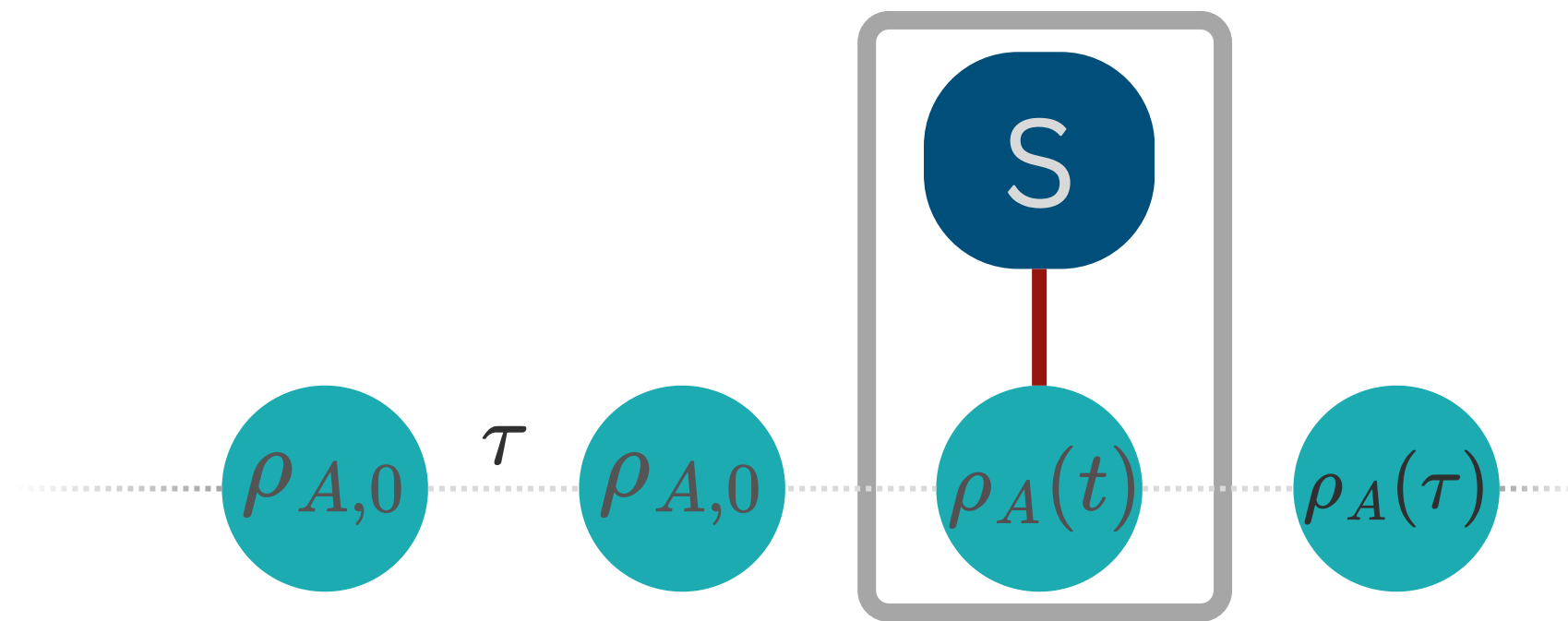


good description of the dynamics



successful at capturing energy exchanges

Problem: matter exchange



Collision models:



good description of the dynamics

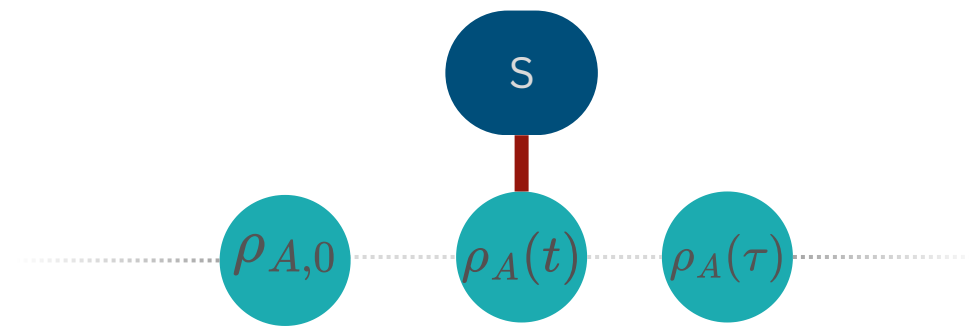


successful at capturing energy exchanges

Problem: can systems that exchange particles be characterised?

Clustering collision models

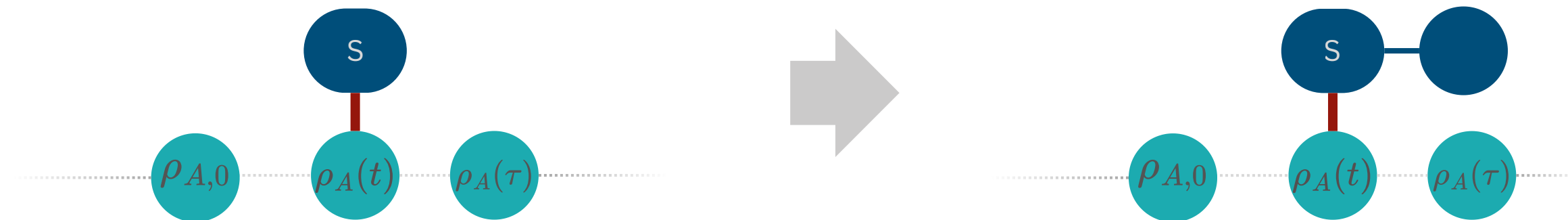
Attachment $\rho_S(\tau) = \mathcal{A}_p \circ \xi[\rho_S(0)]$



