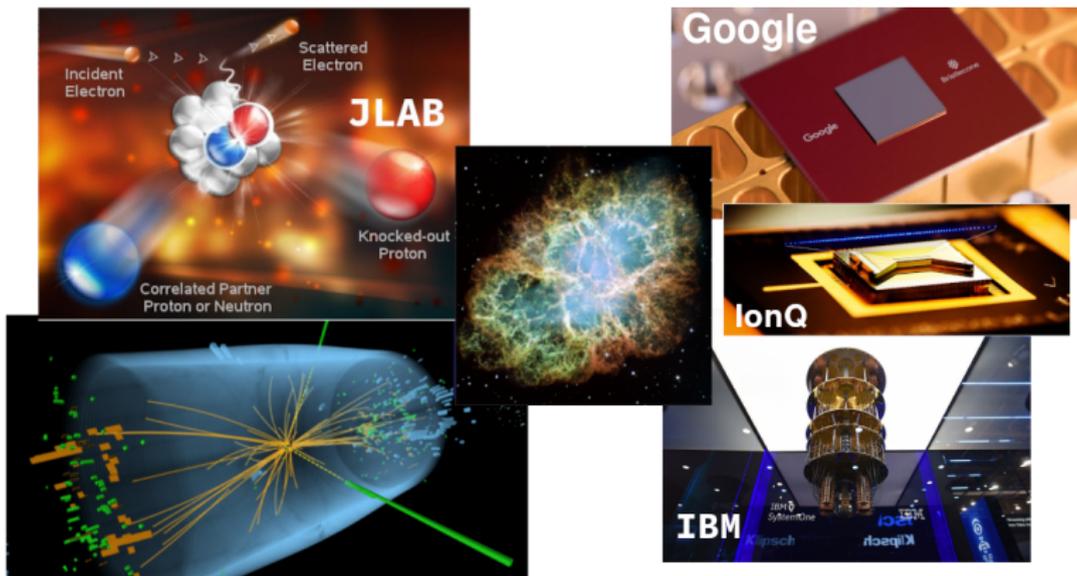


# Quantum Simulation for Nuclear and Neutrino Physics

Alessandro Roggero



Trento Institute for  
Fundamental Physics  
and Applications

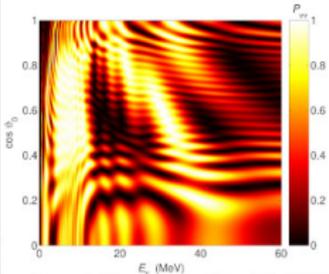
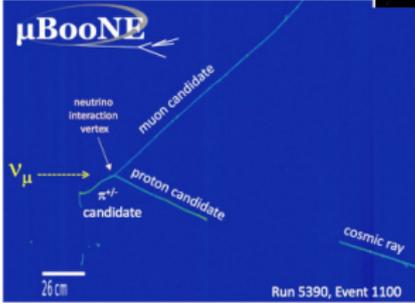
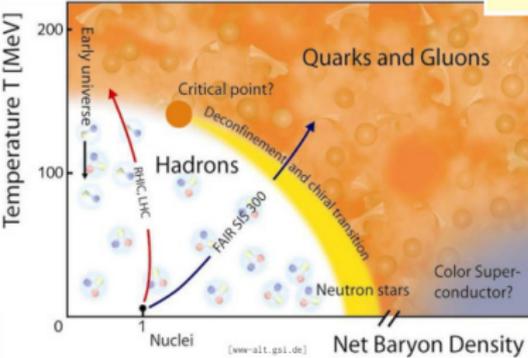
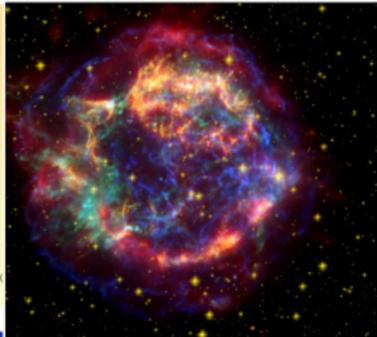
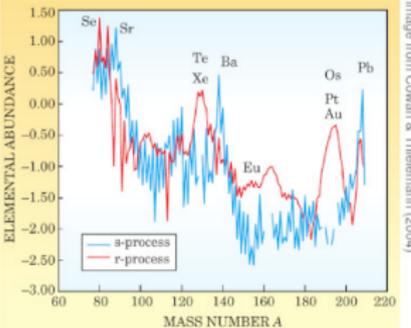
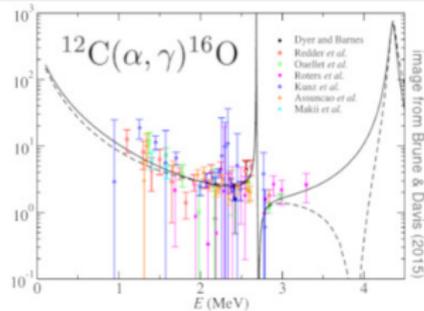
MPA Summer School

Chiemsee – 10-15 Sep, 2023



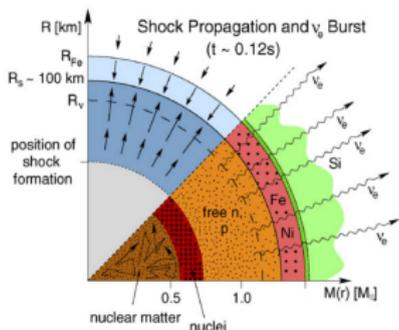
# Plan for these lectures

## Quantum Simulation for Nuclear Physics

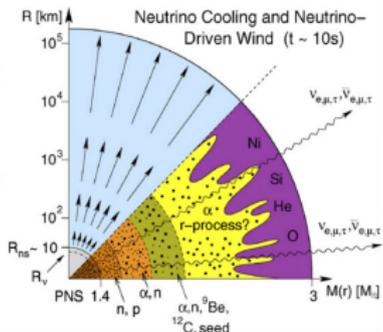


# Neutrino's roles in supernovae

- efficient energy transport away from the shock region (burst)

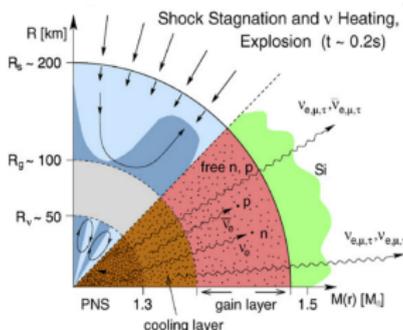


- regulation of electron fraction in  $\nu$ -driven wind (nucleosynthesis)



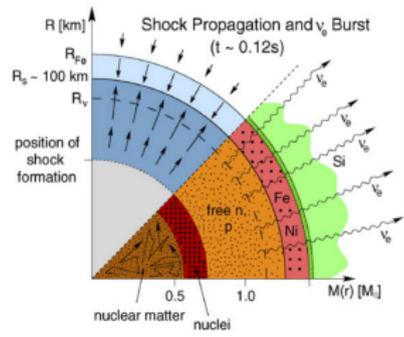
figures from Janka et al. (2007)

- energy deposition to revive the stalled shock (explosion)

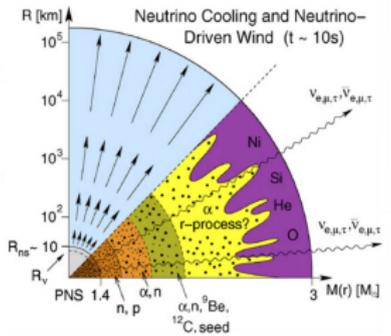


# Neutrino's roles in supernovae

- efficient energy transport away from the shock region (burst)

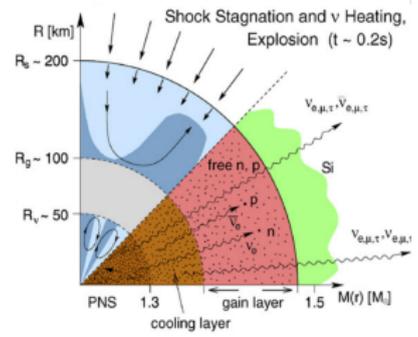


- regulation of electron fraction in  $\nu$ -driven wind (nucleosynthesis)

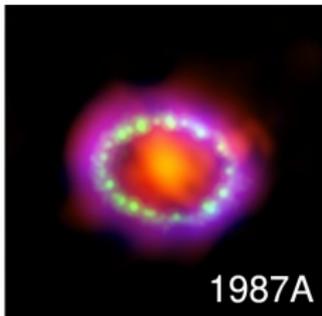
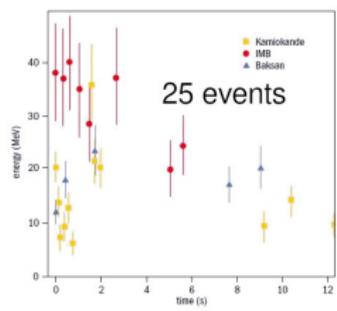


figures from Janka et al. (2007)

- energy deposition to revive the stalled shock (explosion)



$\approx 10^{58}$  neutrinos emitted in few sec.



