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1 Introduction

In this project, we consider the nonlinear Schrödinger equation (NLS) with focusing parameter ϵ ,

$$i\partial_t\psi = -\epsilon\Delta\psi + V(x)\psi - m(x)|\psi|^2\psi,$$

where $\psi(t, x) : \mathbb{R} \times \mathbb{R}^2 \rightarrow \mathbb{C}$ is the wavefunction, $V(x)$ is the trap potential, $m(x)$ is the mass and s-wave scattering length term, and $\epsilon > 0$. We would like to consider two kinds of trap potentials and domains in this problem.

First, we introduce nonlinear Schrödinger equations and the two kinds of trap potentials first.

1.1 Nonlinear Schrödinger Equations

The Schrödinger equation is a basic equation in quantum theory. The study of Schrödinger equations plays an important role in modern physics. In 1926, an Austrian physicist Erwin Schrödinger constructed the Schrödinger equation for explaining the active of particles. The famous Schrödinger equation is the form: $\mathcal{H}\psi = i\hbar\frac{\partial\psi}{\partial t}$, where \mathcal{H} is a Hamiltonian operator, ψ is the wavefunction, \hbar represents Planck's constant over 2π , i is the imaginary unit. The Schrödinger equation is an equation to describe the possible distribution of atoms. Solutions to Schrödinger's equation describe not only atomic and subatomic systems, electrons and atoms, but also macroscopic systems, possibly even the whole universe.

The nonlinear Schrödinger equation (NLS) is a nonlinear version of Schrödinger's equation in theoretical physics. In mathematical point of view, the NLS equation is a Schrödinger equation with nonlinear term. The nonlinear Schrödinger equation is the partial differential equation

$$i\partial_t\psi = -\frac{1}{2}\partial_x^2\psi + c|\psi|^2\psi$$

for the complex field ψ . There are many kinds of nonlinear Schrödinger equations. One general form of such equations [12] would be

$$i\partial_t\psi + \Delta\psi = f(\psi, \bar{\psi}).$$

These equations (particularly the cubic NLS equation) arise as model equations from several areas of physics: nonlinear optics, quantum condensates.

Here, we review some properties of NLS equations. Let's consider the equation

$$i\partial_t\psi + \Delta\psi = \pm|\psi|^2\psi.$$

The sign “ $-$ ” on the right hand side is focusing nonlinearity, and the sign “ $+$ ” is defocusing. When we add the potential part to the NLS equation, then it becomes

$$i\partial_t\psi + \Delta\psi = -m|\psi|^{p-1}\psi + V\psi, \tag{1}$$

where V is real and time independent. The equation (1) is sometimes referred to as a Gross-Pitaevskii equation. We can look at the dynamics of the Bose-Einstein condensate from the time-dependent Gross-Pitaevskii equation.

The behavior of ψ is determined by the potential V of the NLS equation. There is a special case of potential $V = \pm|x|^2$, this can be used to model a confining magnetic trap for Bose-Einstein condensation.

1.2 Trap Potential V

In the previous subsection, we have introduced some property of NLS equation. Now, we introduce the trap potential. Trap potential is to trap atoms on the minima potential. The following are two forms of trap potentials in this study, which are optical lattices and elliptic form.

1.2.1 Optical Lattices

An optical lattice is essentially an artificial crystal of light. A periodic intensity pattern that is formed by the interference of two or more laser beams. The shape of optical lattices looks like an egg carton in Fig. 1. It is called an optical lattice, since the periodic arrangement of trapping sites resembles a crystalline lattice. In Fig. 1, atoms are cooled and congregate in the minimal potential. The well depth and the periodicity are two important parts to affect the potential shape.

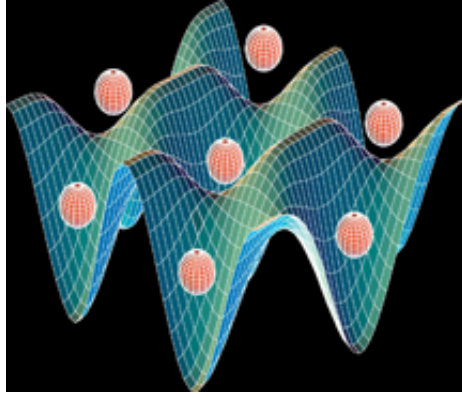


Figure 1: Optical Lattices.

