

# Report of the Annual Meeting

## Numerical Results of the Complex Gross-Pitaevskii Equation

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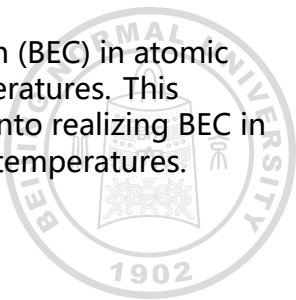
# Background

- Bose-Einstein Condensation (BEC) was predicted by Bose and Einstein in 1924-1925. The condensate was obtained experimentally for the first time in 1995 in a system consisting of about half a million alkali atoms cooled down to nanokelvin-level temperatures.



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- Achieving Bose-Einstein Condensation (BEC) in atomic systems requires extremely low temperatures. This challenge has prompted exploration into realizing BEC in atomless systems at relatively higher temperatures.

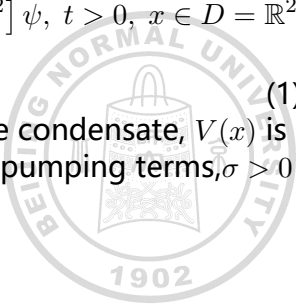


## Background

One possible candidate is a system of exciton-polaritons. Various mathematical models have been proposed for this new condensate. One of them, called complex GP equation, is explored there:

$$i\psi_t = -\Delta\psi + V(x)\psi + |\psi|^2\psi + i[\omega(x) - \sigma|\psi|^2]\psi, \quad t > 0, \quad x \in D = \mathbb{R}^2, \\ \psi(x, 0) = \psi_0(x), \quad x \in D. \quad (1)$$

Here,  $\psi = \psi(x, t)$  is the wave function of the condensate,  $V(x)$  is the trapping potential,  $\omega = \omega(x) \geq 0$  is the pumping terms,  $\sigma > 0$  is the decaying terms.



## Stationary radial solutions

We first explore the stationary solutions. Inserting  $\psi(x, t) = \exp(-i\mu t)\phi(x)$  into equation (1) leads to the following equation for  $\phi$ :

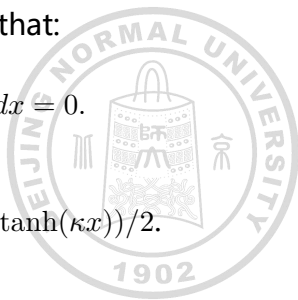
$$\mu\phi(x) = [-\Delta + V(x) + |\phi(x)|^2 + i(\omega(x) - \sigma|\phi(x)|^2)]\phi(x). \quad (2)$$

To make  $\mu$  to be real, it must be required that:

$$\int_D (\omega(x) - \sigma|\phi(x)|^2) |\phi(x)|^2 dx = 0.$$

We assume

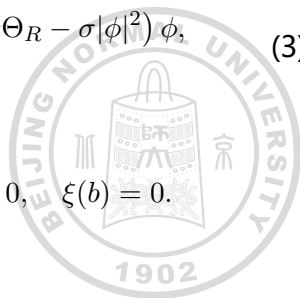
$$V(x) = |x|^2, \omega(x) = \alpha\Theta(R - |x|), \Theta(x) = (1 + \tanh(\kappa x))/2.$$



## Stationary radial solutions

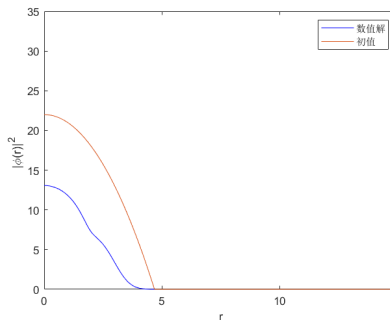
Now we solve the stationary radial solution of equation (2)(i.e.  $r = |x| \geq 0$ ). After some computation, we get the following equation:

$$\begin{aligned}\phi' &= \varphi, \\ \varphi' &= -\frac{1}{r}\varphi - \mu\phi + r^2\phi + |\phi|^2\phi + i(\alpha\Theta_R - \sigma|\phi|^2)\phi, \\ \xi' &= (\alpha\Theta_R - \sigma|\phi(r)|^2)|\phi(r)|^2r, \\ \mu' &= 0, \\ \phi(b) &= 0, \quad \varphi(0) = 0, \quad \text{Im}(\phi(0)) = 0, \quad \xi(b) = 0.\end{aligned}\tag{3}$$

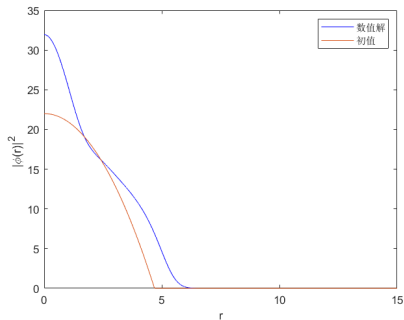


## Stationary radial solutions

We solve system (3) with high accuracy by using a collocation method. The MATLAB function *bvp5c* is used:



(a)  $R=2$



(b)  $R=8$

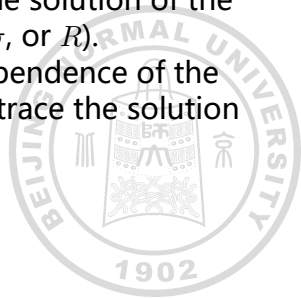
**Figure 1:** numerical solution of collocation method

# Numerical continuation

Let us rewrite system (3) as

$$F(u, \lambda) = 0, \quad (4)$$

where  $u \in \mathbb{B}$  ( $\mathbb{B}$  is a real Banach space) is the solution of the system,  $\lambda$  is a real parameter ( $\lambda$  can be  $\alpha$ ,  $\sigma$ , or  $R$ ). The idea in this section is to study the dependence of the solution,  $u(\lambda)$ , on the parameter,  $\lambda$ , i.e., to trace the solution branches  $[u(\lambda), \lambda]$  of (4).





## Numerical continuation

Suppose that  $(u_0, \lambda_0)$  is a solution of the discretized problem (4) and that the directional derivative  $\dot{u}_0 = du_0/d\lambda$  is known. Then the solution  $u_1$  at  $\lambda_1 = \lambda_0 + \Delta\lambda$  can be computed as

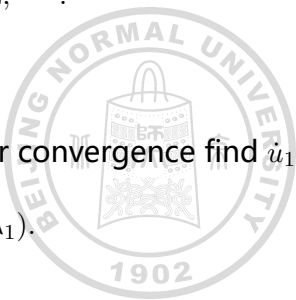
$$\begin{cases} F_u(u_1^i, \lambda_1)\Delta u_1^i = -F(u_1^i, \lambda_1), \\ u_1^{i+1} = u_1^i + \Delta u_1^i, \quad i = 0, 1, 2, \dots \end{cases}$$

with

$$u_1^0 = u_0 + \Delta\lambda \dot{u}_0,$$

where  $F_u(u, \lambda)$  is the Jacobian matrix. After convergence find  $\dot{u}_1$  from

$$F_u(u_1, \lambda_1)\dot{u}_1 = -F_\lambda(u_1, \lambda_1).$$



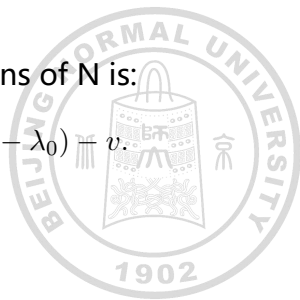
# Numerical continuation

To avoid folds, we can introduce a pseudo-arclength continuation method. This leads to the augmented system

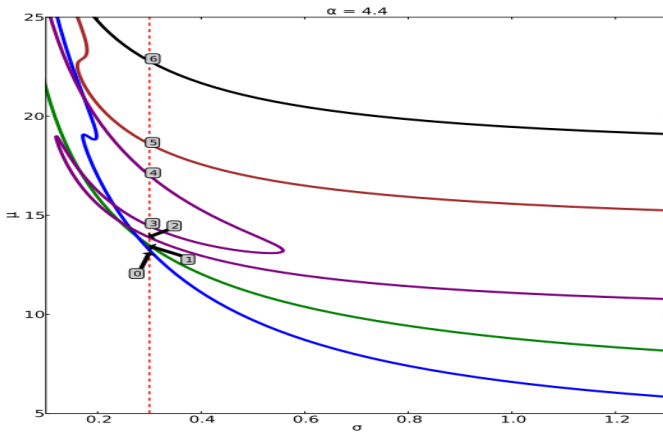
$$\begin{cases} F(u(v), \lambda(v)) = 0, \\ N(u(v), \lambda(v), v) = 0. \end{cases}$$

One of the most frequently used definitions of  $N$  is:

$$N(u, \lambda, v) \equiv \dot{u}_0^T(u - u_0) + \dot{\lambda}_0(\lambda - \lambda_0) - v.$$



# Numerical continuation



**Figure 2:** The chemical  $\mu$  of the system (3) for different values of  $\sigma$  with  $\alpha = 4.4$  and  $R = 2$  fixed

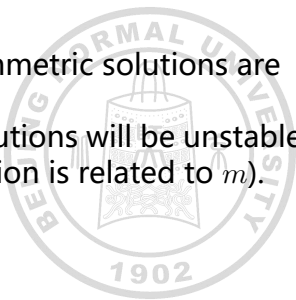
# Linear stability analysis

Consider the linear stability of the radially symmetric solutions  $\psi(r, t) = \exp(-i\mu t)\phi(r)$  under the small perturbations:

$$\psi(r, \theta, t) = \exp(-i\mu t)[\phi(r) + \varepsilon(u(r) \exp(-i(m\theta + \omega t)) + v^*(r) \exp(i(m\theta + \omega^* t)))],$$

with  $\varepsilon \ll 1, m = 1, 2, 3, \dots$ . The radially symmetric solutions are linearly unstable if  $\text{Im}(\omega) > 0$ .