# Neutrino oscillations in astrophysical environments

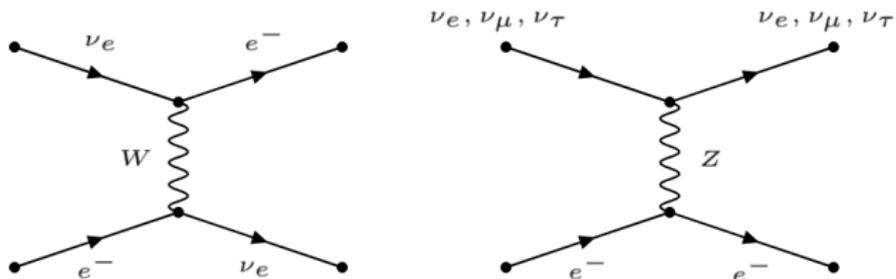
We know that neutrinos can display flavor oscillations in vacuum, does it matter in a core-collapse supernova?

- energy deposition behind shock and in the wind proceeds through charge-current reactions (large differences in  $\nu_e - \nu_{\mu/\tau}$ )

# Neutrino oscillations in astrophysical environments

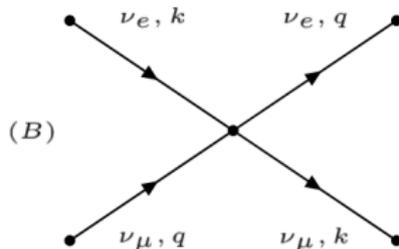
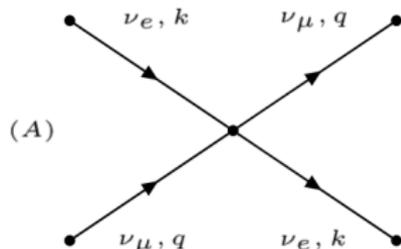
We know that neutrinos can display flavor oscillations in vacuum, does it matter in a core-collapse supernova?

- energy deposition behind shock and in the wind proceeds through charge-current reactions (large differences in  $\nu_e - \nu_{\mu/\tau}$ )
- neutrino oscillation rates can get enhanced through elastic forward scattering with high density external matter (MSW effect)



# Neutrino-neutrino forward scattering

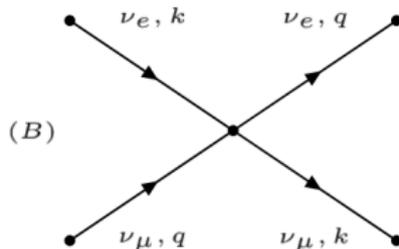
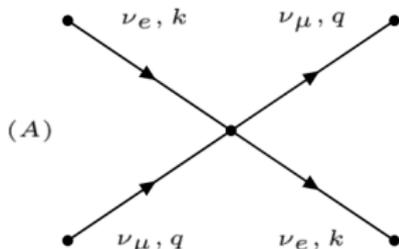
Fuller, Qian, Pantaleone, Sigl, Raffelt, Sawyer, Carlson, Duan, . . .



- diagonal contribution (A) does not impact flavor mixing
- off-diagonal term (B) equivalent to flavor/momentum exchange between two neutrinos
  - total flavor is conserved

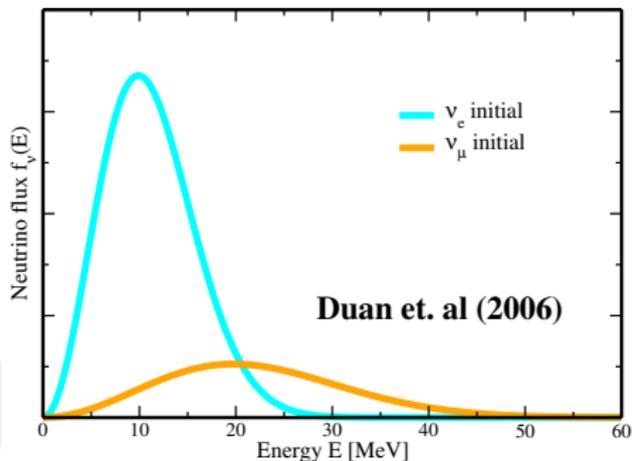
# Neutrino-neutrino forward scattering

Fuller, Qian, Pantaleone, Sigl, Raffelt, Sawyer, Carlson, Duan, ...



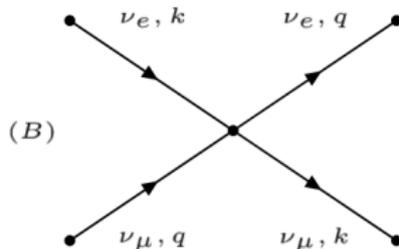
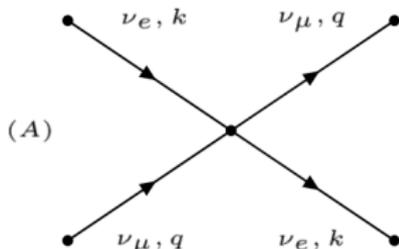
- diagonal contribution (A) does not impact flavor mixing
- off-diagonal term (B) equivalent to flavor/momentum exchange between two neutrinos
  - total flavor is conserved

Important effect if initial distributions are strongly flavor dependent



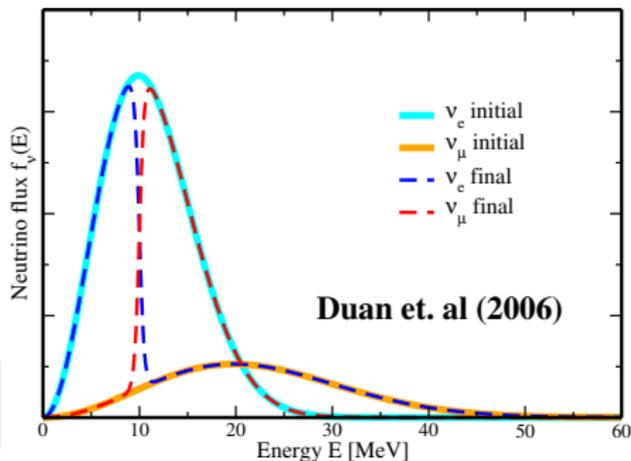
# Neutrino-neutrino forward scattering

Fuller, Qian, Pantaleone, Sigl, Raffelt, Sawyer, Carlson, Duan, ...



- diagonal contribution (A) does not impact flavor mixing
- off-diagonal term (B) equivalent to flavor/momentum exchange between two neutrinos
  - total flavor is conserved

Important effect if initial distributions are strongly flavor dependent



## Two-flavor approximation and the iso-spin Hamiltonian

Consider two active flavors ( $\nu_e, \nu_x$ ) and encode flavor amplitudes for a neutrino with momentum  $p_i$  into an  $SU(2)$  iso-spin:

$$|\Phi_i\rangle = \cos(\eta_i)|\nu_e\rangle + \sin(\eta_i)|\nu_x\rangle \equiv \cos(\eta_i)|\uparrow\rangle + \sin(\eta_i)|\downarrow\rangle$$

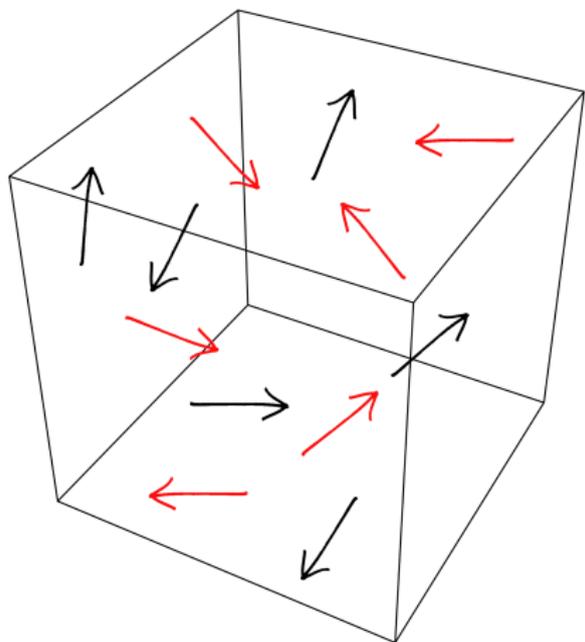
A system of  $N$  interacting neutrinos is then described by the Hamiltonian

$$H = \sum_i \frac{\Delta m^2}{4E_i} \vec{B} \cdot \vec{\sigma}_i + \lambda \sum_i \sigma_i^z + \frac{\mu}{2N} \sum_{i<j} (1 - \cos(\phi_{ij})) \vec{\sigma}_i \cdot \vec{\sigma}_j$$

- vacuum oscillations:  $\vec{B} = (\sin(2\theta_{mix}), 0, -\cos(2\theta_{mix}))$
- interaction with matter:  $\lambda = \sqrt{2}G_F\rho_e$
- neutrino-neutrino interaction:  $\mu = \sqrt{2}G_F\rho_\nu$ 
  - dependence on momentum direction:  $\cos(\phi_{ij}) = \frac{\vec{p}_i}{\|\vec{p}_i\|} \cdot \frac{\vec{p}_j}{\|\vec{p}_j\|}$

for a full derivation, see e.g. Pehlivan et al. PRD(2011)

## Finite size effects and thermodynamic limit



$$H = \sum_{i=1}^N \vec{B}_i \cdot \vec{\sigma}_i + \frac{\mu}{2N} \sum_{i < j}^N v_{ij} \vec{\sigma}_i \cdot \vec{\sigma}_j$$

- the quantum system is defined in some finite volume  $V$
- we have a finite number  $N$  of neutrinos within the box
- the neutrino density  $\rho_\nu$  (and thus  $\mu$ ) is given by  $N/V$