Clustering collision models

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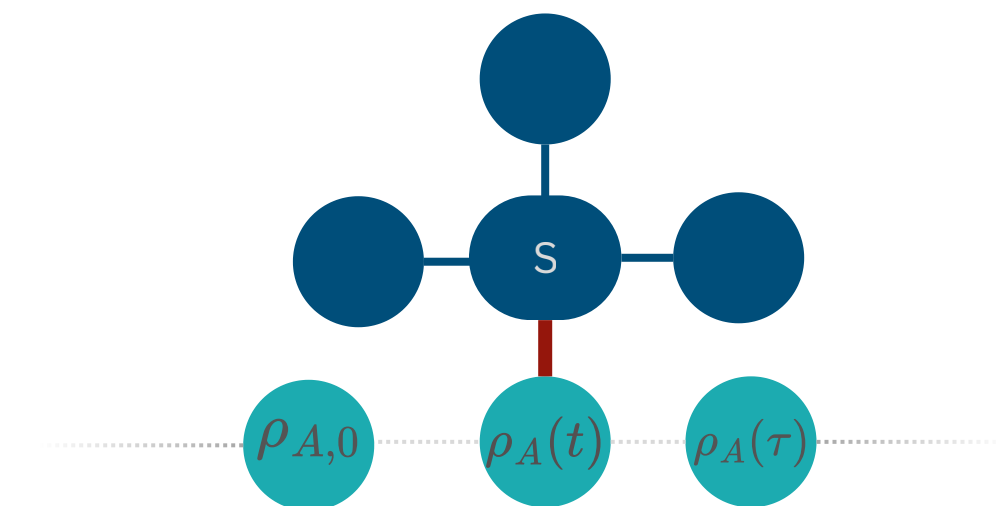
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Clustering collision models

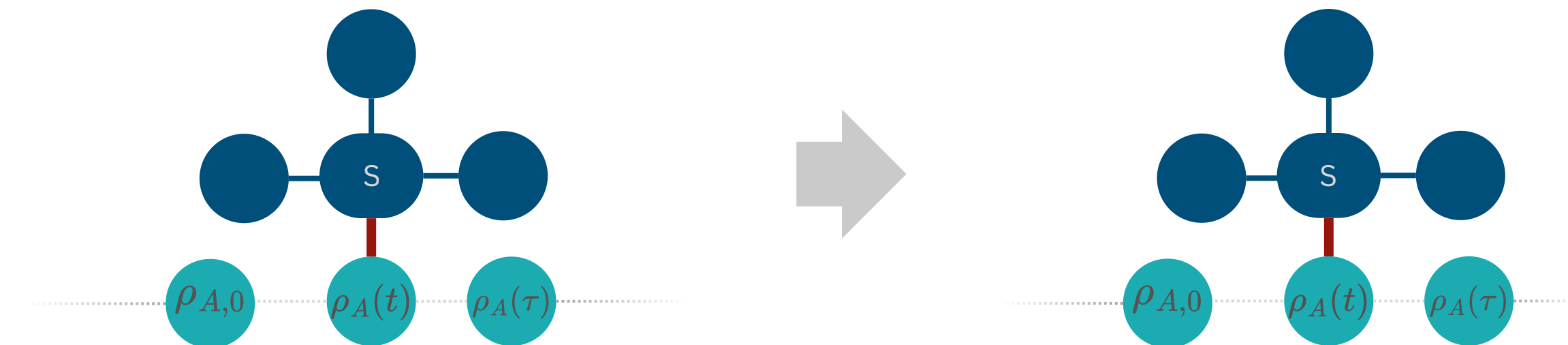
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Clustering collision models