Besides trapping cold atoms, optical lattices are also use in sorting microscopic particles [11] recently.

1.2.2 Elliptic Potential

Let $d \times d$ matrix M be symmetric, positive definite. Let $\mathbf{v} \in \mathbb{R}^d$ and $c \in \mathbb{R}$ be arbitrary. The elliptic potential [2] defines as

$$V(\mathbf{u}) = \frac{1}{2}\mathbf{u}^T M \mathbf{u} + \mathbf{u}^T \mathbf{v} + c,$$

where \mathbf{u} is a column vector in \mathbb{R}^d . Because M is symmetric and positive definite, $M^{\frac{1}{2}}$ exists. Therefore, we can also write $V(\mathbf{u}) = \frac{1}{2}\|M^{\frac{1}{2}}\mathbf{u}\|^2 + \mathbf{u}^T \mathbf{v} + c$. In \mathbb{R}^2 , the elliptic potential can be rewritten as

$$V(x, y) = ax^2 + bxy + cy^2 + dx + ey + f, \quad a, b, c, d, e, f \in \mathbb{R}$$

with $b^2 - 4ac < 0$. The shape of this potential looks like a bowl.

1.3 Goals

In this project, we consider nonlinear Schrödinger equations with focusing parameter ϵ ,

$$i\partial_t\psi = -\epsilon\Delta\psi + V(x)\psi - m(x)|\psi|^2\psi, \quad (2)$$

where $\psi(t, x) : \mathbb{R} \times \mathbb{R}^2 \rightarrow \mathbb{C}$ is the wavefunction, $V(x)$ is the trap potential, $m(x)$ is the mass and s-wave scattering length term, and $\epsilon > 0$. Such equation occurs in many physics, including nonlinear optics, quantum physics, and water waves.

In this project, we consider two kinds of different trap potentials and domains as follows:

Case 1. (ω case)

$$V(x) = (\omega_1 x_1)^2 + (\omega_2 x_2)^2,$$

where $\omega_1, \omega_2 \geq 0$ and $\omega_1, \omega_2 = 1, 2, \dots, 5$, and $m(x) = 1 + \frac{1}{2} \sin(2\pi x_1)$ with unit disk domain $\Omega = \{(x_1, x_2) | x_1^2 + x_2^2 < 1\}$.

Case 2. (μ case)

$$V(x) = 1 + \sin^2(\mu_1 x_1) + \sin^2(\mu_2 x_2),$$

where $\mu_1, \mu_2 = 1, 2, \dots, 5$ and $m(x) = 1 + \frac{1}{2} \cos(2\pi x_1)$ with square domain $\Omega = \{(x_1, x_2) | -1 < x_1 < 1, -1 < x_2 < 1\}$.

Here, we have two kinds of trap potentials and function $m(x)$ in Fig. 3 and Fig. 2, respectively. We focus on different situation in ω case and μ case. We concentrate on the well depth of the potential by varying the parameters ω_1 and ω_2 for the elliptic trap potential. And then we control the periodic for the optical lattices trap potential by changing the two parameters, μ_1 and μ_2 . For keeping the shapes of trap potentials, we use different types of domains for ω case and μ case.

The NLS equation has a special solution: $\psi(t, x) = e^{i\lambda t}\phi(x)$. The aim of this project is to find out the spectra of linearized operator which arises when the equation (2) is linearized around the special solution ψ . The goal is to study the stability of the solution by the spectrum of its linear operator.

There are many literature on NLS equation. The necessary condition for orbital stability and instability of single-spike bound state can be obtained in [9, 10].

Next, in Section 2 we will present the mathematical analysis about the NLS equation. And then in Section 3 we will discuss the numerical method. At last, some numerical results will be shown in Section 4.