For astrophysically relevant predictions need to understand how the system behaves when  $V \rightarrow \infty$  and  $N \rightarrow \infty$  while keeping  $\rho_\nu = N/V$  constant

## The mean-field approximation

The equations of motion for the polarization vector  $\vec{P}_i = \langle \vec{\sigma}_i \rangle$  are

$$\frac{d}{dt} \vec{P}_i = \left( \frac{\Delta m^2}{4E_i} \vec{B} + \lambda \hat{z} \right) \times \vec{P}_i + \frac{\mu}{2N} \sum_{j \neq i} (1 - \cos(\phi_{ij})) \langle \vec{\sigma}_j \times \vec{\sigma}_i \rangle$$

## The mean-field approximation

The equations of motion for the polarization vector  $\vec{P}_i = \langle \vec{\sigma}_i \rangle$  are

$$\frac{d}{dt} \vec{P}_i = \left( \frac{\Delta m^2}{4E_i} \vec{B} + \lambda \hat{z} \right) \times \vec{P}_i + \frac{\mu}{2N} \sum_{j \neq i} (1 - \cos(\phi_{ij})) \langle \vec{\sigma}_j \times \vec{\sigma}_i \rangle$$

The mean-field approximation replaces  $\langle \vec{\sigma}_j \times \vec{\sigma}_i \rangle$  with  $\langle \vec{\sigma}_j \rangle \times \langle \vec{\sigma}_i \rangle$  so that

$$\frac{d}{dt} \vec{P}_i = \left( \frac{\Delta m^2}{4E_i} \vec{B} + \lambda \hat{z} + \frac{\mu}{2N} \sum_{j \neq i} (1 - \cos(\phi_{ij})) \vec{P}_j \right) \times \vec{P}_i$$

## The mean-field approximation

The equations of motion for the polarization vector  $\vec{P}_i = \langle \vec{\sigma}_i \rangle$  are

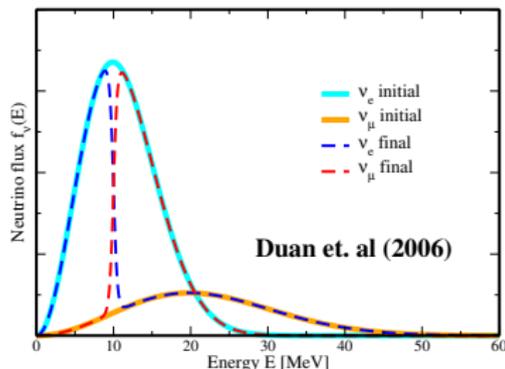
$$\frac{d}{dt} \vec{P}_i = \left( \frac{\Delta m^2}{4E_i} \vec{B} + \lambda \hat{z} \right) \times \vec{P}_i + \frac{\mu}{2N} \sum_{j \neq i} (1 - \cos(\phi_{ij})) \langle \vec{\sigma}_j \times \vec{\sigma}_i \rangle$$

The mean-field approximation replaces  $\langle \vec{\sigma}_j \times \vec{\sigma}_i \rangle$  with  $\langle \vec{\sigma}_j \rangle \times \langle \vec{\sigma}_i \rangle$  so that

$$\frac{d}{dt} \vec{P}_i = \left( \frac{\Delta m^2}{4E_i} \vec{B} + \lambda \hat{z} + \frac{\mu}{2N} \sum_{j \neq i} (1 - \cos(\phi_{ij})) \vec{P}_j \right) \times \vec{P}_i$$

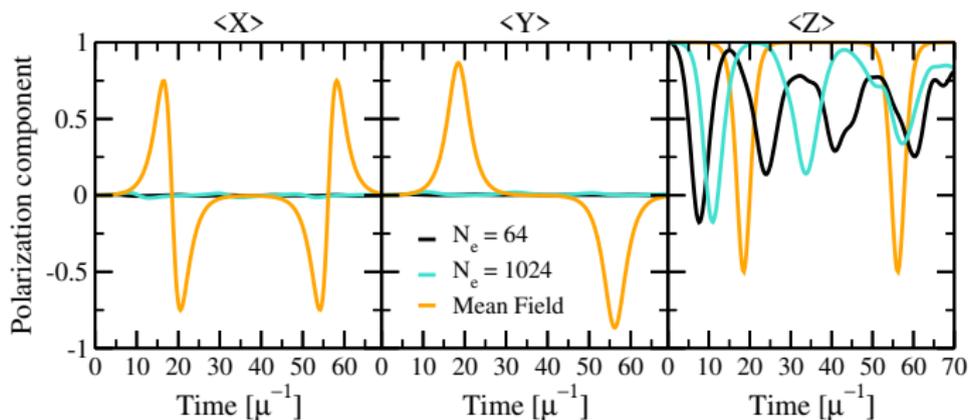
In this way we obtain a closed system of  $3N$  coupled differential equations

- efficient solutions for systems containing  $N \approx \mathcal{O}(10^{4-5})$  neutrino amplitudes [ $\approx \mathcal{O}(100)$  energies and  $\approx \mathcal{O}(100)$  angles]



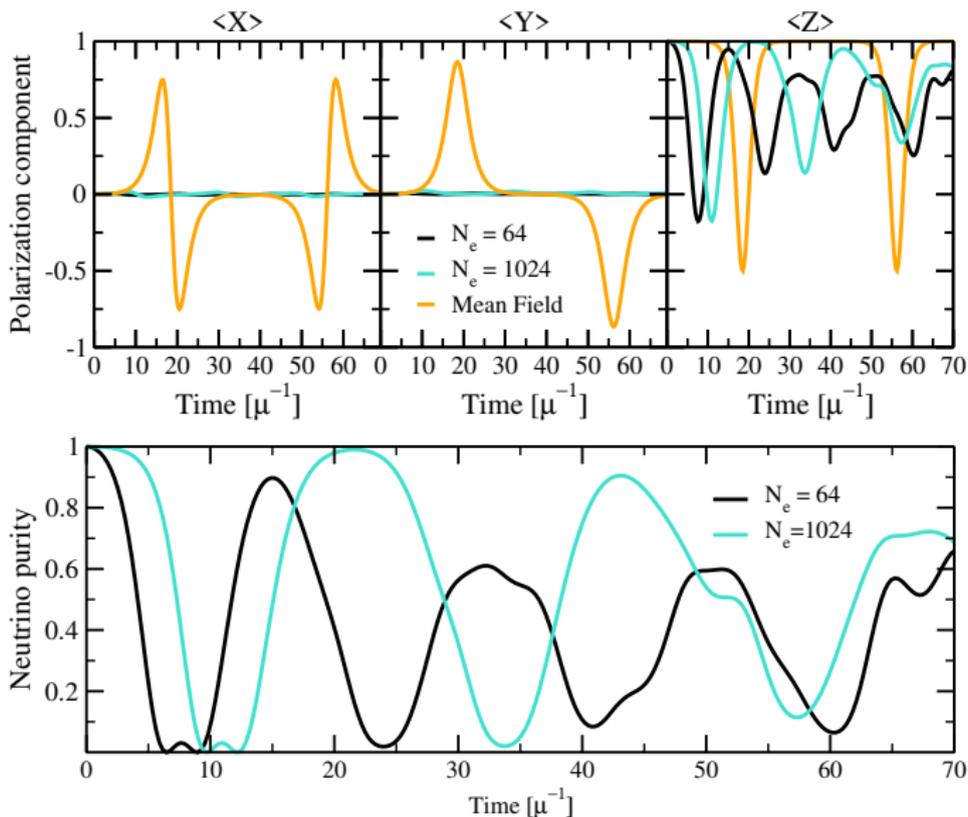
# Beyond mean field effects: a simple example

J. Martin, AR, H. Duan, J. Carlson, V. Cirigliano PRD(2022)



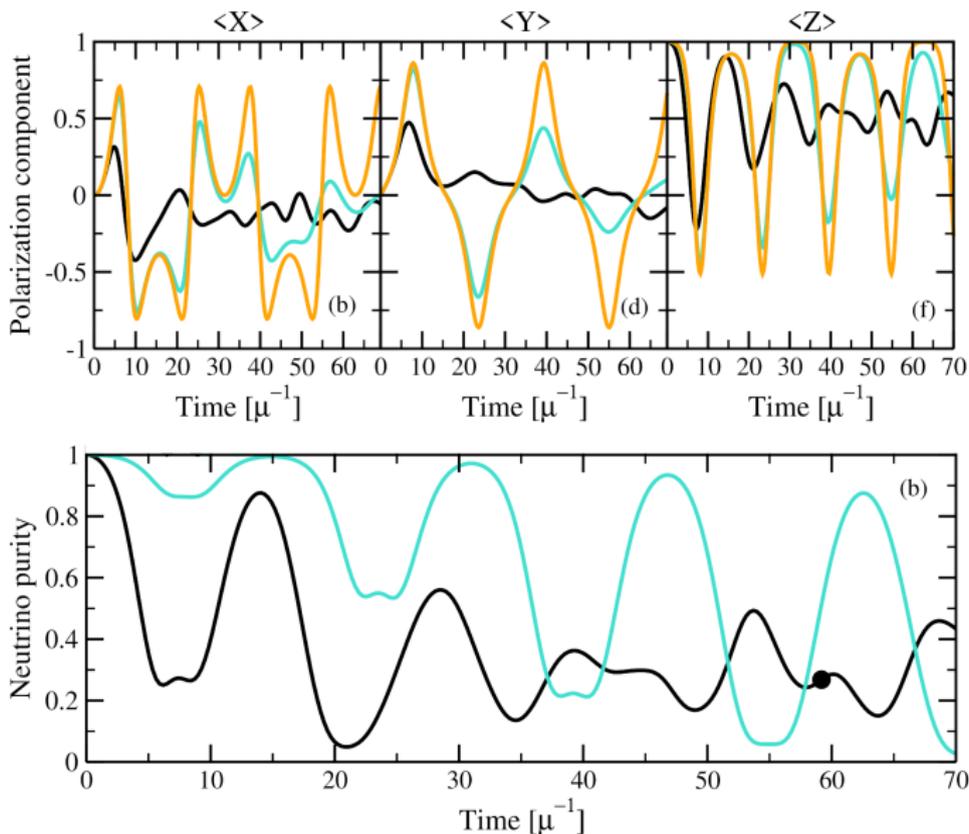
# Beyond mean field effects: a simple example