The results in literature shows that, the solutions will be unstable when  $R \gtrsim 4.4$  or  $R \lesssim 0.6$  (the specific situation is related to  $m$ ).



# Strange-splitting Fourier spectral method

We now investigate the dynamics of equation (1) by Strange-splitting Fourier spectral method. The situation of one dimension can be stated as:

$$\begin{aligned} i\psi_t &= -\psi_{xx} + V(x)\psi + |\psi|^2\psi + i(\alpha\Theta_R - \sigma|\psi|^2)\psi, \\ \psi(x, 0) &= \psi_0(x), \quad a \leq x \leq b, \\ \psi(a, t) &= \psi(b, t), \quad \psi_x(a, t) = \psi_x(b, t), \quad t > 0. \end{aligned} \tag{5}$$

Consider  $h = \Delta x = (b - a)/M > 0$  as the mesh size, where  $M$  is an even positive integer and  $\tau = \Delta t > 0$  as the time step. Define  $\Psi_j^n$  as the numerical approximation of  $\psi(x_j, t_n)$ .

# Strange-splitting Fourier spectral method

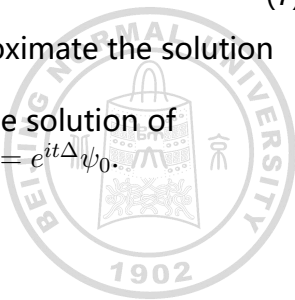
We now separate equation (5) into:

$$i\psi_t = V(x)\psi + |\psi|^2\psi + i(\alpha\Theta_R - \sigma|\psi|^2)\psi, \quad (6)$$

$$i\psi_t = -\psi_{xx}. \quad (7)$$

We combine equation (6) and (7) to approximate the solution of (5) on  $[t_n, t_{n+1}]$ .

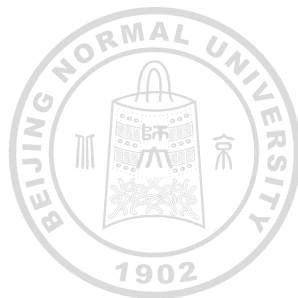
Equation (6) can be solved directly, and the solution of equation (7) can be represented as  $\psi(x, t) = e^{it\Delta}\psi_0$ .



# Strange-splitting Fourier spectral method

- Step1: From  $t_n$  to  $t_n + \tau/2$ , solve equation (6):

$$\Psi_j^{(1)} = \begin{cases} \Psi_j^n U_j^{(1)} e^{i\theta_j^{(1)}}, & \Theta_j > 0, \\ \Psi_j^n W_j^{(1)} e^{i\phi_j^{(1)}}, & \Theta_j = 0. \end{cases}$$



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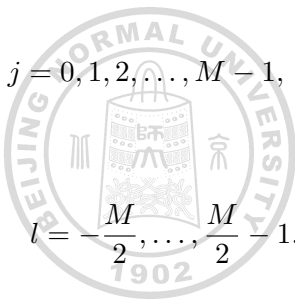
$$\Psi_j^{(1)} = \begin{cases} \Psi_j^n U_j^{(1)} e^{i\theta_j^{(1)}}, & \Theta_j > 0, \\ \Psi_j^n W_j^{(1)} e^{i\phi_j^{(1)}}, & \Theta_j = 0. \end{cases}$$

- Step2: From  $t_n$  to  $t_n + \tau$  solve equation (7), and use  $\Psi_j^{(1)}$  as initial condition:

$$\Psi_j^{(2)} = \frac{1}{M} \sum_{l=-M/2}^{M/2-1} e^{-i\omega_l^2 \tau} \hat{\Psi}_l^{(1)} e^{i\omega_l(x_j-a)}, \quad j=0,1,2,\dots,M-1,$$

where,

$$\hat{\Psi}_l^{(1)} = \sum_{j=0}^{M-1} \Psi_j^{(1)} e^{-i\omega_l(x_j-a)}, \quad \omega_l = \frac{2\pi l}{b-a}, \quad l = -\frac{M}{2}, \dots, \frac{M}{2} - 1.$$



