



# Path Integrals Without Integrals

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# Path integral formalism (1)

- Amplitudes for transition from an initial state  $|\alpha\rangle$  to a final state  $|\beta\rangle$  in imaginary time  $T$  can be written as

$$A(\alpha, \beta; T) = \langle \beta | e^{-T\hat{H}} | \alpha \rangle$$

- Dividing the evolution into  $N$  time steps  $\epsilon = T/N$ , we get

$$A(\alpha, \beta; T) = \int dq_1 \cdots dq_{N-1} A(\alpha, q_1; \epsilon) \cdots A(q_{N-1}, \beta; \epsilon),$$

- Approximate calculation of short-time amplitudes leads to

$$A_N(\alpha, \beta; T) = \frac{1}{(2\pi\epsilon)^{MdN/2}} \int dq_1 \cdots dq_{N-1} e^{-S_N}$$

- Hagen Kleinert, *Path Integrals in Quantum Mechanics, Statistics, Polymer Physics, and Financial Markets*, 5<sup>th</sup> edition, World Scientific, Singapore, 2009.



## Path integral formalism (2)

- Continual amplitude  $A(\alpha, \beta; T)$  is obtained in the limit  $N \rightarrow \infty$  of the discretized amplitude  $A_N(\alpha, \beta; T)$ ,

$$A(\alpha, \beta; T) = \lim_{N \rightarrow \infty} A_N(\alpha, \beta; T)$$

- Discretized amplitude  $A_N$  is expressed as a multiple integral of the function  $e^{-S_N}$ , where  $S_N$  is called discretized action
- For a theory defined by the Lagrangian  $L = \frac{1}{2} \dot{q}^2 + V(q)$ , (naive) discretized action is given by

$$S_N = \sum_{n=0}^{N-1} \left( \frac{\delta_n^2}{2\epsilon} + \epsilon V(\bar{q}_n) \right),$$

where  $\delta_n = q_{n+1} - q_n$ ,  $\bar{q}_n = \frac{q_{n+1} + q_n}{2}$ .



# Discretized effective actions

- Discretized actions can be classified according to the speed of convergence of discretized path integrals
- Improved discretized actions have been earlier constructed, mainly tailored for calculation of partition functions
  - generalizations of the Trotter-Suzuki formula
  - improvements in the short-time propagation
  - expansion of the propagator by the number of derivatives
- Li-Broughton effective potential (1987)

$$V^{LB} = V + \frac{1}{24}\epsilon^2 (\nabla V)^2$$

in the left prescription gives  $1/N^4$  convergence for calculation of partition functions

- This cannot be extended to higher orders, nor such an approach was developed for general transition amplitudes



# Ideal discretization

- Ideal discretized action  $S^*$  is defined as the action giving exact continual amplitudes  $A_N = A$  for any discretization
- From the completeness relation

$$A(\alpha, \beta; T) = \int dq_1 \cdots dq_{N-1} A(\alpha, q_1; \epsilon) \cdots A(q_{N-1}, \beta; \epsilon),$$

it follows that the ideal short-time discretized action  $S_n^*$  is given by

$$A(q_n, q_{n+1}; \epsilon) = \frac{1}{(2\pi\epsilon)^{Md/2}} e^{-S_n^*}$$

where  $M$  is the number of particles,  $d$  dimensionality, and

$$S_n^* = \frac{\delta_n^2}{2\epsilon} + \epsilon W_n(\bar{q}_n, \delta_n; \epsilon),$$

and  $W$  is the (ideal) effective potential



# Improving effective actions (1)

- We start from Schrödinger's equation for the short-time amplitude  $A(q, q'; \epsilon)$

$$\left[ \frac{\partial}{\partial \epsilon} - \frac{1}{2} \sum_{i=1}^M \Delta_i + V(q) \right] A(q, q'; \epsilon) = 0$$
$$\left[ \frac{\partial}{\partial \epsilon} - \frac{1}{2} \sum_{i=1}^M \Delta'_i + V(q') \right] A(q, q'; \epsilon) = 0$$

- Here  $\Delta_i$  and  $\Delta'_i$  are  $d$ -dimensional Laplacians over initial and final coordinates of the particle  $i$ , while  $q$  and  $q'$  are  $d \times M$  dimensional vectors representing positions of all particles at the initial and final time.