2 Mathematical Analysis

2.1 Perturbation in NLS

Recall the nonlinear Schrödinger equation:

$$i\partial_t\psi = -\epsilon\Delta\psi + V(x)\psi - m(x)|\psi|^2\psi. \quad (3)$$

We consider the special solutions of NLS equation (3): $\psi(t, x) = e^{i\lambda t}\phi(x)$, where $\lambda = 0, 1$, and general case. $\phi(x)$ is a real-valued function independent of time. Substitution $\psi(t, x) = e^{i\lambda t}\phi(x)$ into (3) and the ϕ satisfy the nonlinear elliptic equation:

$$-\epsilon\Delta\phi + V\phi - m|\phi|^2\phi = -\lambda\phi,$$

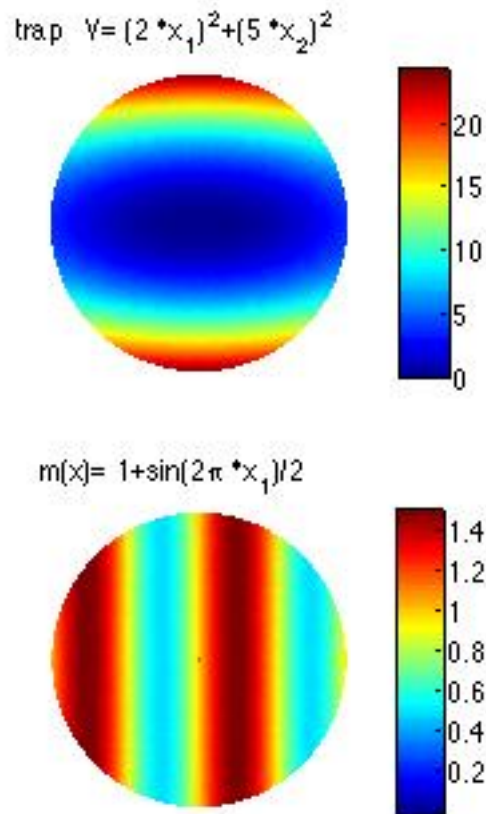


Figure 2: Elliptic potential and $m(x) = 1 + \frac{1}{2} \sin(2\pi x_1)$.

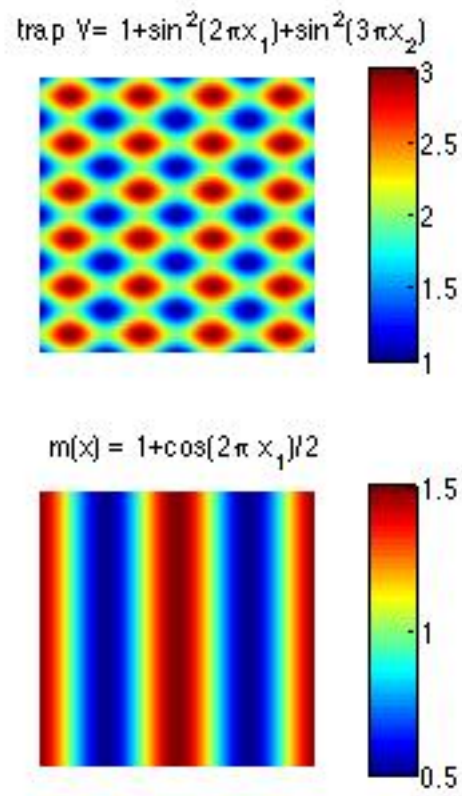


Figure 3: Optical lattices potential and $m(x) = 1 + \frac{1}{2} \cos(2\pi x_1)$.

that is

$$(-\epsilon\Delta + V - m|\phi|^2)\phi = -\lambda\phi \quad (4)$$

or

$$(-\epsilon\Delta + (V + \lambda) - m|\phi|^2)\phi = 0. \quad (5)$$

Equation (4) and (5) will be used in different solution form and algorithm.

To study the stability of the special solution form, we consider solutions of the NLS equation of the form

$$\psi(t, x) = e^{i\lambda t}(\phi(x) + h(t, x)). \quad (6)$$

The perturbation $h(t, x) \in \mathbb{R}$ satisfies an equation:

$$\partial_t h = \mathcal{L}h + (\text{nonlinear terms}),$$

where

$$\mathcal{L}h = -i\{(-\epsilon\Delta + (V + \lambda) - 3m\phi^2)h\}. \quad (7)$$