J. Martin, AR, H. Duan, J. Carlson, V. Cirigliano PRD(2022)



# Beyond mean field effects: a simple example

J. Martin, AR, H. Duan, J. Carlson, V. Cirigliano PRD(2022)

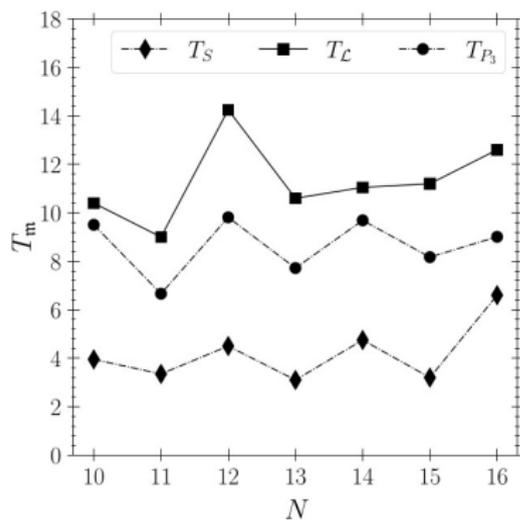
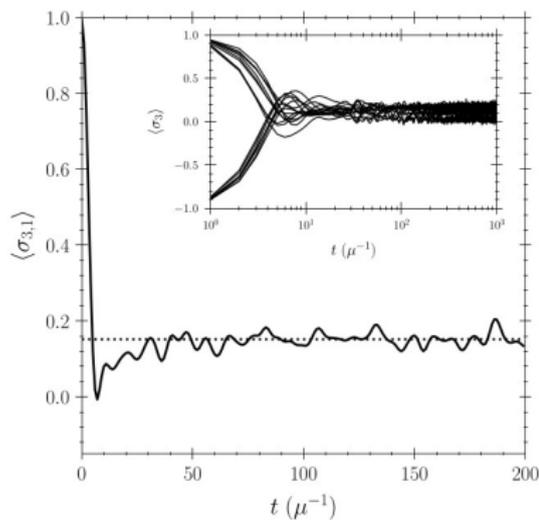


# Beyond mean field effects: a (less) simple example

J. Martin, D. Neill, AR, H. Duan, J. Carlson, arXiv:2307.16793

$$H = \frac{\mu}{2N} \sum_{i < j} (1 - \vec{v}_i \cdot \vec{v}_j) \vec{\sigma}_i \cdot \vec{\sigma}_j$$

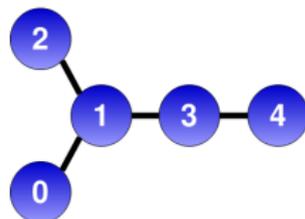
- spherical sym.:  $\vec{v}_i \cdot \vec{v}_j \rightarrow \text{const.}$
- axial sym.:  $\vec{v}_i \cdot \vec{v}_j \rightarrow v_i^z v_j^z$



Interacting neutrino systems can thermalize!

# Quantum simulation of collective neutrino oscillations

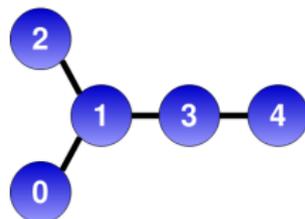
$$H = \sum_i \omega_i \vec{B} \cdot \vec{\sigma}_i + \frac{\mu}{2N} \sum_{i < j} J_{ij} \vec{\sigma}_i \cdot \vec{\sigma}_j$$



- with only 2 flavors direct map to spin 1/2 degrees of freedom (qubits)
- only one- and two-body interactions  $\Rightarrow$  only  $\mathcal{O}(N^2)$  terms
- all-to-all interactions are difficult with reduced connectivity

# Quantum simulation of collective neutrino oscillations

$$H = \sum_i \omega_i \vec{B} \cdot \vec{\sigma}_i + \frac{\mu}{2N} \sum_{i < j} J_{ij} \vec{\sigma}_i \cdot \vec{\sigma}_j$$



- with only 2 flavors direct map to spin 1/2 degrees of freedom (qubits)
- only one- and two-body interactions  $\Rightarrow$  only  $\mathcal{O}(N^2)$  terms
- all-to-all interactions are difficult with reduced connectivity

## SWAP network



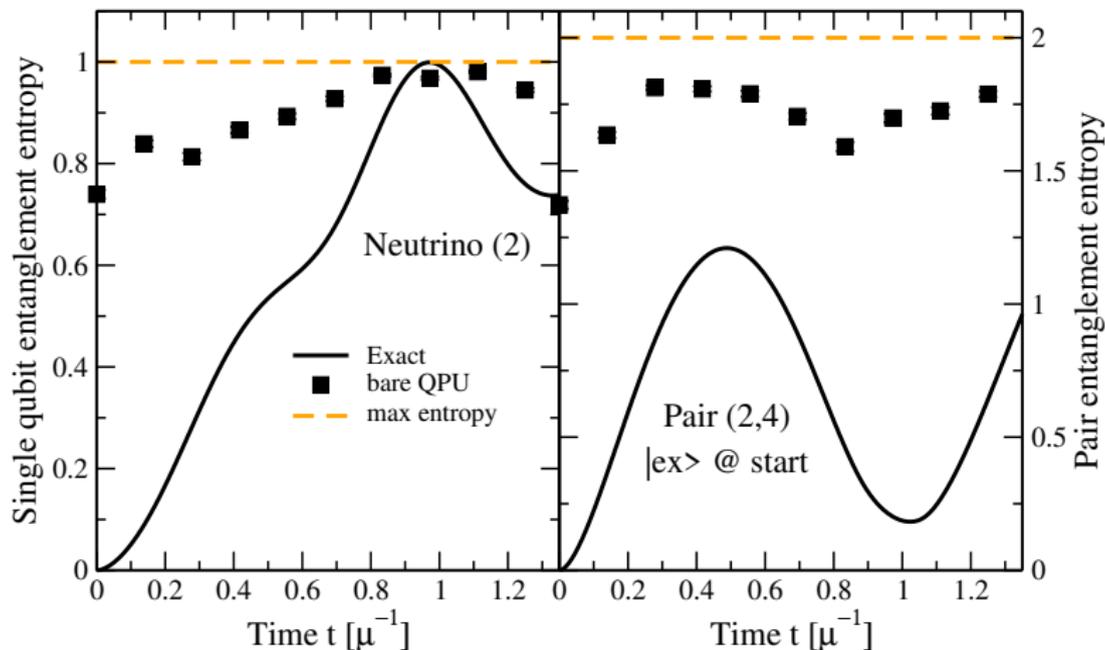
- SWAP qubits every time we apply time-evolution to neighboring terms
- in  $N$  steps we perform full evolution using only  $\binom{N}{2}$  two qubit gates
  - NOTE: final order will be reversed

Kivlichan et al. PRL (2018)

B.Hall, AR, A.Baroni, J.Carlson PRD(2021)

# Entanglement evolution and error mitigation with $N = 4$

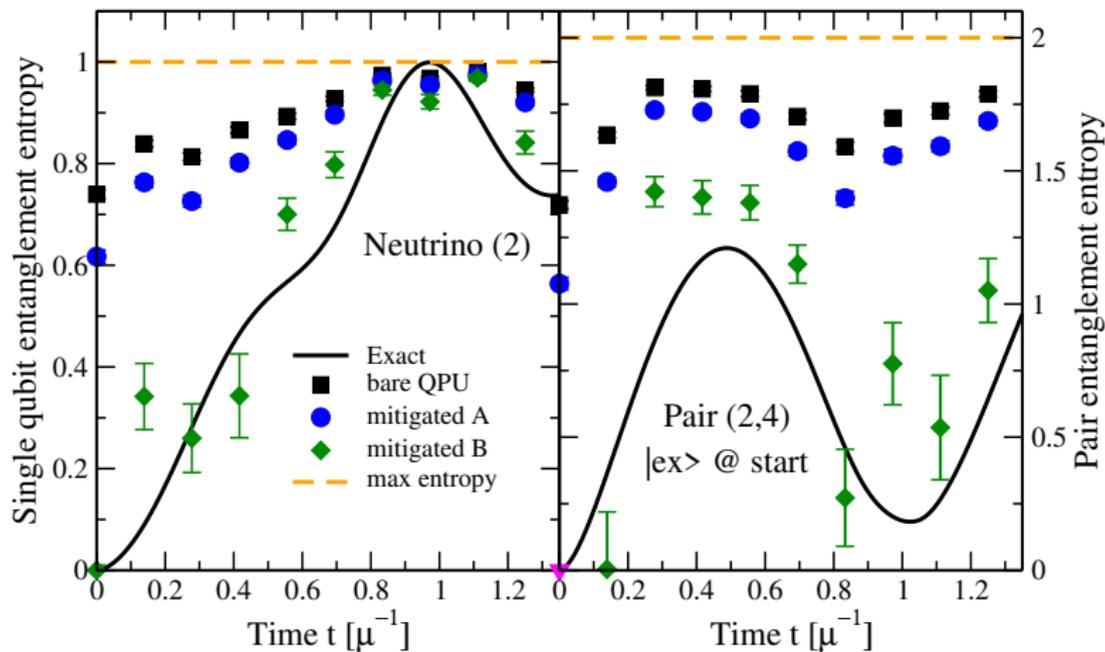
B.Hall, AR, A.Baroni, J.Carlson PRD(2021)