## Improving effective actions (2)

- If we express short-time amplitude  $A(q, q'; \epsilon)$  by the ideal discretized effective potential  $W$

$$A(q, q'; \epsilon) = \frac{1}{(2\pi\epsilon)^{Md/2}} \exp \left[ -\frac{\delta^2}{2\epsilon} - \epsilon W \right]$$

we obtain equation for the effective potential in terms of  $x = \delta/2$ ,  $\bar{x} = (q + q')/2$ ,  $V_{\pm} = V(\bar{x} \pm x)$

$$W + x \cdot \partial W + \epsilon \frac{\partial W}{\partial \epsilon} - \frac{1}{8} \epsilon \bar{\partial}^2 W - \frac{1}{8} \epsilon \partial^2 W + \frac{1}{8} \epsilon^2 (\bar{\partial} W)^2 + \frac{1}{8} \epsilon^2 (\partial W)^2 = \frac{V_+ + V_-}{2}$$





# Recursive relations (1)

- The effective potential is given as a power series

$$W(x, \bar{x}; \epsilon) = \sum_{m=0}^{\infty} \sum_{k=0}^m W_{m,k}(x, \bar{x}) \epsilon^{m-k},$$

where systematics in  $\epsilon$ -expansion is ensured by  $\epsilon \propto x^2$ , and

$$W_{m,k}(x, \bar{x}) = x_{i_1} x_{i_2} \cdots x_{i_{2k}} c_{m,k}^{i_1, \dots, i_{2k}}(\bar{x})$$

- Coefficients  $W_{m,k}$  are obtained from recursive relations

$$\begin{aligned} 8(m+k+1)W_{m,k} &= \bar{\partial}^2 W_{m-1,k} + \partial^2 W_{m,k+1} \\ &- \sum_{l=0}^{m-2} \sum_r (\bar{\partial} W_{l,r}) \cdot (\bar{\partial} W_{m-l-2,k-r}) \\ &- \sum_{l=1}^{m-2} \sum_r (\partial W_{l,r}) \cdot (\partial W_{m-l-1,k-r+1}) \end{aligned}$$

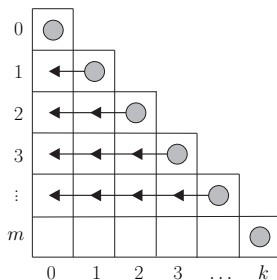


## Recursive relations (2)

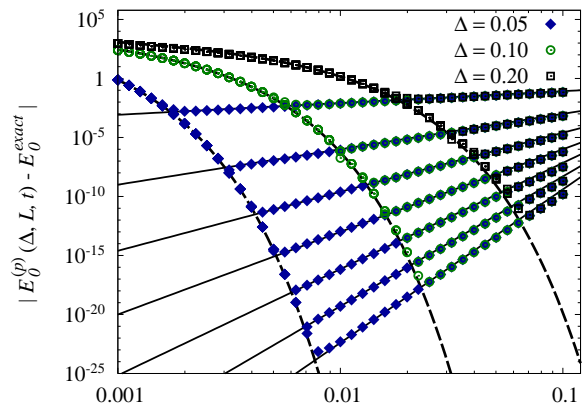
- Diagonal coefficients are easily obtained from recursive relations

$$W_{m,m} = \frac{1}{(2m+1)!} (x \cdot \bar{\partial})^{2m} V$$

- Off-diagonal coefficients are obtained by applying recursive relations in the following order

