Mass-imbalanced mixture of a few strongly interacting fermions driven through a critical point

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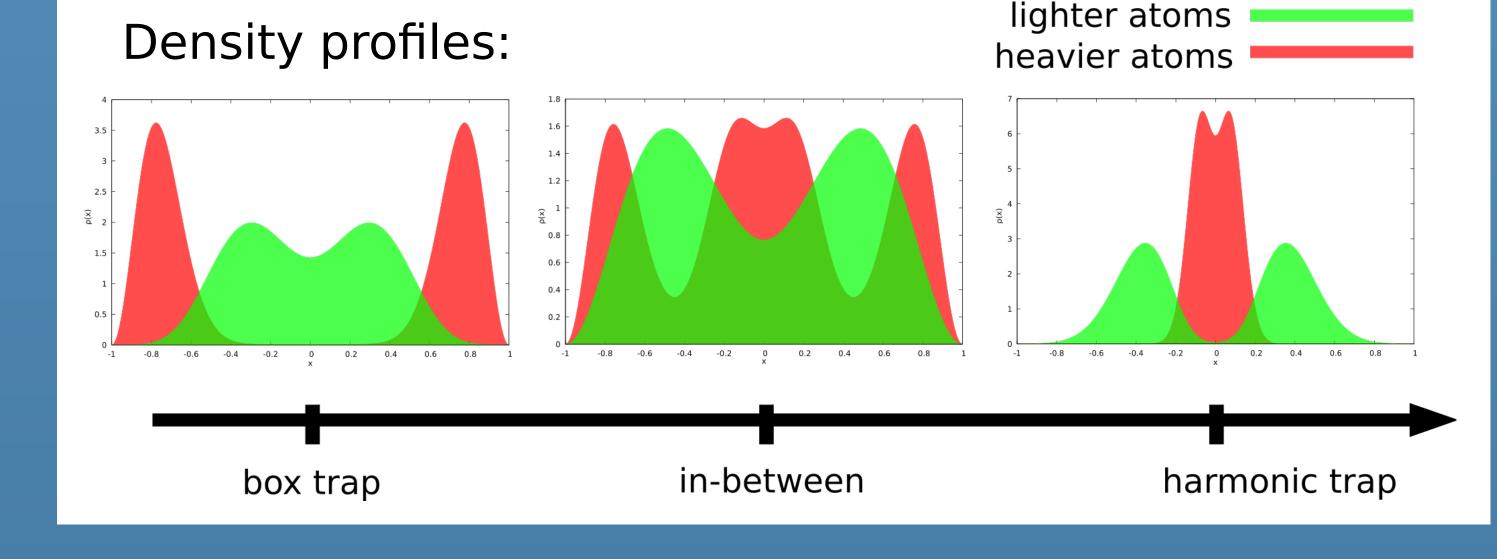
Abstract:

In a mass-imbalanced mixture of a few ultracold fermionic atoms with strong repulsive interactions, a spatial arrangement of the components depends on the shape of the external confinement. When the mixture is initially prepared in a one-dimensional box trap and then the harmonic potential is slowly turned on, the system undergoes structural transition. Finite-time quench through this

transition is analyzed, using an exact diagonalization method.

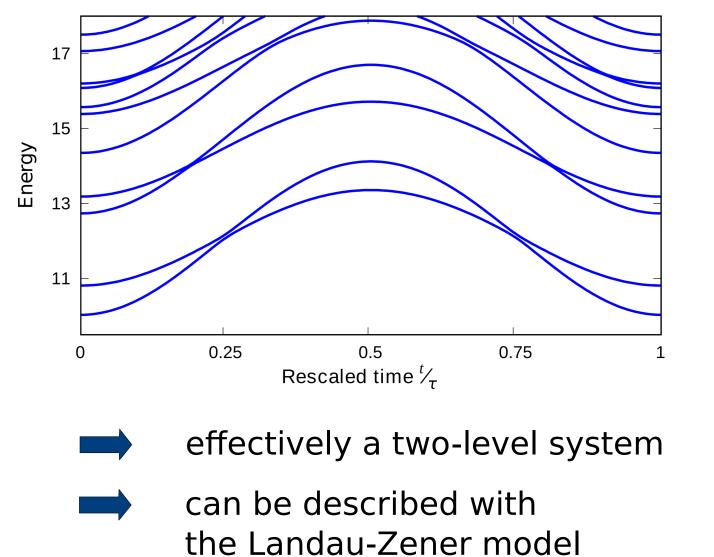
Analyzed transition

- strong repulsive interactions induce the separation of the components
- in a box trap, lighter atoms are localized in the center
- in a harmonic trap, heavier atoms are localized in the center
- transition between the traps is analogical to quantum phase transition