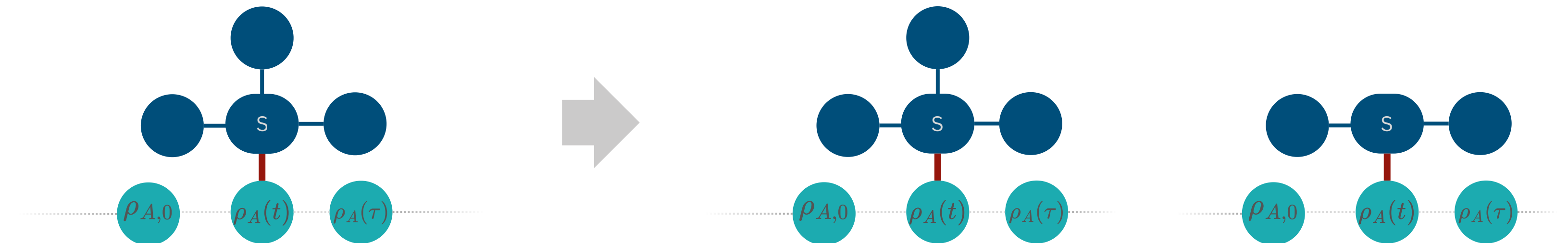
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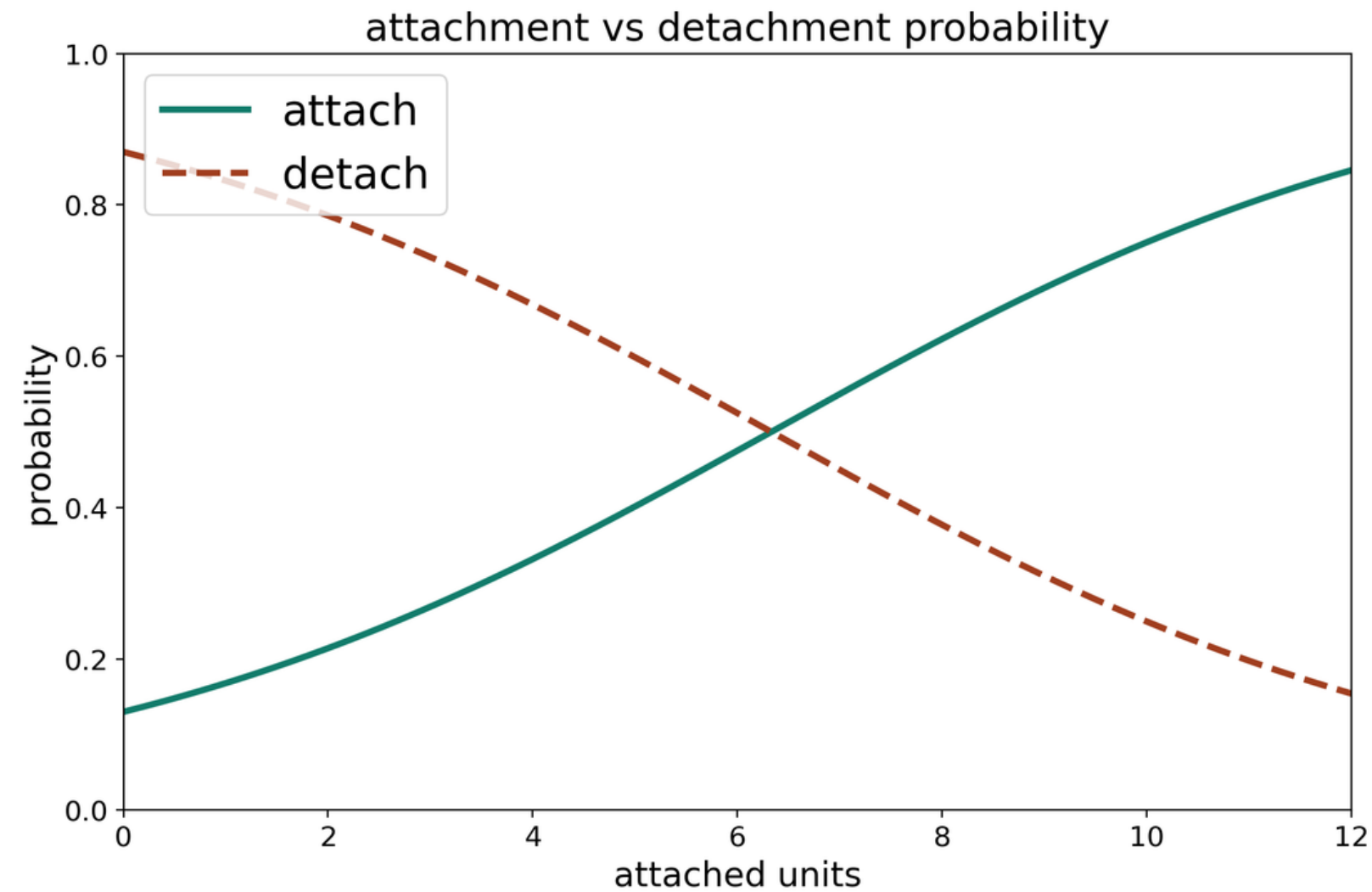