Remark. To explain how to obtain (7):

in equation (3), replace $\psi(x)$ by $e^{i\lambda t}(\psi(x) + h(t, x))$ then the left hand side of (3):

$$\begin{aligned} i\partial_t \psi &= i\partial_t(e^{i\lambda t}(\phi + h)) \\ &= i(i\lambda e^{i\lambda t}(\phi + h) + e^{i\lambda t}\partial_t h) \\ &= e^{i\lambda t}(-\lambda(\phi + h) + i\partial_t h). \end{aligned}$$

And the right hand side of (3):

$$\begin{aligned} -\epsilon\Delta\psi + V(x)\psi - m(x)|\psi|^2\psi &= e^{i\lambda t}(-\epsilon\Delta\phi - \epsilon\Delta h + V(x)\psi + h\phi - m|\phi + h|^2(\phi + h)) \\ &= e^{i\lambda t}\{-\epsilon\Delta\phi + V\phi - \epsilon\Delta h + h\phi \\ &\quad -m(|\phi|^2\phi + \phi^2(\bar{h} + h) + \phi h\bar{h} + \phi^2 h + \phi(h + \bar{h})h + h^2\bar{h})\} \\ &= e^{i\lambda t}(-\epsilon\Delta\phi + V\phi - m|\phi|^2\phi) \\ &\quad + e^{i\lambda t}(-\epsilon\Delta h + Vh - m|\phi|^2(h + \bar{h}) - m|\phi|^2 h) + e^{i\lambda t}O(h^2). \end{aligned}$$

Notice that $|\phi + h|^2 = (\phi + h)\overline{(\phi + h)} = |\phi|^2 + \phi(\bar{h} + h) + h\bar{h}$. Then

$$\begin{aligned} e^{i\lambda t}(-\lambda(\phi + h) + i\partial_t h) &= e^{i\lambda t}(-\epsilon\Delta\phi + V\phi - m|\phi|^2\phi) \\ &\quad + e^{i\lambda t}(-\epsilon\Delta h + Vh - m|\phi|^2(h + \bar{h}) - m|\phi|^2 h) + O(h^2). \end{aligned}$$

So

$$\begin{aligned} i\partial_t h &= (-\epsilon\Delta\phi + (V + \lambda)\phi - m|\phi|^2\phi) \\ &\quad + (-\epsilon\Delta h + (V + \lambda)h - m|\phi|^2(h + \bar{h}) - m|\phi|^2 h) + O(h^2) \\ &= -\epsilon\Delta h + (V + \lambda)h - m|\phi|^2(h + \bar{h}) - m|\phi|^2 h + O(h^2). \end{aligned}$$

If we consider $h(t, x)$ as a complex perturbation, ie. $h(t, x) \in \mathbb{C}$, then

$$\begin{aligned} \mathcal{L}h &= \mathcal{L}(\text{Re } h + i \text{Im } h) \\ &= -i\{(-\epsilon\Delta + (V + \lambda) - m\phi^2) \text{Re } h + i(-\epsilon\Delta + (V + \lambda) - m\phi^2) \text{Im } h - 2m\phi^2 \text{Re } h\} \\ &= (-\epsilon\Delta + (V + \lambda) - m\phi^2) \text{Im } h + i\{-(-\epsilon\Delta + (V + \lambda) - 3m\phi^2) \text{Re } h\}. \end{aligned}$$

And we rewrite \mathcal{L} as a matrix acting on $\begin{bmatrix} \operatorname{Re} h \\ \operatorname{Im} h \end{bmatrix}$,

$$\mathcal{L} = \begin{bmatrix} 0 & L_- \\ -L_+ & 0 \end{bmatrix}, \quad (8)$$

where

$$L_+ = -\epsilon\Delta + (V + \lambda) - 3m\phi^2, \quad L_- = -\epsilon\Delta + (V + \lambda) - m\phi^2. \quad (9)$$

L_+ and L_- are self-adjoint.

In the following Lemma, we will explain how we could rewrite the linearization term as (8). Are eigenvalues and eigenvectors of (8) and the original form the same or not?