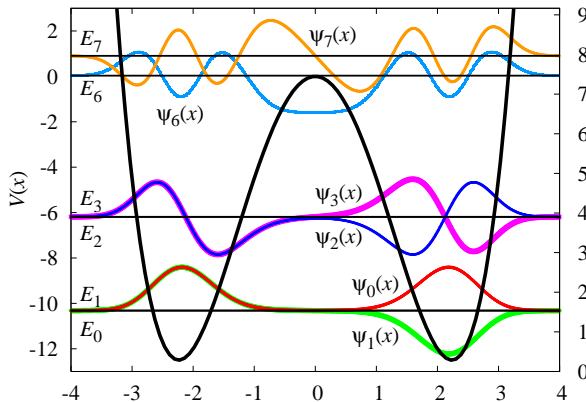


# Diagonalization of the evolution operator



$|E_0^{(p)}(\Delta, L, t) - E_0^{exact}|$  as a function of  $t$  calculated using level  $p = 1, 3, 5, 7, 9, 11, 13$  effective action for the quartic anharmonic potential, with  $m = \omega = 1, g = 48, \Delta = 0.05, L = 4$ .

# Energy eigenvalues and eigenstates



The double-well potential, its energy eigenvalues and eigenfunctions  $\psi_k(x)$  for  $k = 0, 1, 2, 3, 6, 7$ , with the parameters  $m = -10$ ,  $\omega = 1$ ,  $g = 12$ ,  $L = 10$ ,  $\Delta = 1.22 \cdot 10^{-3}$ ,  $t = 0.1$ .



## Effective actions: many-body $p=4$ result

$$\begin{aligned}
 S_N^{(p=4)} = & \sum \left\{ \epsilon \left( \frac{1}{2} \frac{\delta_i \delta_i}{\epsilon^2} + V \right) \right. \\
 & + \frac{\epsilon^2}{12} \partial_{k,k}^2 V + \frac{\epsilon \delta_i \delta_j}{24} \partial_{i,j}^2 V \\
 & - \frac{\epsilon^3}{24} \partial_i V \partial_i V + \frac{\epsilon^3}{240} \partial_{i,i,j,j}^4 V + \frac{\epsilon^2 \delta_i \delta_j}{480} \partial_{i,j,k,k}^4 V + \frac{\epsilon \delta_i \delta_j \delta_k \delta_l}{1920} \partial_{i,j,k,l}^4 V \\
 & + \frac{\epsilon^4}{6720} \partial_{i,i,j,j,k,k}^6 V - \frac{\epsilon^4}{120} \partial_i V \partial_{i,k,k}^3 V - \frac{\epsilon^4}{360} \partial_{i,j}^2 V \partial_{i,j}^2 V \\
 & - \frac{\epsilon^3 \delta_i \delta_j}{480} \partial_k V \partial_{k,i,j}^3 V + \frac{\epsilon^3 \delta_i \delta_j}{13440} \partial_{i,j,k,k,l,l}^6 V - \frac{\epsilon^3 \delta_i \delta_j}{1440} \partial_{i,k}^2 V \partial_{k,j}^2 V \\
 & \left. + \frac{\epsilon^2 \delta_i \delta_j \delta_k \delta_l}{53760} \partial_{i,j,k,l,m,m}^6 V + \frac{\epsilon \delta_i \delta_j \delta_k \delta_l \delta_m \delta_n}{322560} \partial_{i,j,k,l,m,n}^6 V \right\}
 \end{aligned}$$



## Effective actions: time-dependent formalism

$$W(\mathbf{x}, \bar{\mathbf{x}}; \varepsilon, \tau) = \sum_{m=0}^{\infty} \sum_{k=0}^m \left\{ W_{m,k}(\mathbf{x}, \bar{\mathbf{x}}; \tau) \varepsilon^{m-k} + W_{m+1/2,k}(\mathbf{x}, \bar{\mathbf{x}}; \tau) \varepsilon^{m-k} \right\},$$