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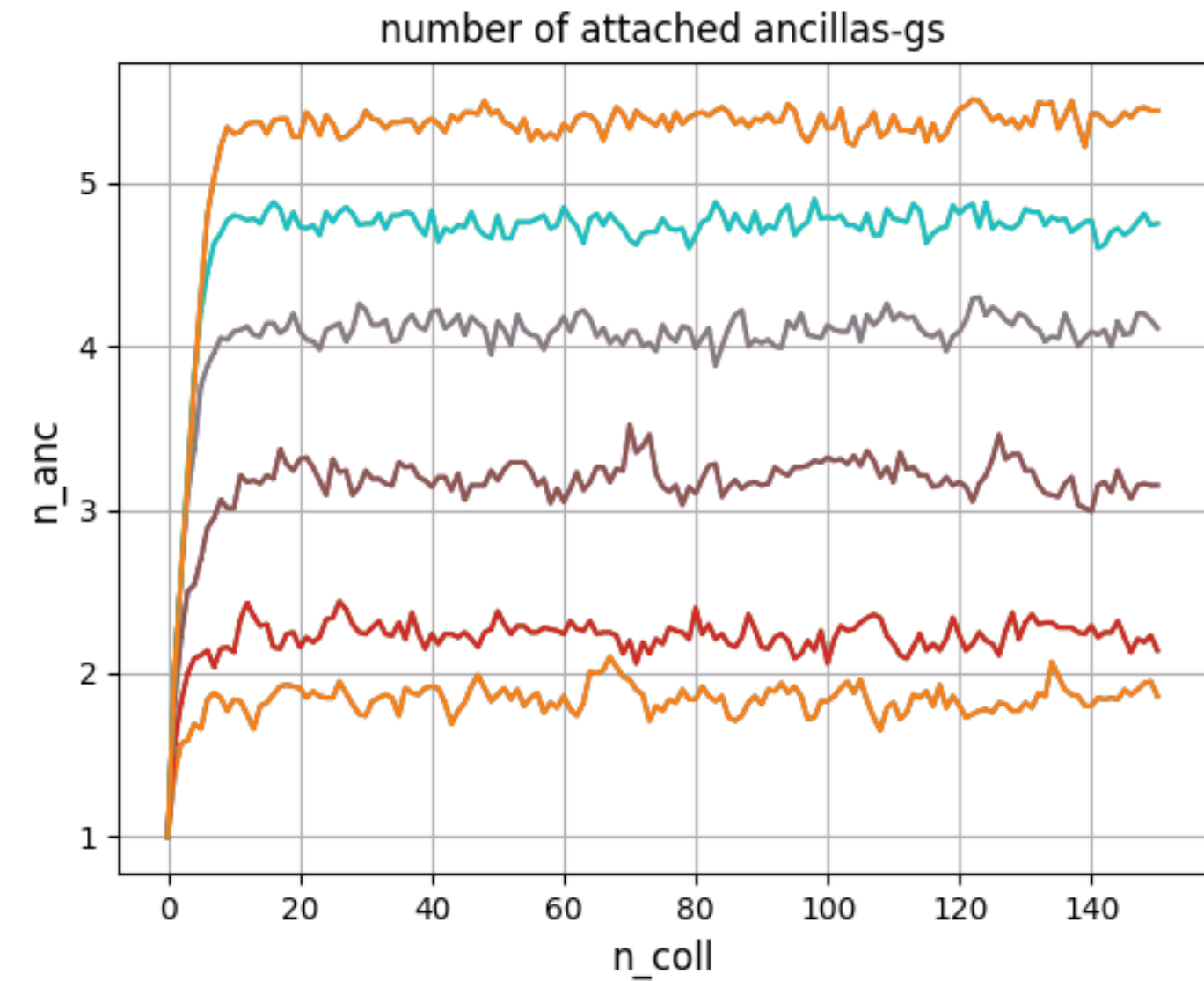
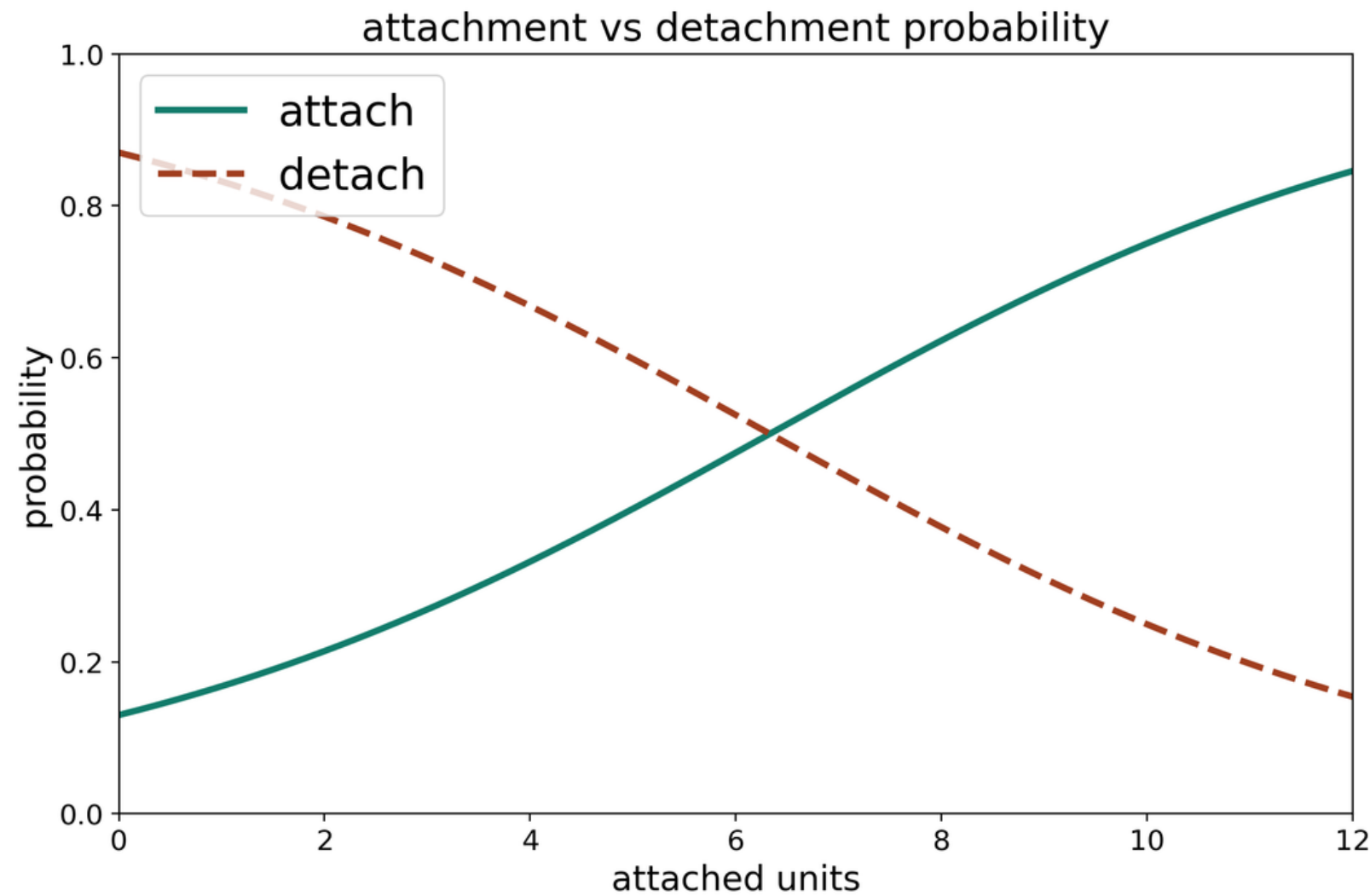
Stabilising number of auxiliary units