Lemma 1. *Let $L = \begin{bmatrix} 0 & L_- \\ -L_+ & 0 \end{bmatrix}$ and $\mathcal{L}h = -\epsilon\Delta h + (V + \lambda)h - m|\phi|^2(h + \bar{h}) - m|\phi|^2h$, where L_+ and L_- are defined by (9). If ρ is an eigenvalue of L and $[u^T, v^T]^T$ be the eigenvector, then ρ is also an eigenvalue of \mathcal{L} and $u + iv$ is an eigenvector of L corresponding to ρ .*

Proof. Since ρ is an eigenvalue of L , that is,

$$L \begin{bmatrix} u \\ w \end{bmatrix} = \rho \begin{bmatrix} u \\ w \end{bmatrix}. \quad (10)$$

So

$$\begin{bmatrix} 0 & L_- \\ -L_+ & 0 \end{bmatrix} \begin{bmatrix} u \\ w \end{bmatrix} = \rho \begin{bmatrix} u \\ w \end{bmatrix}.$$

Then

$$-(-\epsilon\Delta + (V + \lambda) - 3m\phi^2)u = \rho w \quad (11)$$

and

$$(-\epsilon\Delta + (V + \lambda) - m\phi^2)w = \rho u. \quad (12)$$

(12) + $i \times$ (11), therefore

$$\begin{aligned} \rho(u + iw) &= -i(-\epsilon\Delta + (V + \lambda) - 3mq^2)u + (-\epsilon\Delta + (V + \lambda) - mq^2)w \\ &= -i\{(-\epsilon\Delta + (V + \lambda) - 3mq^2)u + i(-\epsilon\Delta + (V + \lambda) - mq^2)w\} \\ &= -i\{(-\epsilon\Delta(u + iw) + (V + \lambda)(u + iw) - mq^2(u + iw) - 2mq^2u)\}. \end{aligned}$$

Hence $u + iw$ is an eigenvector of \mathcal{L} corresponding to ρ . □

2.2 Stability

Our aim of this project is to study the stability of these solutions. We show the stability by solving eigenvalue problem. So in this subsection, we recall the definition of some stability, and the relation between eigenvalues and stability [4, 5].

At first, we give some notations and definitions. Consider an ODE

$$\frac{dx}{dt} = \dot{x} = f(x); \quad x \in \mathbb{R}^n, \quad (13)$$

where $f : \mathbb{R}^n \rightarrow \mathbb{R}^n$ and $x = [x_1, x_2, \dots, x_n]^T$. If $f(x^*) = 0$ for all t , the point x^* is called an equilibrium point.

Definition 1. x^* is a Lyapunov stable equilibrium if for every neighborhood U of x^* there is a neighborhood $V \subseteq U$ of x^* such that every solution $x(t)$ with $x(0) = x_0 \in V$ is defined and remains in U for all $t \geq 0$.

Definition 2. If V can be chosen above so that, in addition to the properties for stability, we have $\lim_{t \rightarrow \infty} x(t) = x^*$ then we say that x^* is asymptotically stable.

An equilibrium is called *neutrally stable* if it is Lyapunov stable but not asymptotically stable.

In studying the stability of x^* , we consider x^* plus a small perturbation $h(t)$, ie, $x(t) = x^* + h(t)$, where $|h(t)| \ll 1$. Substitute $x(t)$ into (13) and expand $f(x)$ by Taylor series: $\dot{x} + \dot{h} = f(x^* + h) = f(x^*) + Df(x^*)h + \mathcal{O}(|h|^2)$. The notation $Df(x^*)$ is the $n \times n$ Jacobian matrix of partial derivative of a vector-valued function f .

The eigenvalues and eigenvectors of the matrix $Df(x^*)$ determine the general solution. In studying stability we want to know whether the solution grows, stays constant, or decay to 0 as $t \rightarrow \infty$. It can be answered by evaluating the eigenvalues.

If λ is a real eigenvalue with eigenvector v , then there is a solution to the linearization equation of the form: $h(t) = cve^{\lambda t}$. If $\lambda = a \pm ib$ is a complex conjugate pair with eigenvectors $v = u \pm iw$ (where u, w are real), then $h_1(t) = e^{at}(u \cos bt - w \sin bt)$ and $h_2(t) = e^{at}(u \sin bt + w \cos bt)$ are two linearly independent solutions. In both cases, the real part of λ almost determines stability. Any solution of the linearized equation can be written as a linear combination of terms of these forms. We can obtain the following conclusions:

1. If all eigenvalues of $Df(x^*)$ have negative real parts, then $|h(t)| \rightarrow 0$ as $t \rightarrow \infty$ for all solutions.
2. If there exists one eigenvalue of $Df(x^*)$ has a positive real part, then there is a solution $h(t)$ with $|h(t)| \rightarrow +\infty$ as $t \rightarrow \infty$.
3. If some pair of complex-conjugate eigenvalues have zero real parts with distinct imaginary parts, then the corresponding solutions for $|h(t)|$ oscillate as $t \rightarrow \infty$ and neither decay nor grow as $t \rightarrow \infty$.