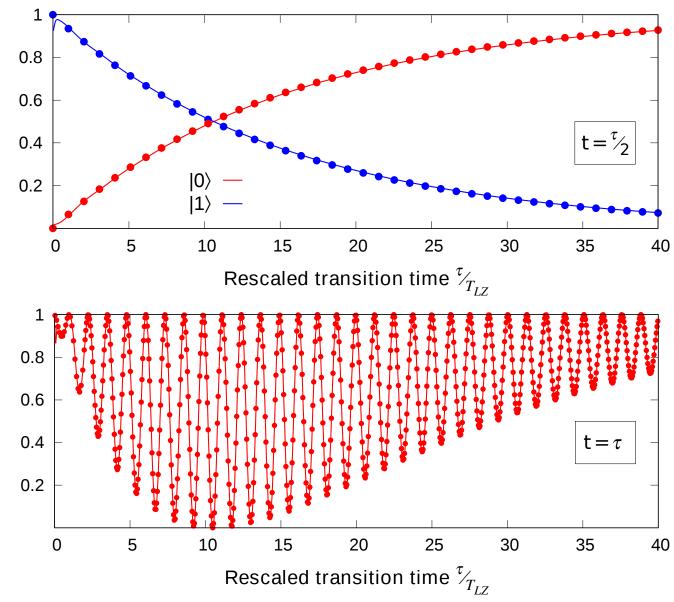


Results: 1+3 mixture

Energy spectrum:



Transition probability:



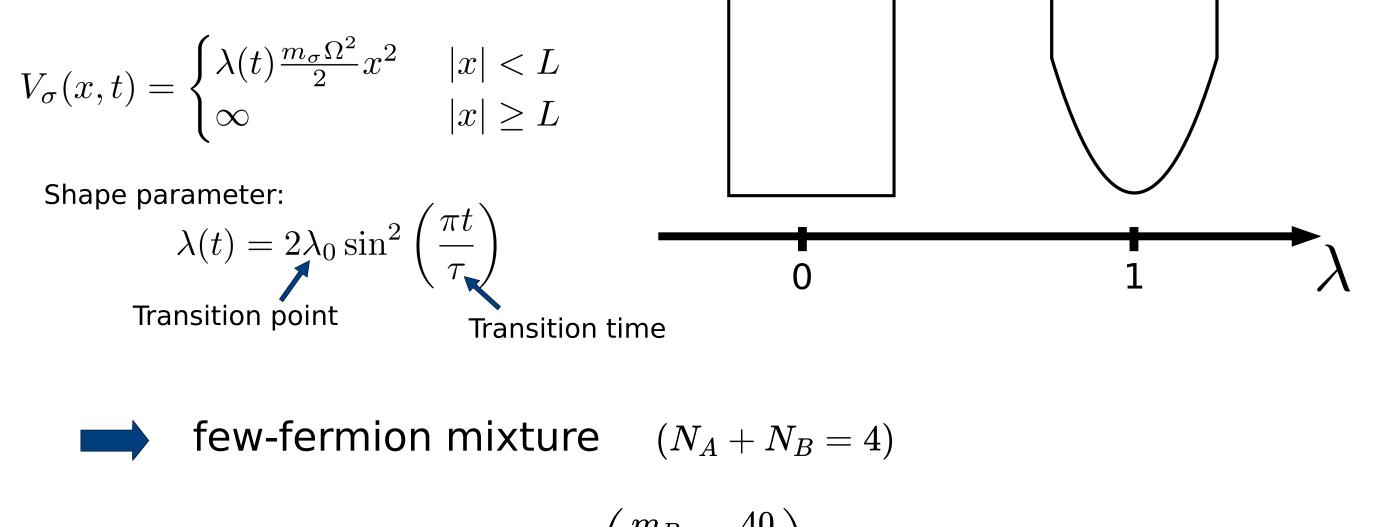
first transition: $\mathcal{P}_1(au/2) = \exp\left(-lpha au
ight)$

second transition: $\mathcal{P}_0(\tau) = \left| \mathcal{P}_0(\tau/2) e^{2i\phi_{\tau/2}} + \mathcal{P}_1(\tau/2) \right|$