$$\begin{aligned} \mathbf{R1} : 8(m+k+1) W_{m,k} &= 8 \frac{\Pi(m,k) (\bar{\mathbf{x}} \cdot \partial)^{2k} V^{(m-k)}}{(2k)! (m-k)! 2^{m-k}} + \bar{\partial}^2 W_{m,k+1} + \partial^2 W_{m-1,k} \\ &- \sum_{l,r} \left\{ \partial W_{l,r} \cdot \partial W_{m-l-2,k-r} + \partial W_{l+1/2,r} \cdot \partial W_{m-l-5/2,k-r-1} \right. \\ &\left. + \bar{\partial} W_{l,r} \cdot \bar{\partial} W_{m-l-1,k-r+1} + \bar{\partial} W_{l+1/2,r} \cdot \bar{\partial} W_{m-l-3/2,k-r} \right\}, \end{aligned}$$

$$\begin{aligned} \mathbf{R2} : 8(m+k+2) W_{m+1/2,k} &= 8 \frac{(1-\Pi(m,k)) (\bar{\mathbf{x}} \cdot \partial)^{2k+1} V^{(m-k)}}{(2k+1)! (m-k)! 2^{m-k}} + \bar{\partial}^2 W_{m+1/2,k+1} \\ &+ \partial^2 W_{m-1/2,k} - \sum_{l,r} \left\{ \partial W_{l,r} \cdot \partial W_{m-l-3/2,k-r} + \partial W_{l+1/2,r} \cdot \partial W_{m-l-2,k-r} \right. \\ &\left. + \bar{\partial} W_{l+1/2,r} \cdot \bar{\partial} W_{m-l-1,k-r+1} + \bar{\partial} W_{l,r} \cdot \bar{\partial} W_{m-l-1/2,k-r+1} \right\}. \end{aligned}$$



# Rotating ideal Bose gases

- Good approximation for weakly-interacting dilute gases
- Bose-Einstein condensates usually realized in harmonic magneto-optical traps
- Fast-rotating Bose-Einstein condensates extensively studied - one of the hallmarks of a superfluid is its response to rotation
- Paris group (J. Dalibard) has recently realized critically rotating BEC of  $3 \cdot 10^5$  atoms of  $^{87}\text{Rb}$  in an axially symmetric trap - we model this experiment
- The small quartic anharmonicity in  $x - y$  plane was used to keep the condensate trapped even at the critical rotation frequency [PRL **92**, 050403 (2004)]

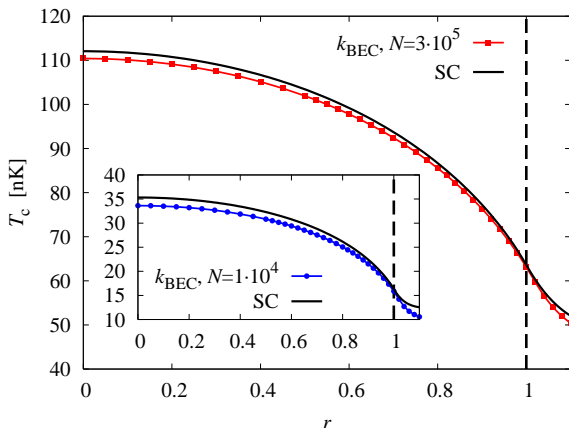


# Path integrals without integrals

- Using the large number of energy eigenvalues and eigenvectors of one-particle states, calculated by the exact diagonalization of the evolution operator, we study global and local properties of condensates
- $V_{BEC} = \frac{M}{2}(\omega_{\perp}^2 - \Omega^2)r_{\perp}^2 + \frac{M}{2}\omega_z^2 z^2 + \frac{k_{BEC}}{4}r_{\perp}^4$ ,  $\omega_{\perp} = 2\pi \times 64.8$  Hz,  $\omega_z = 2\pi \times 11.0$  Hz,  $k_{BEC} = 2.6 \times 10^{-11}$  Jm<sup>-4</sup>
- Typical values of the dimensionless inverse temperature  $\beta_{\text{eff}} = \hbar\omega_{\perp}/k_B T \lesssim 0.1$  represent already short (imaginary) times of propagation
- Hence, one-time-step (analytic) approximation to the calculation of BEC properties in the path integral formalism can be applied