- Dechoerence with environment leads to increase in measured entropy

# Entanglement evolution and error mitigation with $N = 4$

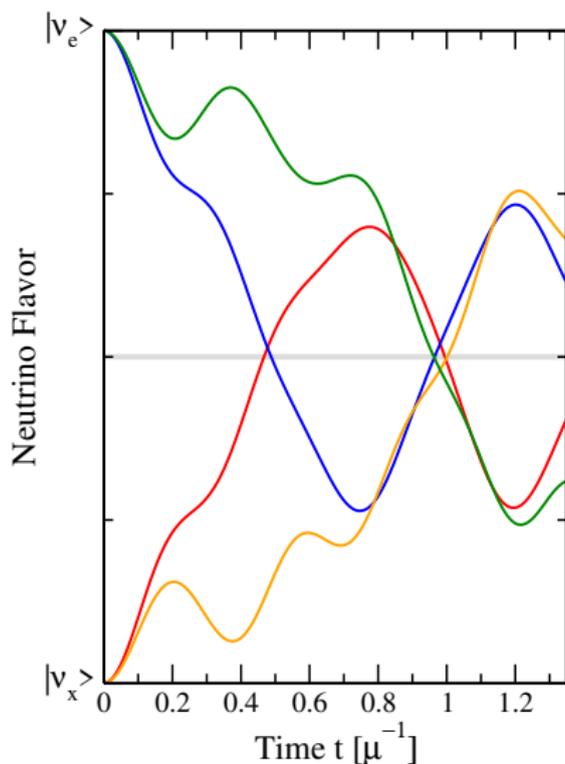
B.Hall, AR, A.Baroni, J.Carlson PRD(2021)



- Dechoerence with environment leads to increase in measured entropy
- Noise impact on observables can be modeled and effect mitigated

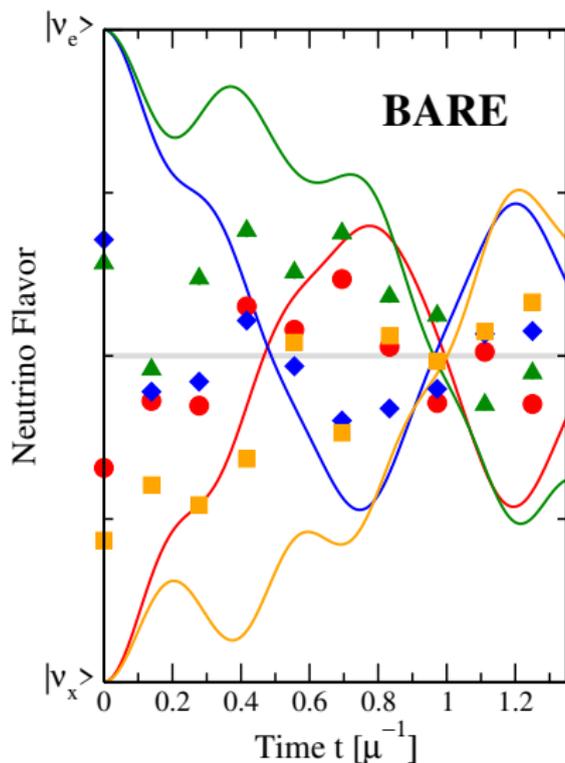
## Accuracy in flavor evolution

Entanglement is useful to understand collective oscillation mechanism but priority is to predict flavor evolution. How's the current (2020) accuracy?



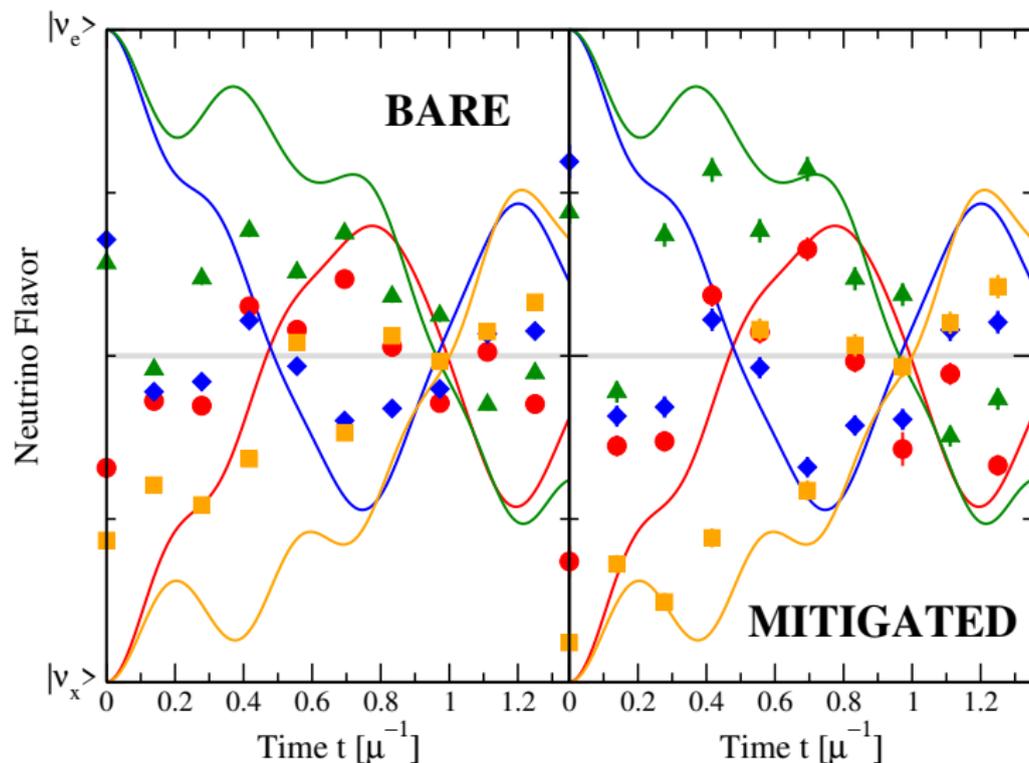
## Accuracy in flavor evolution

Entanglement is useful to understand collective oscillation mechanism but priority is to predict flavor evolution. How's the current (2020) accuracy?



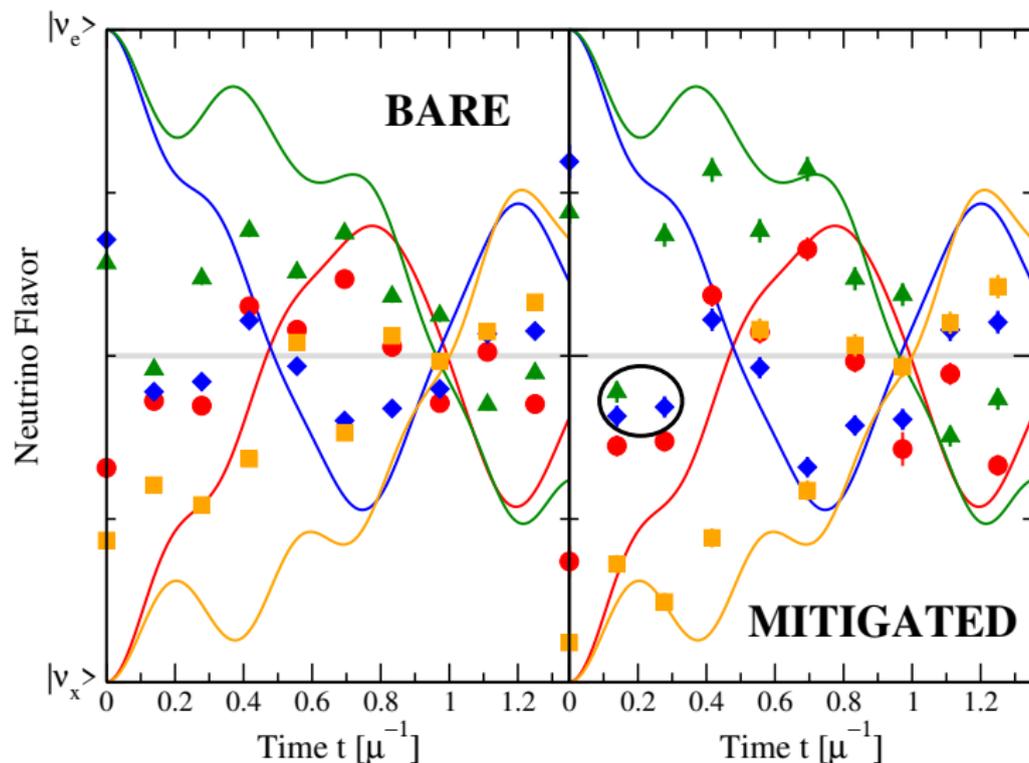
## Accuracy in flavor evolution

Entanglement is useful to understand collective oscillation mechanism but priority is to predict flavor evolution. How's the current (2020) accuracy?