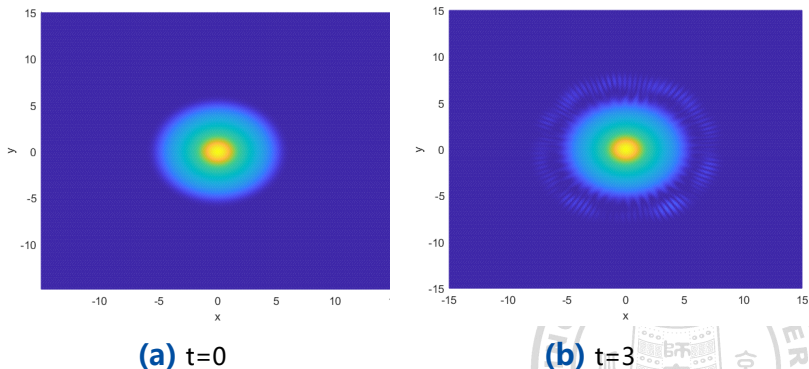


# Strange-splitting Fourier spectral method

- Step3: From  $t_n + \tau/2$  to  $t_n + \tau$  solve equation (6), and use  $\Psi_j^{(2)}$  as initial condition to get  $\Psi^{n+1}$ .

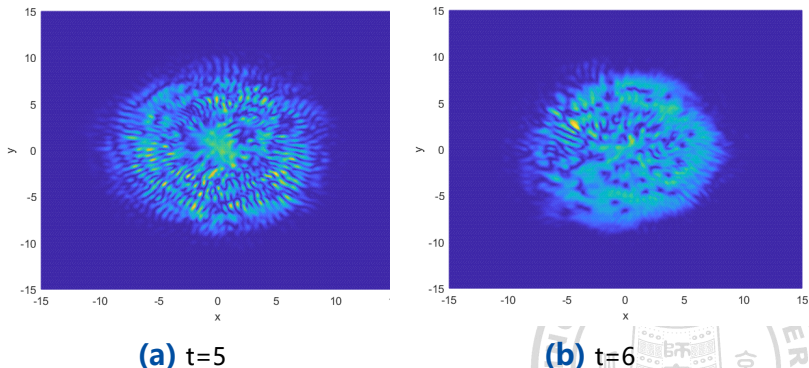


## Numerical results



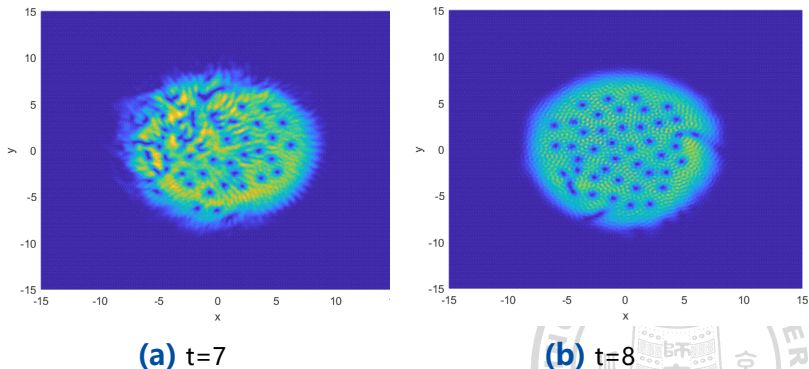
**Figure 3:** Simulation results for  $R = 8$ . The plots show the density distribution at different times. The dark blue areas represent regions where the density becomes zero.

## Numerical results



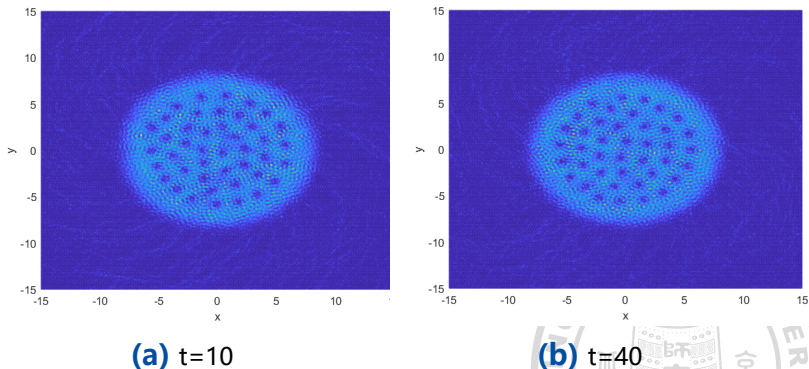
**Figure 4:** Simulation results for  $R = 8$ . The plots show the density distribution at different times. The dark blue areas represent regions where the density becomes zero.

## Numerical results



**Figure 5:** Simulation results for  $R = 8$ . The plots show the density distribution at different times. The dark blue areas represent regions where the density becomes zero.

## Numerical results



**Figure 6:** Simulation results for  $R = 8$ . The plots show the density distribution at different times. The dark blue areas represent regions where the density becomes zero.