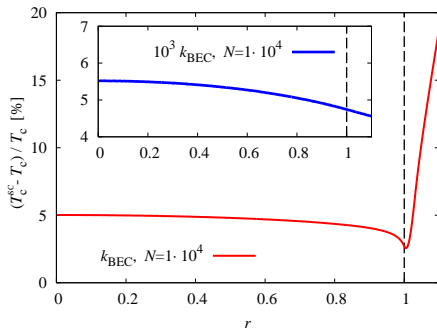
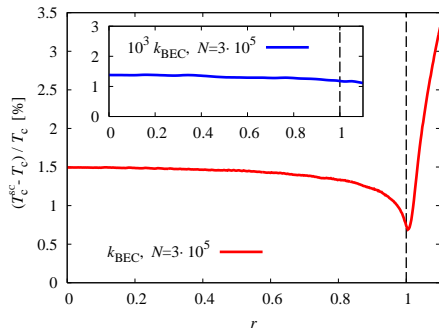
# Condensation temperature (1)



$T_c$  of a condensate in an anharmonic trap for different rotation frequencies  $r = \Omega/\omega_{\perp}$ , obtained with  $p = 21$  effective action. SC calculation: S. Kling and A. Pelster, PRA **76**, 023609 (2007).



## Condensation temperature (2)



Relative error of SC approximation for  $T_c$  of a condensate in an anharmonic trap for different rotation frequencies  $r = \Omega / \omega_{\perp}$ . Numerical results are obtained with  $p = 21$  effective action.



# Density profiles and time-of-flight graphs (1)

- Density profile is given in terms of the diagonal two-point propagator  $n(\mathbf{r}) = \rho(\mathbf{r}, \mathbf{r}) = \langle \hat{\Psi}^\dagger(\mathbf{r}) \hat{\Psi}(\mathbf{r}) \rangle$ , and for the ideal Bose gas

$$n(\mathbf{r}) = N_0 |\psi_0(\mathbf{r})|^2 + \sum_{n \geq 1} N_n |\psi_n(\mathbf{r})|^2$$

- In typical BEC experiments, a trapping potential is switched off and gas is allowed to expand freely during a short time of flight  $t$  (of the order of 10 ms)

$$n(\mathbf{r}, t) = N_0 |\psi_0(\mathbf{r}, t)|^2 + \sum_{n \geq 1} N_n |\psi_n(\mathbf{r}, t)|^2$$

where

$$\psi_n(\mathbf{r}, t) = \int \frac{d^3 \mathbf{k} d^3 \mathbf{R}}{(2\pi)^3} e^{-i\omega_{\mathbf{k}} t + i\mathbf{k} \cdot \mathbf{r} - i\mathbf{k} \cdot \mathbf{R}} \psi_n(\mathbf{R})$$



## Density profiles and time-of-flight graphs (2)

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Evolution of the  $x - y$  density profile of over-critically rotating ( $\Omega/\omega_{\perp} = 1.05$ ) condensate at  $T = 10$  nK  $< T_c = 55.3$  nK. The linear size of the profile is  $54 \mu\text{m}$ .



# Conclusions

- New method for numerical calculation of path integrals for a general non-relativistic many-body quantum theory
- In the numerical approach, discretized effective actions of level  $p$  provide substantial speedup of Monte Carlo algorithm from  $1/N$  to  $1/N^p$
- If the time of propagation/inverse temperature is small, analytic one-time-step approximation can be used: path integrals without integrals
- The derived results used to study properties of quantum systems by numerical diagonalization of the space-discretized evolution operator
- Numerical study of properties of (fast-rotating) ideal Bose-Einstein condensates
  - Condensation temperature and ground-state occupancy
  - Density profiles and time-of-flight graphs



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- Mathematica and PIMC codes: <http://www.scl.rs/speedup/>



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