## Comment on "Exact Quantum Dynamics of a Bosonic Josephson Junction"

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Sakmann *et al.* challenge the two-mode Bose-Hubbard (TMBH) and Gross-Pitaevskii (GP) approximations for double-well Bose-Einstein condensates [1]. They find interesting deviations from the predictions of the TMBH model for dimensionless interaction parameter values as low as  $\Lambda \approx 1.5$ . Moreover, deviations from the GP approximation are obtained despite the validity of the 1D weak-interaction (1DWI) criterion.

In this comment we clarify that the value of  $\Lambda$  by itself is not necessarily relavant to the TMBH modelling validity question. The relevant dimensionless parameters of this problem are both  $\nu$  and  $\Lambda$  (see definitions below). In the specific setup of [1], if the interaction strength is increased, both  $\nu$  and  $\Lambda$  become larger, hance larger deviations are observed. Below we illuminate that  $\Lambda$  can be experimentally controlled *independently* of  $\nu$ . Hence we conclude that the relevance of  $\Lambda$  to the validity of the TMBH model (for a fixed value of  $\nu$ ) has not been established, and remains an interesting *open question*.

**Parameters and approximations.**— The experimental parameters of the 1D double-well system are the axial trap-frequency  $\omega$ , the barrier transmission coefficient T, and the atom number N. The 1D interaction strength is  $\lambda_0 = 2\hbar\omega_{\perp}a_s$ , where  $a_s$  is the s-wave scattering length, and  $\omega_{\perp}$  is the transverse trap frequency. The atoms mass is  $\mathbf{m}$ . These parameters define three characteristic length-scales: the axial trap size  $L = \sqrt{\hbar/m\omega}$ , the healing length  $l_c = \sqrt{\hbar/2m\lambda_0 n}$ , and the mean distance between atoms d = 1/n, where n = N/2L is the average atom density.

The conditions for TMBH validity [2–5] and for 1DWI [6] are  $l_c \gg L$  and  $l_c \gg d$  respectively:

$$\nu \equiv (L/l_c)^2 = \lambda_0 n/\hbar\omega \ll 1,$$
 [TMBH] (1)

$$\gamma \equiv (d/l_c)^2 = \mathbf{m}\lambda_0/\hbar^2 n \ll 1, \quad [1\text{DWI}] \quad (2)$$

If the TMBH approximation is valid, its effective parameters are the tunnel-splitting  $J \approx \hbar \omega \sqrt{T}$ , and the interaction-strength  $UN \approx \lambda_0 n$ . With these we have

$$\Lambda = NU/(2J), \qquad \nu = NU/(\hbar\omega), \qquad \gamma = \nu/N^2 \quad (3)$$

The interaction parameter  $\Lambda$  distinguishes between three *interaction regimes* [2, 3]: Rabi ( $\Lambda < 1$ ); Josephson ( $1 < \Lambda < N^2$ ); and Fock ( $\Lambda > N^2$ ). For  $\Lambda > 2$  preparations with all particles in one well exhibit self-trapping. We rewrite Eq.(1) and Eq.(2) as follows:

$$\nu \ll 1 \qquad \Longleftrightarrow \qquad \Lambda \ll [\hbar \omega / J] ,$$
 (4)

$$\gamma \ll 1 \qquad \Longleftrightarrow \qquad \Lambda \ll N^2[\hbar\omega/J] .$$
 (5)

Attainability of the Josephson regime.- It is now clear that the numerical results in [1] do not preclude the experimental realization of the strong-interaction regimes of the TMBH model. Ref. [1] only considers approaching these regimes via increasing  $\lambda \approx \lambda_0 N$ , keeping  $\omega$  and J fixed. Hence both  $\Lambda$  and  $\nu$  become larger. However,  $\Lambda$ can be increased without damaging Eq.(4), by either of the following strategies: (i) Decreasing J by increasing trap separation - resulting in a higher barrier between the harmonic traps while keeping fixed UN, with minor variation of the band-gap and the density in each trap; (ii) Increasing  $\omega$  by tightening the traps - increases n as  $\sqrt{\omega}$  as well as raises the barrier height, thus increasing  $\Lambda$ . Since *n* scales as  $\sqrt{\omega}$ , the TMBH small parameter  $\nu$  will decrease as  $1/\sqrt{\omega}$ , indicating improvement of TMBH validity. The  $\Lambda \gg 1$  TMBH regime can thus be realized [8].

The GP condition.- While  $\gamma \ll 1$  indicates phasecoherence *in each well* it does not guarantee its extension *across both wells*. Pushing this criterion ad-absurdum, it is clear that for N > 1, condition (4) automatically guarantees condition (5), leading to the false conclusion that the GP approximation is valid whenever the TMBH model is. Yet the TMBH model clearly has a quantum domain where GP fails [1–5, 7].

The classical GP limit of the TMBH model is only properly attained by taking the limit  $N \to \infty$  while maintaining fixed  $\Lambda$  [5, 7]. Quantum-classical correspondence is thus obtained over a timescale  $t_b$ , the quantum break time, which grows with increasing N. It might be argued that substantial deviations from GP are shown in Fig. 1d of [1] despite large N, but a closer inspection reveals they are typical: If all particles are initially prepared in one well and  $\Lambda \approx 2$ , then  $t_b$  only grows as  $\log(N)$  [5, 7], precisely as demonstrated in Fig. 1c,d.

- K. Sakmann, A.I. Streltsov, O.E. Alon, and L.S. Cederbaum, Phys. Rev. Lett. **103**, 220601 (2009).
- [2] A.J. Leggett, Rev. Mod. Phys. 73, 307 (2001).
- [3] R. Gati, M.K. Oberthaler, J. Phys. B 40, 61(R) (2007).
- [4] G.J. Milburn, J. Corney, E.M. Wright, and D.F. Walls, Phys. Rev. A 55, 4318 (1997).
- [5] J.R. Anglin, A. Vardi, Phys. Rev. A 64, 013605 (2001).
- [6] D.S. Petrov, G.V. Shlyapnikov, J.T.M. Walraven, Phys. Rev. Lett. 85, 3745 (2000).
- [7] A. Vardi and J.R. Anglin, Phys. Rev. Lett. 86, 568 (2001).
- [8] Note this procedure does not apply to 3D traps where the density scales as  $L^{-3}$  and the two mode limit is approached at large L. In this case, a variation of N will also be required.