Moreover, if x^* is an equilibrium of $\dot{x} = f(x)$ and all the eigenvalues of the matrix $Df(x^*)$ have strictly negative real parts, then x^* is asymptotically stable. If at least one eigenvalue has strictly positive real part, then x^* is unstable.

We had discussed the solution of the ODE is an equilibrium. The statement of stability may be extend to non-constant orbits of ODE.

Here, let $\mathcal{O}_t(x) = x(t)$ and the initial value $x(0) = x_0$. Then the set $\mathcal{O}(x) = \{\mathcal{O}_t(x) : 0 \leq t\}$ is called orbit.

Definition 3. Let two orbits $\mathcal{O}(x)$ and $\mathcal{O}(\hat{x})$. If there is a reparameterization of time $\hat{t}(t)$ such that $|\mathcal{O}_t(\hat{x}) - \mathcal{O}_{\hat{t}(t)}(\hat{x})| < \epsilon$ for all $t \geq 0$, then we say two orbits $\mathcal{O}(x)$ and $\mathcal{O}(\hat{x})$ are ϵ -close.

Definition 4. An orbit $\mathcal{O}(x)$ is orbitally stable if for any $\epsilon > 0$, there is a neighborhood V of x so that, for all $\hat{x} \in V$, $\mathcal{O}(x)$ and $\mathcal{O}(\hat{x})$ are ϵ -close. If additionally V may be chosen so that, for all $\hat{x} \in V$, there exists a constant $\tau(\hat{x})$ so that $|\mathcal{O}_t(\hat{x}) - \mathcal{O}_{t-\tau(\hat{x})}(\hat{x})| < \epsilon$ as $t \rightarrow \infty$, then $\mathcal{O}_t(x)$ is asymptotically stable.

The linearization skill in general orbit is similar to the the pervious linearization method. Also discuss the eigenvalues of the linearization operator.

3 Numerical Method

In this section, we discretized the Laplace operator by finite-difference method first, and we will show it on two kinds of domain respectively. Then we present the numerical method for solving $\phi(x)$ from (5) first. And finally, we will describe a numerical method to compute the spectra of \mathcal{L} from (7).

Now, we recall our main question: to study the stability of the special solution form. For our solution form $\psi(t, x) = e^{i\lambda t}\phi(x)$, we will solving the time-inedpent term $\phi(x)$.

By (5), $\phi = \phi(x_1, x_2)$ satisfies

$$-\epsilon \left[\frac{\partial^2}{\partial x_1^2} \phi + \frac{\partial^2}{\partial x_2^2} \phi \right] + (V(x_1, x_2) + \lambda)\phi - m(x_1, x_2)\phi^3 = 0. \quad (14)$$

The natural boundary condition is

$$\lim_{|x| \rightarrow \infty} \phi(x_1, x_2) = 0. \quad (15)$$

Consider the same equation (5) in the unit disk domain $\Omega = \{(x_1, x_2) : x_1^2 + x_2^2 < 1\}$, applying the polar coordinate transformation,

$$x_1 = r \cos \theta, x_2 = r \sin \theta,$$

where $r = \sqrt{x_1^2 + x_2^2}$, $\theta = \tan^{-1}(x_2/x_1)$. We can rewrite (14), $\phi = \phi(r, \theta)$ in the polar coordinate system satisfies

$$-\epsilon \left[\frac{1}{r} \frac{\partial}{\partial r} \left(r \frac{\partial \phi}{\partial r} \right) + \frac{1}{r^2} \frac{\partial^2 \phi}{\partial \theta^2} \right] + (V(r, \theta) + \lambda)\phi - m(r, \theta)\phi^3 = 0, \quad (16)$$

with $0 < r < 1$, $0 \leq \theta < 2\pi$. And the boundary condition is

$$\lim_{r \rightarrow \infty} \phi(r, \theta) = 0. \quad (17)$$

Here, we have two equations and boundary conditions for solving ϕ in each domain.