Question: how do we define the probabilities in the stochastic operators?



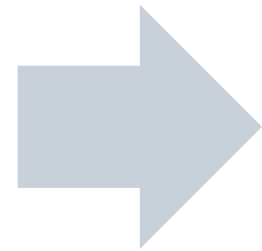
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Non-Markovianity in CluCMs

How do we distinguish two quantum states?

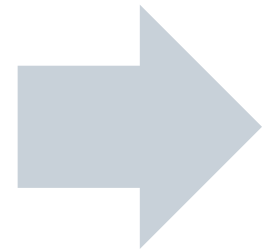


One answer is **TRACE DISTANCE**

$$D(\rho, \sigma) = \frac{1}{2} \|\rho - \sigma\|_1$$

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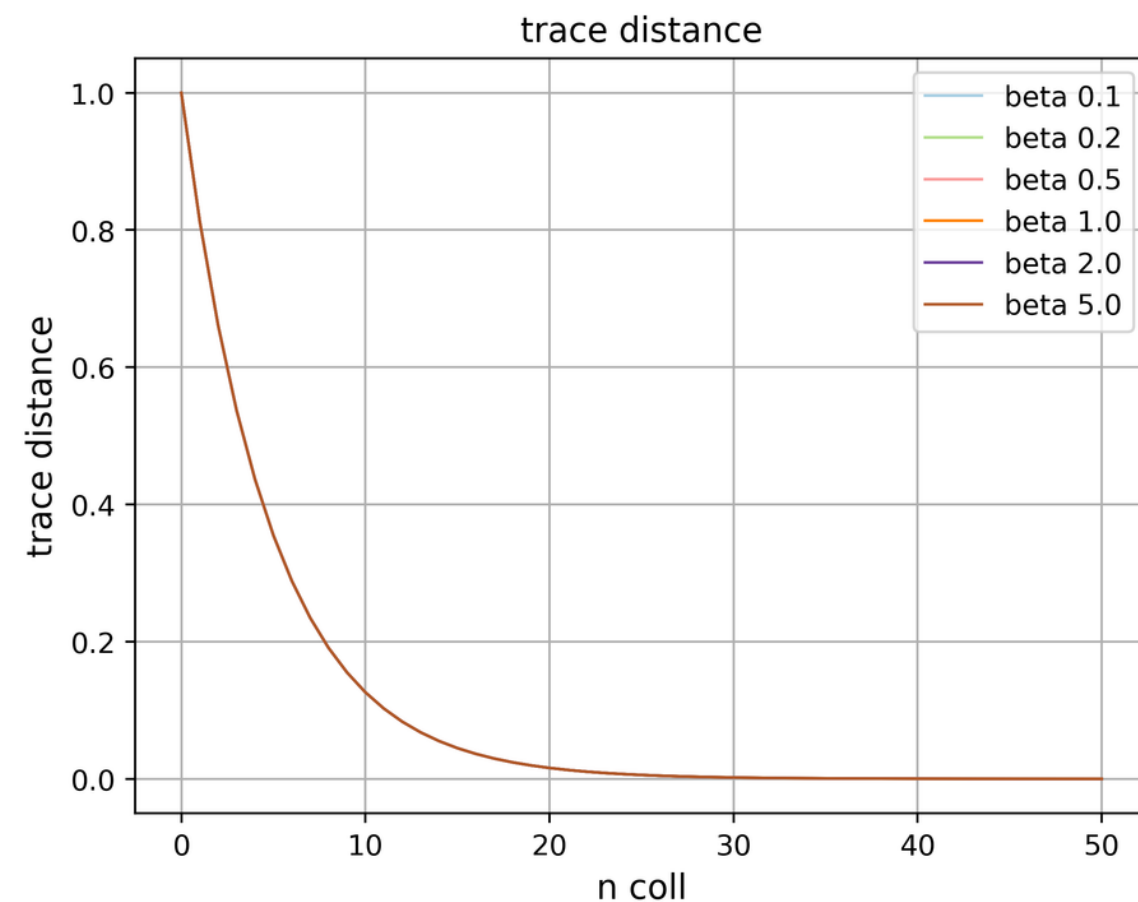
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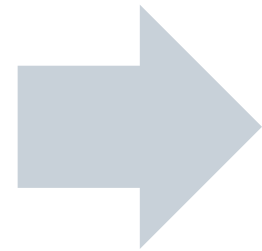
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Standard Lindblad dynamics :