## Accuracy in flavor evolution

Entanglement is useful to understand collective oscillation mechanism but priority is to predict flavor evolution. How's the current (2020) accuracy?



## Fidelity of quantum hardware is improving fast

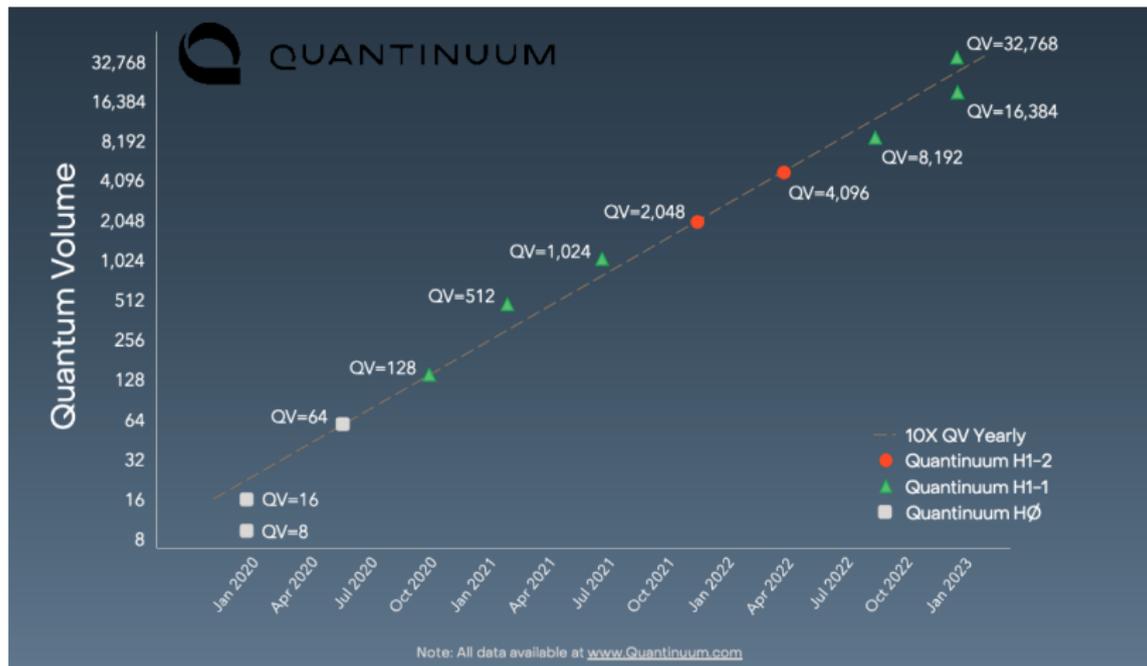
The device used for the previous results was Vigo with a QV of 16

$QV = 2^n \approx$  we can run  $n$  full layers on  $n$  qubits with fidelity  $\geq 66\%$

# Fidelity of quantum hardware is improving fast

The device used for the previous results was Vigo with a QV of 16

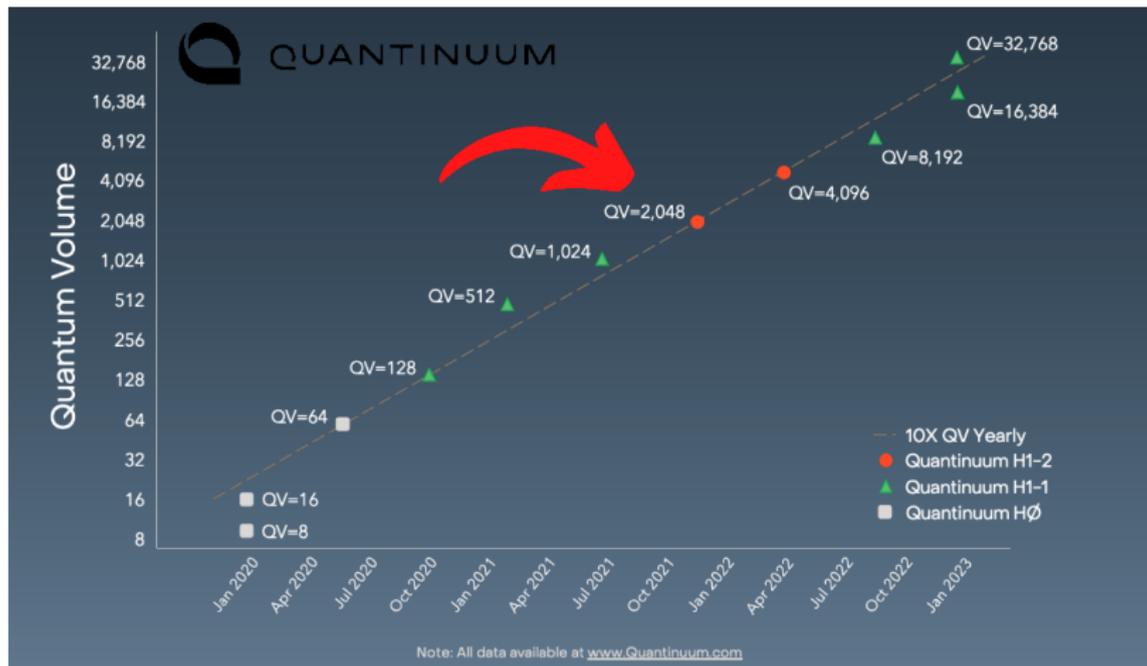
$QV = 2^n \approx$  we can run  $n$  full layers on  $n$  qubits with fidelity  $\geq 66\%$



# Fidelity of quantum hardware is improving fast

The device used for the previous results was Vigo with a QV of 16

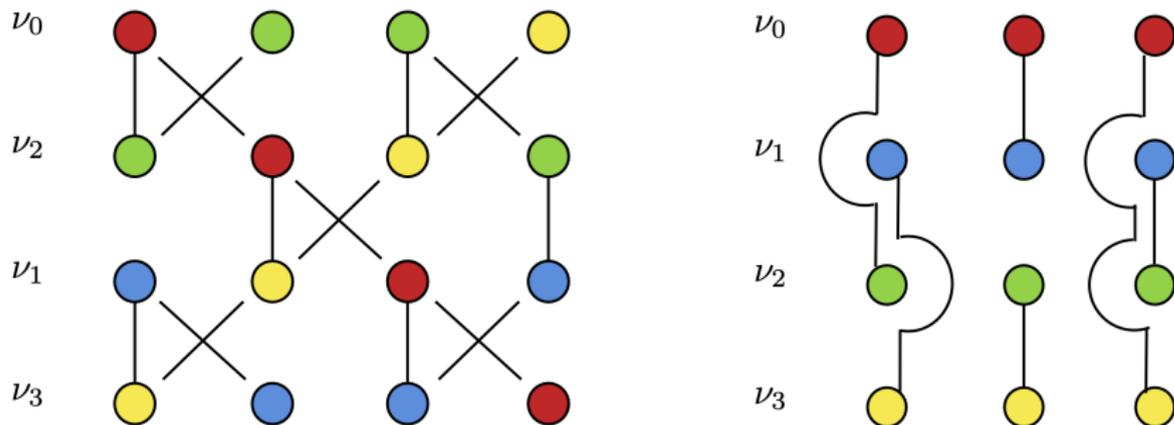
$QV = 2^n \approx$  we can run  $n$  full layers on  $n$  qubits with fidelity  $\geq 66\%$



# Connectivity advantage with trapped ions

V.Amitrano, AR, P.Luchi, F.Turro, L.Vespucci, F.Pederiva, arXiv:2207.03189 (2022)

- all-to-all connectivity allows a reduction in circuit depth and the possibility of exploring different orderings for the decomposition

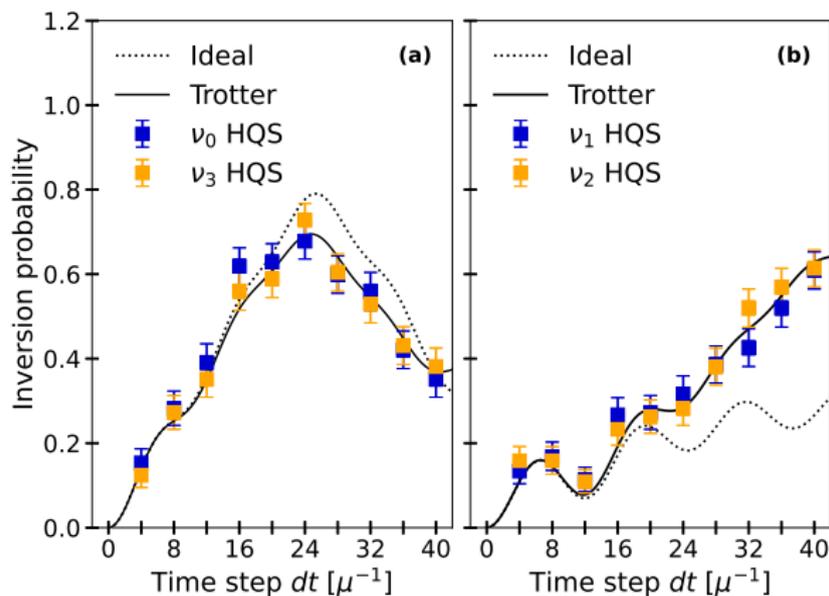


- removing SWAPs allows for a big **reduction in number of rotations**
- very low infidelities:**  $\approx 5 \times 10^{-5}$  one-qubit,  $\approx 3 \times 10^{-3}$  two-qubit