Hamiltonian of the system

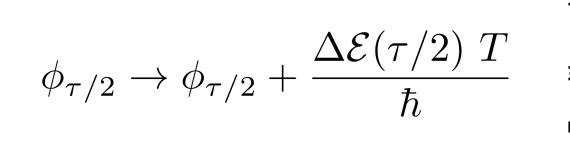
$$\hat{\mathcal{H}} = \sum_{\sigma \in \{A,B\}} \int \mathrm{d}x \,\hat{\Psi}_{\sigma}^{\dagger}(x) \left[-\frac{\hbar^2}{2m_{\sigma}} \frac{\mathrm{d}^2}{\mathrm{d}x^2} + V_{\sigma}(x) \right] \hat{\Psi}_{\sigma}(x) + g \int \mathrm{d}x \,\hat{\Psi}_{A}^{\dagger}(x) \hat{\Psi}_{B}^{\dagger}(x) \hat{\Psi}_{B}(x) \hat{\Psi}_{A}(x)$$

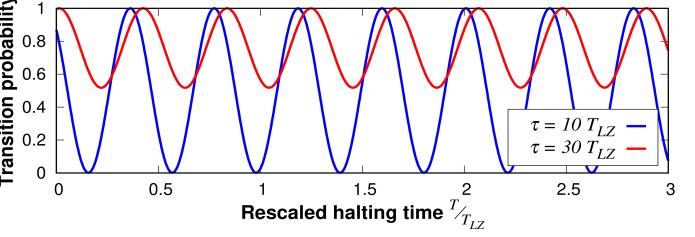
Confining potential:



 $\Rightarrow \text{ mass-imbalanced } \left(\frac{m_B}{m_A} = \frac{40}{6}\right)$ $\Rightarrow \text{ strong interactions } \left(g = 20\frac{\hbar^2}{m_A L}\right)$

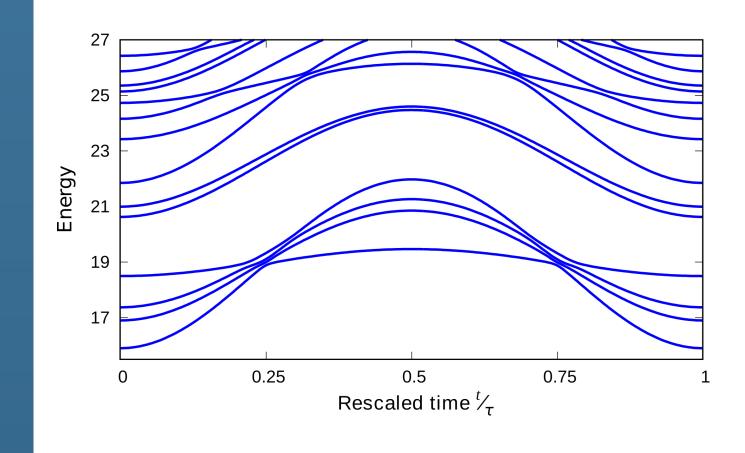
Halting protocol:





Results: 2+2 mixture

Energy spectrum:



- effectively a four-level system
- for long transition times effectively a two-level system

Transition probability: $\begin{array}{c} f = \frac{1}{2} \\ f = \frac{1}{2} \\$

Numerical method

The exact diagonalization of the many-body Hamiltonian:

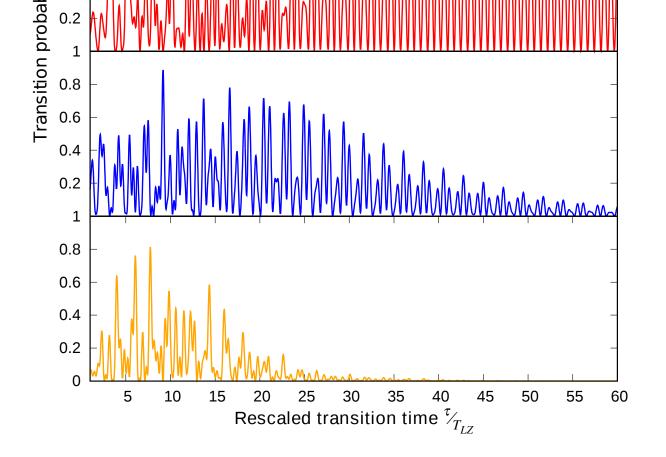
Fock basis is constructed from eigenstates of the non-interacting system

$$\hat{\Psi}_{\sigma}(x) = \sum_{n}^{M} \phi_{n\sigma}(x) \hat{a}_{n\sigma}$$

The Hamiltonian is expressed as a matrix in this basis

Required eigenstates are obtained (Lanczos method)

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References:

D. Pęcak and T. Sowiński, Phys. Rev, A 94, 042118 (2016)
D. Włodzyński, D. Pęcak and T. Sowiński, Phys. Rev, A 101, 023604 (2020)
D. Włodzyński and T. Sowiński, arXiv:2103.01798