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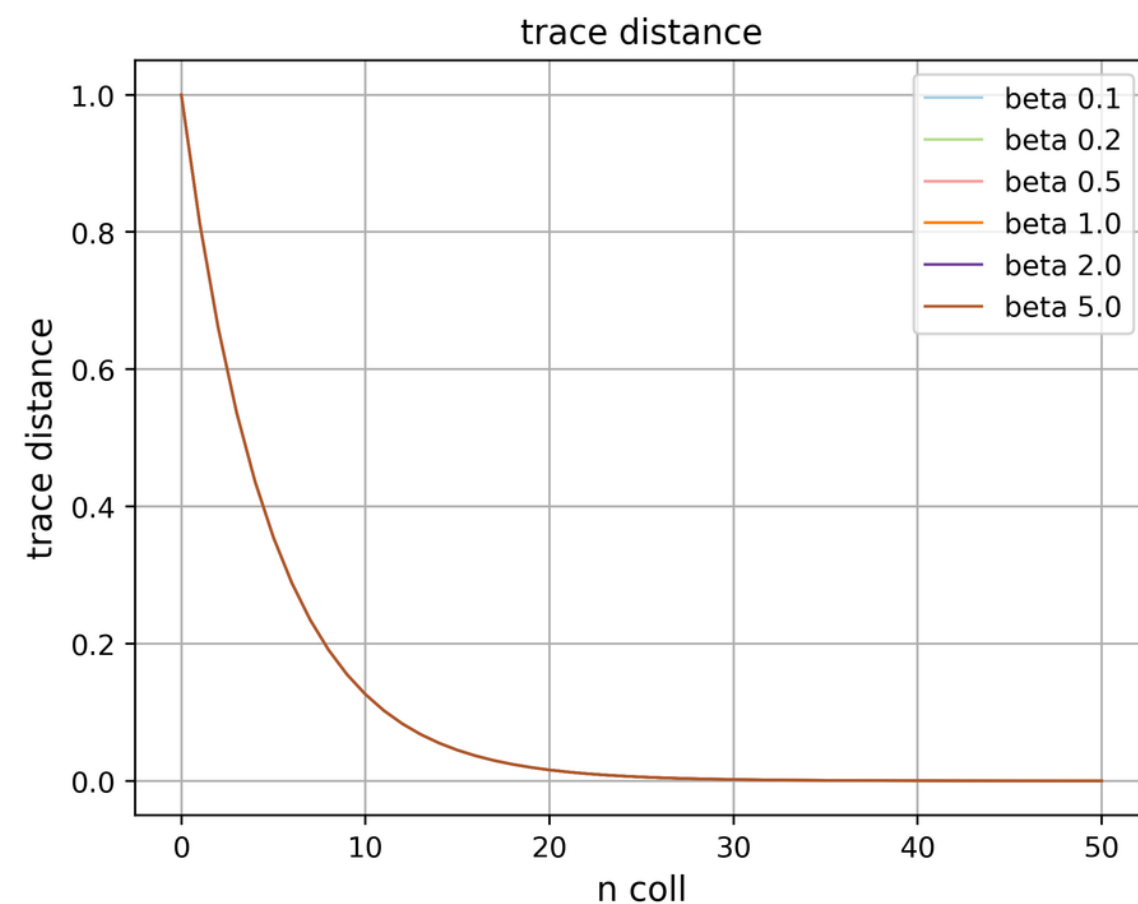
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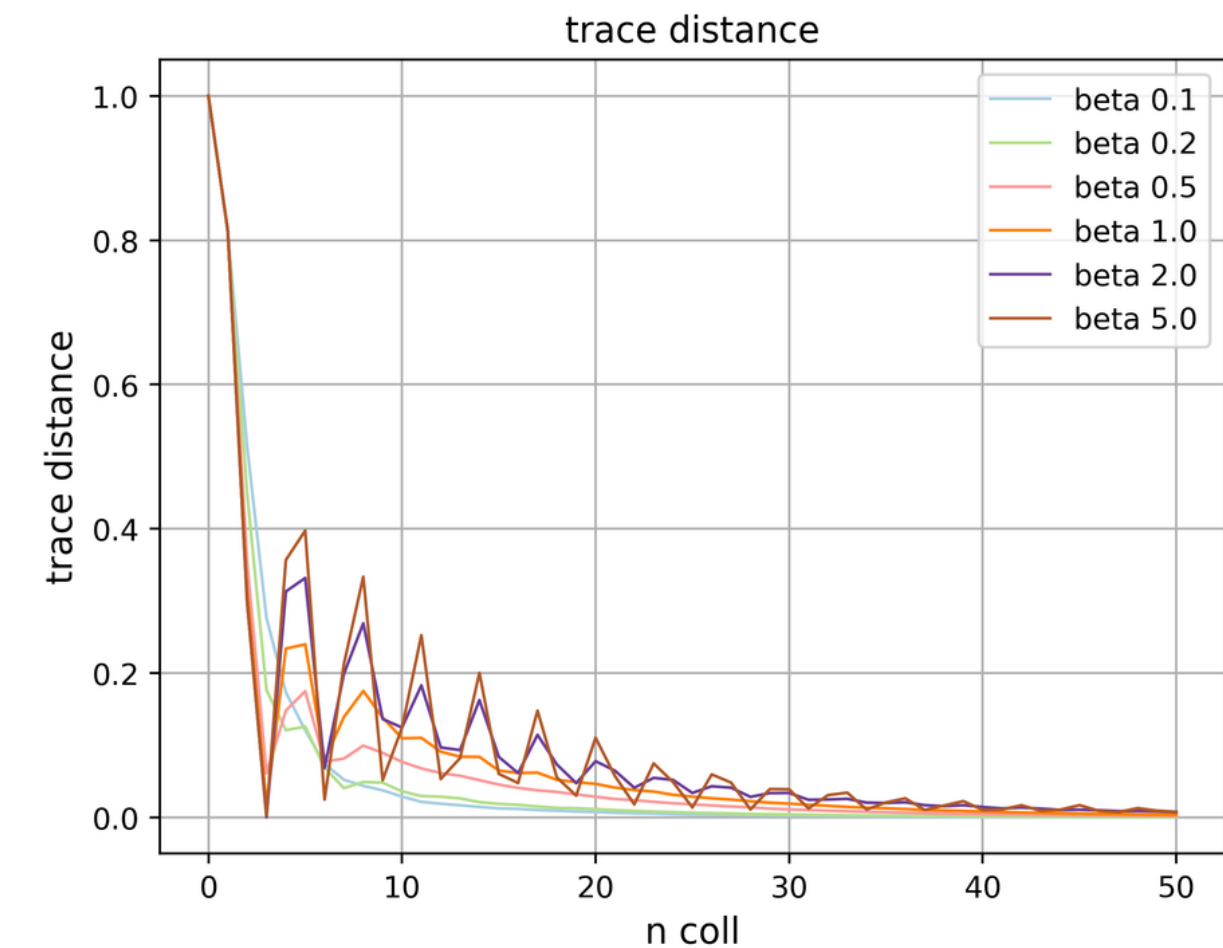
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A second detachment mechanism: natural decay