# Recent progress in porting the scheme to trapped ions

V.Amitrano, AR, P.Luchi, F.Turro, L.Vespucci, F.Pederiva, arXiv:2207.03189 (2022)

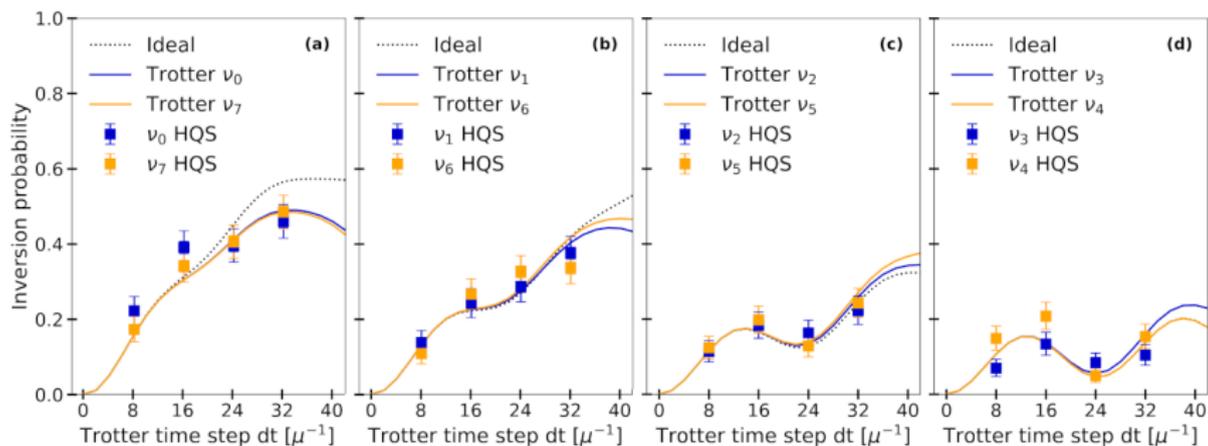
$N = 4$  neutrinos, one time step



# Recent progress in porting the scheme to trapped ions II

V.Amitrano, AR, P.Luchi, F.Turro, L.Vespucci, F.Pederiva, arXiv:2207.03189 (2022)

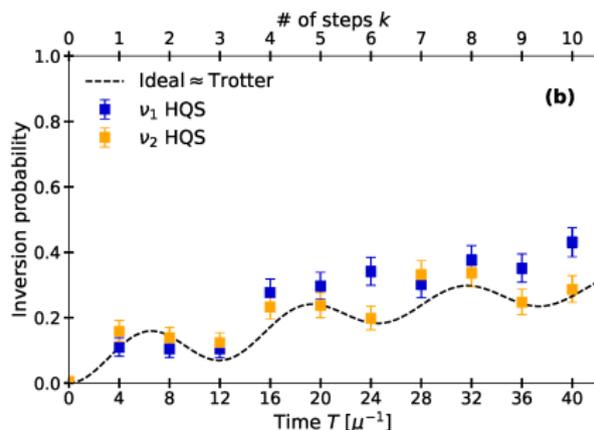
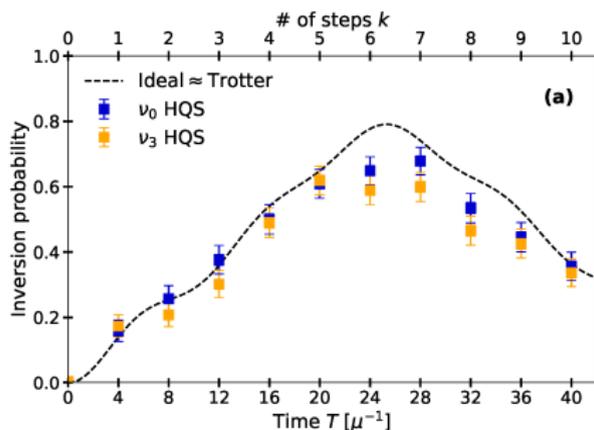
$N = 8$  neutrinos, one time step



# Recent progress in porting the scheme to trapped ions III

V.Amitrano, AR, P.Luchi, F.Turro, L.Vespucci, F.Pederiva, arXiv:2207.03189 (2022)

$N = 4$  neutrinos, multiple time steps



Last two points required:  $\approx 350$  two-qubit gates over 8 qubits

## Scaling to large system sizes

In most cases the entire cost of the simulation comes from time evolution since the initial state preparation is trivial if we start in a product state

$$H = \sum_i \omega_i \vec{B} \cdot \vec{\sigma}_i + \frac{\mu}{2N} \sum_{i < j} J_{ij} \vec{\sigma}_i \cdot \vec{\sigma}_j = H_\nu + H_{\nu\nu}$$

In simpler case where  $\omega_i = \omega$  then  $[H_\nu, H_{\nu\nu}] = 0$  (2-body dominates)

## Scaling to large system sizes

In most cases the entire cost of the simulation comes from time evolution since the initial state preparation is trivial if we start in a product state

$$H = \sum_i \omega_i \vec{B} \cdot \vec{\sigma}_i + \frac{\mu}{2N} \sum_{i < j} J_{ij} \vec{\sigma}_i \cdot \vec{\sigma}_j = H_\nu + H_{\nu\nu}$$

In simpler case where  $\omega_i = \omega$  then  $[H_\nu, H_{\nu\nu}] = 0$  (2-body dominates)

- First order Trotter:  $r = \mathcal{O}\left(\frac{T^2}{\epsilon} \sum_{ijkl} \|[H_{\nu\nu}^{ij}, H_{\nu\nu}^{kl}]\| \right) = \mathcal{O}\left(\frac{T^2 \mu^2}{\epsilon} N\right)$
- QSP [Low&Chuang(2016)]:  $r = \mathcal{O}\left(T \lambda_H + \log\left(\frac{1}{\epsilon}\right)\right) = \mathcal{O}\left(T \mu N + \log\left(\frac{1}{\epsilon}\right)\right)$

For both schemes the gate cost of one step is  $\mathcal{O}(N^2) \Rightarrow$  QSP better

## Scaling to large system sizes

In most cases the entire cost of the simulation comes from time evolution since the initial state preparation is trivial if we start in a product state

$$H = \sum_i \omega_i \vec{B} \cdot \vec{\sigma}_i + \frac{\mu}{2N} \sum_{i < j} J_{ij} \vec{\sigma}_i \cdot \vec{\sigma}_j = H_\nu + H_{\nu\nu}$$

In simpler case where  $\omega_i = \omega$  then  $[H_\nu, H_{\nu\nu}] = 0$  (2-body dominates)

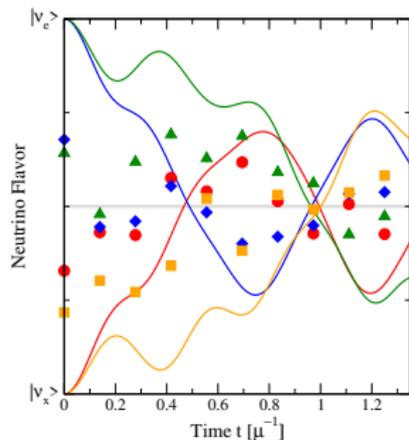
- First order Trotter:  $r = \mathcal{O}\left(\frac{T^2}{\epsilon} \sum_{ijkl} \|[H_{\nu\nu}^{ij}, H_{\nu\nu}^{kl}]\| \right) = \mathcal{O}\left(\frac{T^2 \mu^2}{\epsilon} N\right)$
- QSP [Low&Chuang(2016)]:  $r = \mathcal{O}\left(T \lambda_H + \log\left(\frac{1}{\epsilon}\right)\right) = \mathcal{O}\left(T \mu N + \log\left(\frac{1}{\epsilon}\right)\right)$

For both schemes the gate cost of one step is  $\mathcal{O}(N^2) \Rightarrow$  QSP better

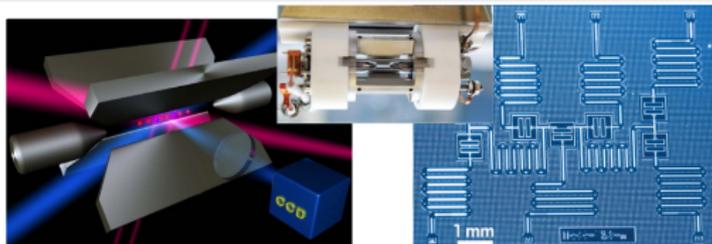
- Second order Trotter:  $r = \mathcal{O}\left(\frac{T^{3/2} \mu^{3/2}}{\sqrt{\epsilon}} \sqrt{N}\right)$

High order Trotter formulas achieve better gate cost for large  $N$ !