3.1 Matrix Form of Laplace Operator

Now, we are going to describe the matrix form of the Laplace operator. In this study, we have two kinds of domains, unit disk domain and square domain. We will show them respectively.

3.1.1 Square Domain

The Laplace operator is a second order differential operator. Here, we consider the function ϕ in \mathbb{R}^2 . Then

$$\Delta\phi = \frac{\partial^2\phi}{\partial x^2} + \frac{\partial^2\phi}{\partial y^2}.$$

Here we use Finite-Difference method for the boundary value problem [1].

First step is to partition x into $-1 = x_0 < x_1 < x_2 < \dots < x_{N-1} < x_N = 1$, and partition y into $-1 = y_0 < y_1 < y_2 < \dots < y_{N-1} < y_N = 1$, where $x_i = -1 + i\Delta x$, $y_j = -1 + j\Delta y$ and $h = \Delta x = \Delta y = 2/N$. We use the Taylor series in the variable x, y about x_i, y_j to generate the centered-difference formula

$$\frac{\partial^2\phi}{\partial x^2} = \frac{f_{i-1,j} - 2f_{i,j} + f_{i+1,j}}{(\Delta x)^2} + O((\Delta x)^2),$$

$$\frac{\partial^2\phi}{\partial y^2} = \frac{f_{i-1,j} - 2f_{i,j} + f_{i+1,j}}{(\Delta y)^2} + O((\Delta y)^2)$$

for each $i = 1, 2, \dots, N-1$ and $j = 1, 2, \dots, N-1$, where $f_{i,j} = \phi(x_i, y_j)$.

In difference-equation form, this results in the finite-difference method, with local truncation error of order $O((\Delta x)^2 + (\Delta y)^2)$. Therefore

$$\begin{aligned} -\Delta\phi &\simeq - \left[\frac{f_{i-1,j} - 2f_{i,j} + f_{i+1,j}}{(\Delta x)^2} + \frac{f_{i,j-1} - 2f_{i,j} + f_{i,j+1}}{(\Delta y)^2} \right] \\ &= - \frac{f_{i-1,j} + f_{i+1,j} - 4f_{i,j} + f_{i,j-1} + f_{i,j+1}}{h^2}. \end{aligned}$$

We can rewrite it as a matrix representation:

$$\frac{1}{h^2} \begin{bmatrix} \hat{\mathbf{A}} & -\mathbf{I} & & 0 \\ -\mathbf{I} & \hat{\mathbf{A}} & -\mathbf{I} & \\ & \ddots & \ddots & \ddots \\ & & -\mathbf{I} & \hat{\mathbf{A}} & -\mathbf{I} \\ 0 & & & -\mathbf{I} & \hat{\mathbf{A}} \end{bmatrix} \begin{bmatrix} \mathbf{f}_1 \\ \mathbf{f}_2 \\ \vdots \\ \mathbf{f}_{N-2} \\ \mathbf{f}_{N-1} \end{bmatrix} \equiv \mathbf{A}\mathbf{F},$$

where

$$\hat{\mathbf{A}} = \begin{bmatrix} 4 & -1 & & 0 \\ -1 & 4 & -1 & \\ & \ddots & \ddots & \ddots \\ & & -1 & 4 & -1 \\ 0 & & & -1 & 4 \end{bmatrix}_{(N-1) \times (N-1)},$$

$\mathbf{f}_j = [f_{1,j}, f_{2,j}, \dots, f_{N-1,j}]^T, j = 1, 2, \dots, N-1$, and \mathbf{I} is an $(N-1)^2 \times (N-1)^2$ identity matrix.