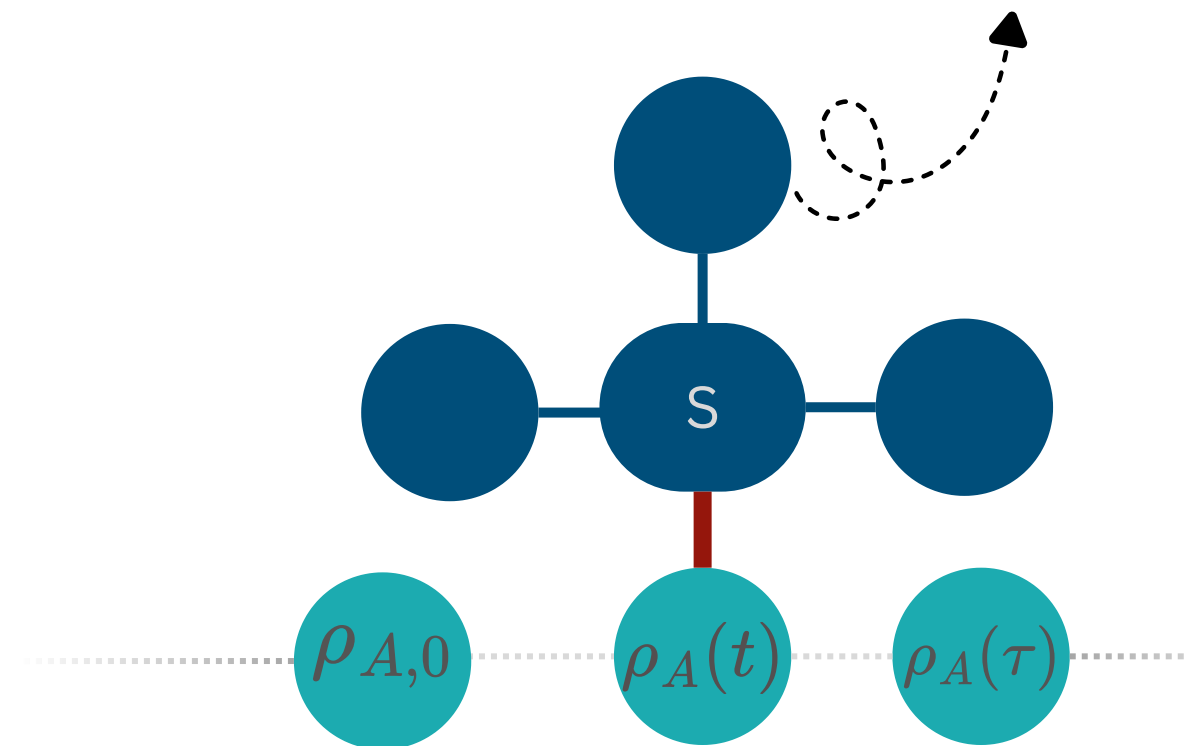
Introducing an average lifetime for the auxiliary units attached to the cluster:

$$\lambda(l, k) = \frac{l^k e^{-l}}{k!}$$

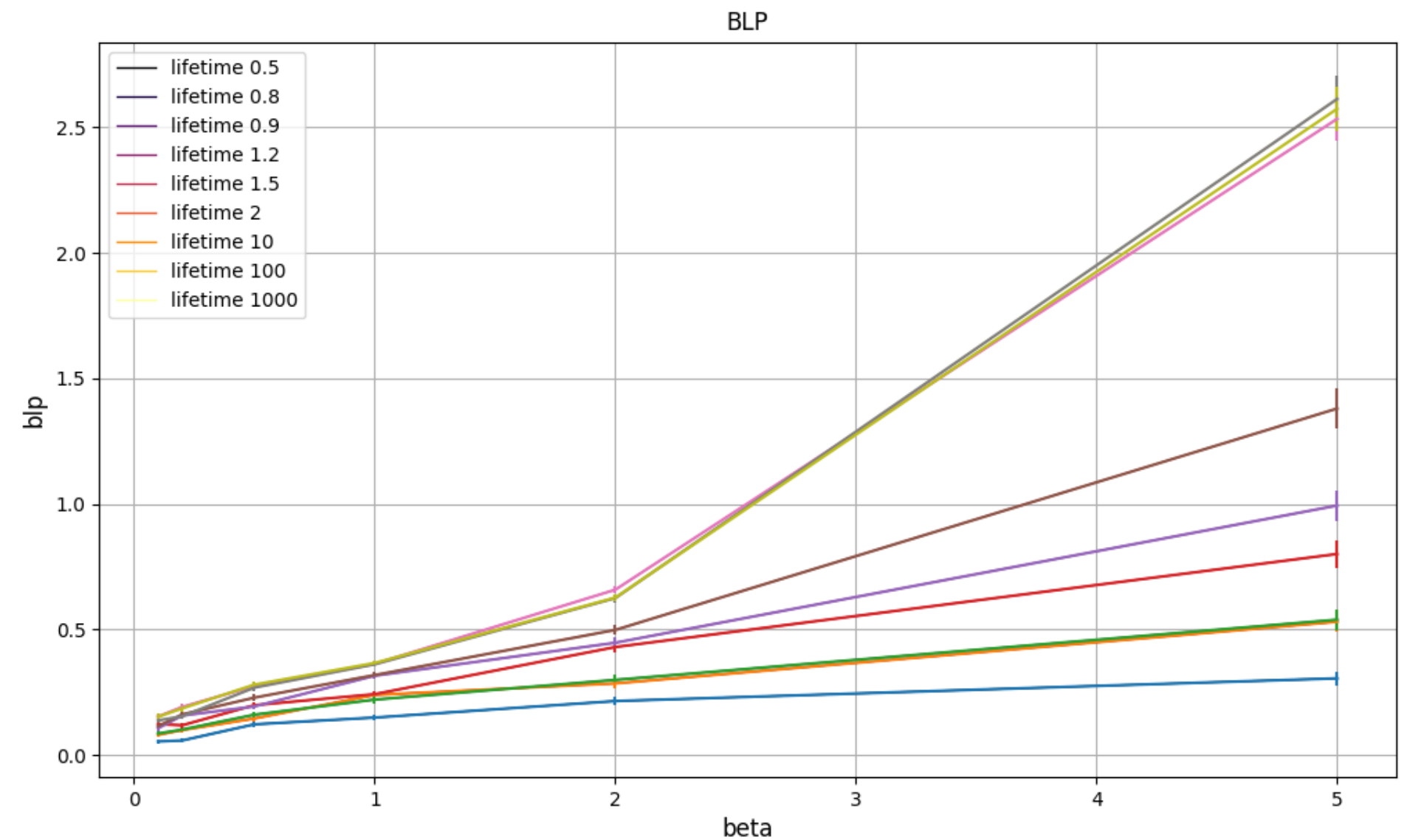
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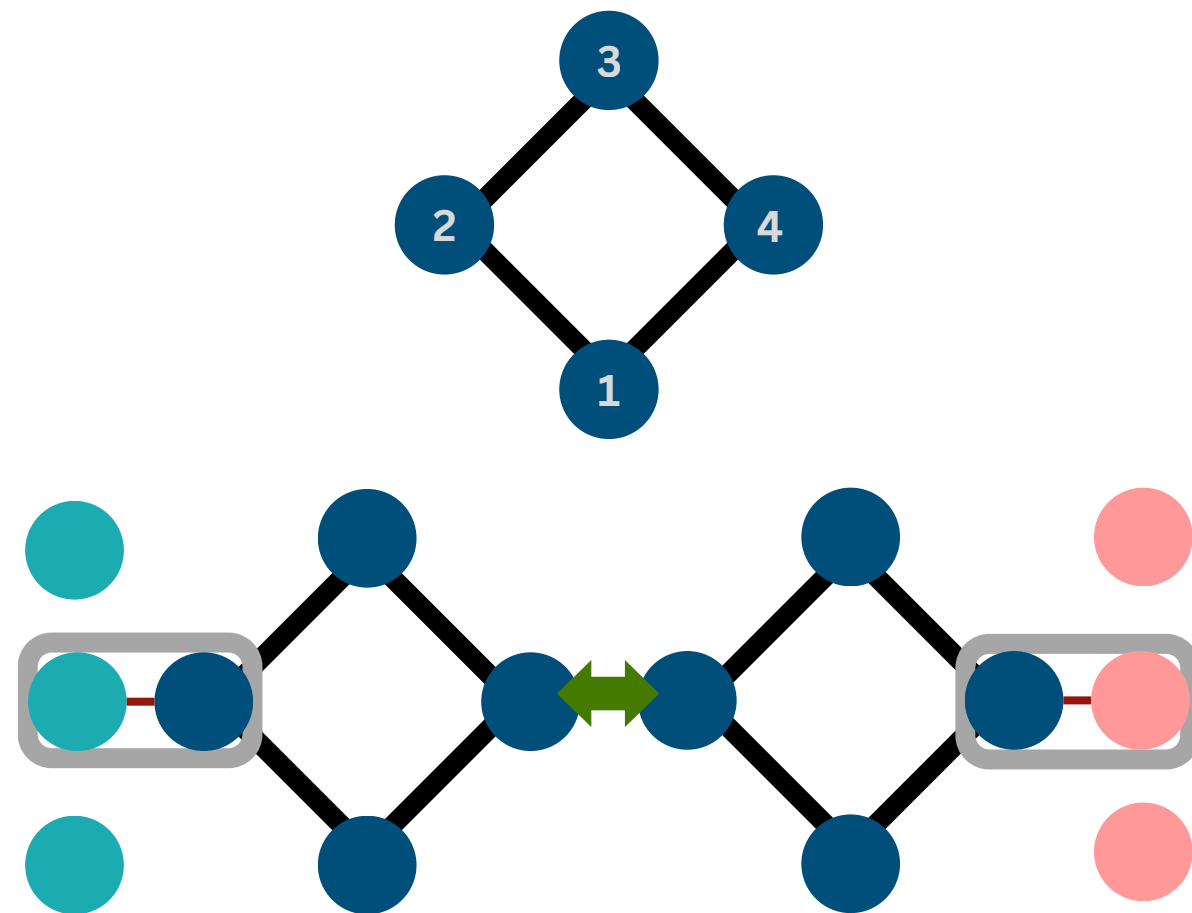
Laine et al. Physical Review A—Atomic, Molecular, and Optical Physics 81.6 (2010): 062115.



Summary and conclusions

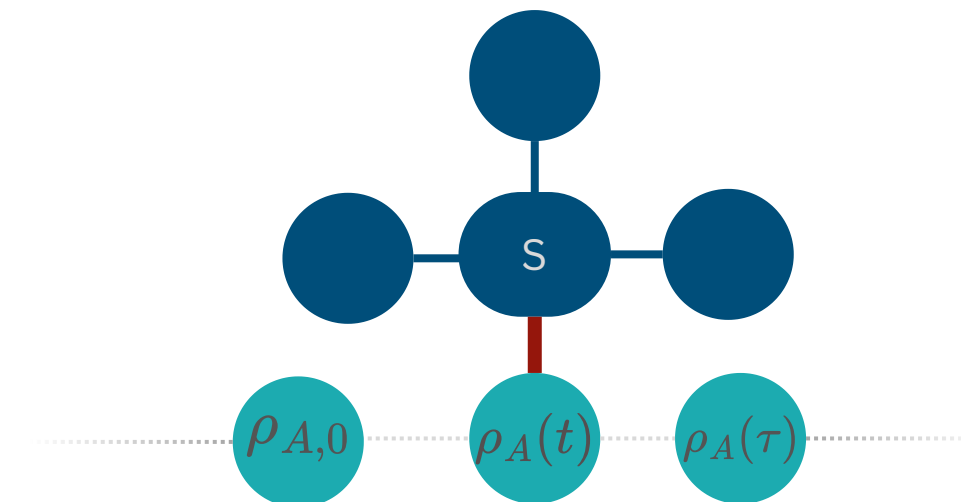
Heisenberg spin ring

- ✓ Dynamics and thermodynamics
- ✓ TN version of the system



Clustering collision models

- ✓ Dynamics of the cluster



Goal: fully characterising energy exchanges