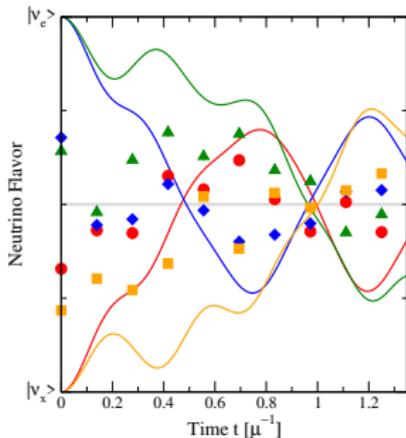
# Current limitations of digital quantum simulations



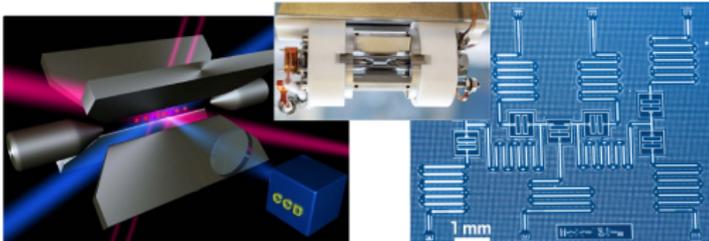
current and near term digital quantum devices have limited fidelity and might not scale much beyond  $N = \mathcal{O}(10)$  neutrinos in next years



# Current limitations of digital quantum simulations



current and near term digital quantum devices have limited fidelity and might not scale much beyond  $N = \mathcal{O}(10)$  neutrinos in next years



## Possible paths to scalability in the meantime

- Analog Quantum Simulators

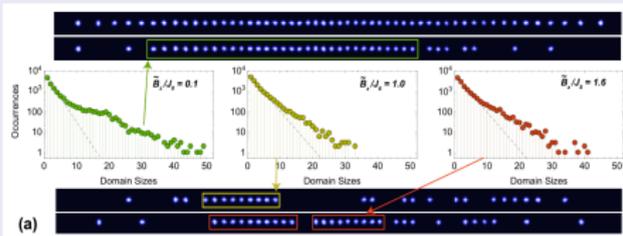
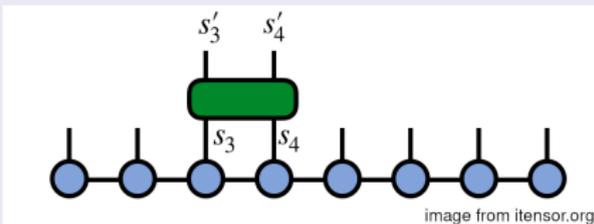


figure from Zhang et al Nature(2017)

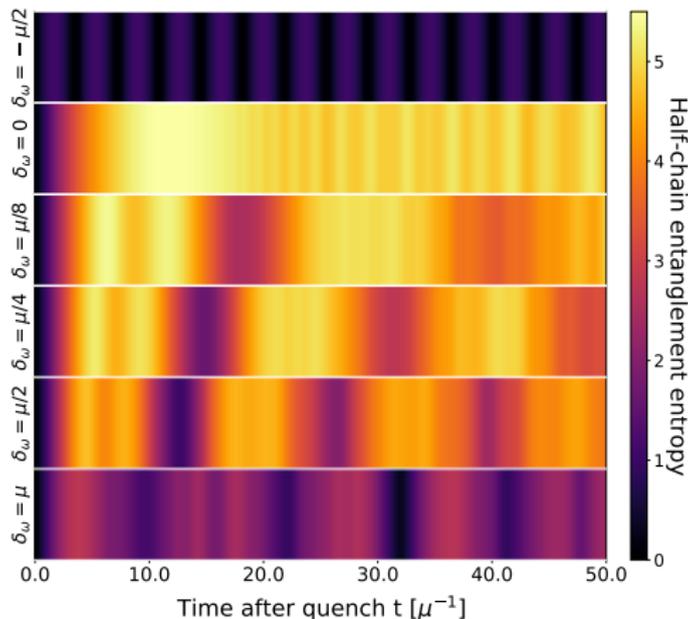
- Describe low entanglement states with Tensor Networks



## Collective oscillations with MPS

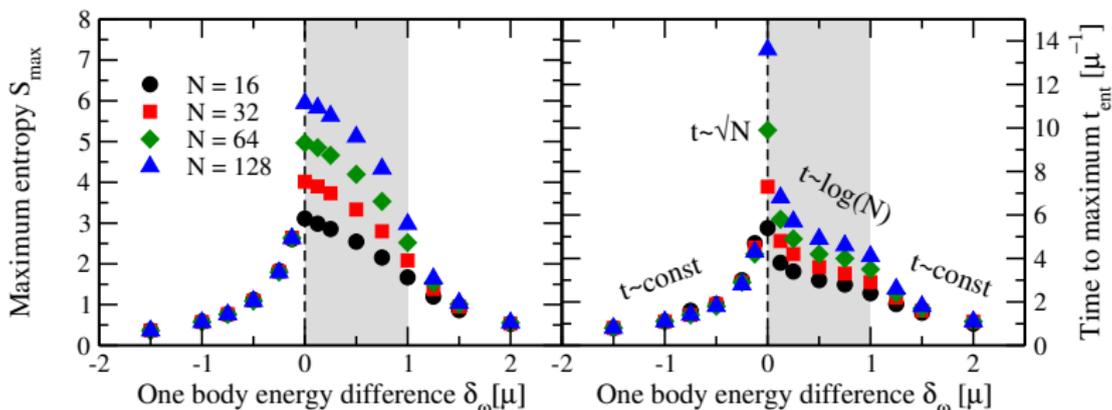
$$H = -\frac{\delta_\omega}{2} \left( \sum_{i \in \{1, \dots, N/2\}} \sigma_i^z - \sum_{i \in \{N/2+1, \dots, N\}} \sigma_i^z \right) + \frac{\mu}{2N} \sum_{i < j} \vec{\sigma}_i \cdot \vec{\sigma}_j ,$$

MF predicts no evolution, MPS has oscillations for  $0 \leq \delta_\omega/\mu \lesssim 1$



# Collective oscillations and entanglement scaling

AR, PRD 104, 103016 (2021) & PRD 104, 123023 (2021)



## Why is this interesting?

- entanglement scaling provides general criterion for appearance of collective modes in full many-body treatment
- entropy scaling as  $\log(N) \Rightarrow$  large ab-initio simulations possible
- MPS method fails when entanglement too large  $\Rightarrow$  we can use this to detect interesting regimes to study on quantum simulators!

## Summary and perspectives

- collective neutrino oscillations are an interesting **strongly coupled** many-body system driven by the **weak interaction** with possible important impact on flavor dynamics in extreme environments
- even the basic 2-flavor model for collective oscillations poses a challenging many-body problem well suited to quantum technologies
  - Hamiltonian is two-local but all-to-all → best suited for trapped-ions
- first calculations on small scale digital devices show promise in studying flavor evolution and achievable fidelity is advancing at a rapid pace (recent  $N = 12$  simulation [Illa & Savage, PRL (2023)])
- analog trapped ion devices are an ideal platform to study mid-size systems as the interactions can be embedded in a natural way
- tensor network methods can help push the boundary of classical simulations and identify interesting regimes to study with simulators
- can the spin-model describe neutrinos in supernovae correctly?

# Error mitigation with zero-noise extrapolation

Li & Benjamin PRX(2017), Temme, Bravy, Gambetta PRL(2017), Endo, Benjamin, Li PRX(2018)

## Zero noise extrapolation

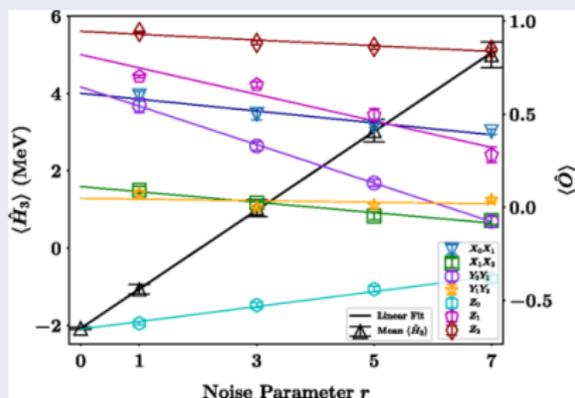
For small enough noise we can write

$$M(\epsilon) = M_0 + \epsilon M_1 + \frac{\epsilon^2}{2} M_2 + \dots$$

Using two points  $\epsilon_2 = \eta \epsilon_1$  we have

$$M_0 \approx M(\epsilon_1) - \frac{M(\epsilon_1) - M(\epsilon_2)}{\eta - 1}$$

picture from Dumitrescu et al. PRL(2018)



- for moderate  $\epsilon$  other parametrizations (like exp) might be more useful

$$M(\epsilon) = M_0 e^{-\alpha \epsilon} \Rightarrow M_0 \approx M(\epsilon_1) \left( \frac{M(\epsilon_2)}{M(\epsilon_1)} \right)^{\frac{\epsilon_1}{\epsilon_1 - \epsilon_2}}$$

In that case it is very beneficial to ensure  $M(\epsilon \rightarrow \infty) \rightarrow 0$  (mitigated B)

## Collective oscillations with MPS

$$H = -\frac{\delta_\omega}{2} \left( \sum_{i \in \{1, \dots, N/2\}} \sigma_i^z - \sum_{i \in \{N/2+1, \dots, N\}} \sigma_i^z \right) + \frac{\mu}{2N} \sum_{i < j} \vec{\sigma}_i \cdot \vec{\sigma}_j,$$

MF predicts no evolution, MPS has oscillations for  $0 \leq \delta_\omega/\mu \lesssim 1$