3.1.2 Unit Disk Domain

Then we consider the equation (5) over a 2-dimensional disk. We use a discretization scheme [8] for equation (16). The Laplace operator is

$$\Delta\phi(r, \theta) = \Delta_r\phi + \Delta_\theta\phi = \frac{1}{r} \frac{\partial}{\partial r} \left(r \frac{\partial\phi}{\partial r} \right) + \frac{1}{r^2} \frac{\partial^2\phi}{\partial\theta^2}.$$

We discretize the Laplace operator in two parts. The first step is to discretize the operator $\Delta_r\phi$ on $r \in [0, 1]$, and then to discretize the operator $\Delta_\theta\phi$ on $\theta \in [0, 2\pi]$.

Firstly, we consider the operator $-\Delta_r\phi$ on $r \in [0, 1]$,

$$-\Delta_r\phi = -\frac{1}{r} \partial_r \left(r \frac{\partial\phi}{\partial r} \right).$$

To partition r into $0 = r_0 < r_{\frac{1}{2}} < r_1 < \dots < r_{N-\frac{1}{2}} < r_N = 1$, $r_j = j\Delta r$ and $\Delta r = \frac{1}{N}$, and to partition θ into $0 = \theta_0 < \theta_1 < \theta_2 < \dots < \theta_{M-1} < \theta_M = 2\pi$, $\theta_j = j\Delta\theta$ and $\Delta\theta = \frac{2\pi}{M}$, and denote $f_{i,j} = \phi(r_i, \theta_j)$. For each θ_j ,

$$-\Delta_r\phi \simeq \left[-\frac{1}{r_{i-\frac{1}{2}}} \frac{r_i(f_{i+\frac{1}{2},j} - f_{i-\frac{1}{2},j}) - r_{i-1}(f_{i-\frac{1}{2},j} - f_{i-\frac{3}{2},j})}{(\Delta r)^2} \right].$$

That is

$$\frac{1}{(\Delta r)^2} \begin{bmatrix} \frac{r_1}{r_{\frac{1}{2}}} & -\frac{r_1}{r_{\frac{1}{2}}} & & & & 0 \\ -\frac{r_1}{r_{\frac{3}{2}}} & \frac{1}{r_{\frac{3}{2}}}(r_1 + r_2) & -\frac{r_2}{r_{\frac{3}{2}}} & & & \\ & \ddots & \ddots & \ddots & & \\ & & -\frac{r_{N-2}}{r_{N-\frac{3}{2}}} & \frac{1}{r_{N-\frac{3}{2}}}(r_{N-2} + r_{N-1}) & -\frac{r_{N-1}}{r_{N-\frac{3}{2}}} & \\ 0 & & & -\frac{r_{N-1}}{r_{N-\frac{1}{2}}} & \frac{1}{r_{N-\frac{1}{2}}}(r_{N-1} + r_N) & \end{bmatrix} \begin{bmatrix} f_{\frac{1}{2},j} \\ f_{\frac{3}{2},j} \\ \vdots \\ f_{N-\frac{3}{2},j} \\ f_{N-\frac{1}{2},j} \end{bmatrix} \equiv \widehat{\mathbf{A}}\mathbf{f},$$

where

$$\widehat{\mathbf{A}} = \frac{1}{(\Delta r)^2} \begin{bmatrix} a_1 & c_1 & & & & 0 \\ b_1 & a_2 & c_2 & & & \\ & \ddots & \ddots & \ddots & & \\ & & & b_{N-2} & a_{N-1} & c_{N-1} \\ 0 & & & & b_{N-1} & a_N \end{bmatrix} \quad \text{and} \quad \mathbf{f}_j = \begin{bmatrix} f_{\frac{1}{2},j} \\ f_{\frac{3}{2},j} \\ \vdots \\ f_{N-\frac{1}{2},j} \end{bmatrix},$$

$$a_i = \frac{1}{r_{i-\frac{1}{2}}}(r_{i-1} + r_i) = \frac{i-1+i}{i-\frac{1}{2}} = 2, \quad i = 1, \dots, N$$

and

$$b_i = -\frac{r_i}{r_{i+\frac{1}{2}}} = -1 + \frac{1}{2i+1}, \quad c_i = -\frac{r_i}{r_{i-\frac{1}{2}}} = -1 - \frac{1}{2i-1}, \quad i = 1, \dots, N-1.$$

Now we consider the operator

$$-\Delta_\theta \phi = -\frac{1}{r^2} \frac{\partial^2 \phi}{\partial \theta^2}$$

on $\theta \in [0, 2\pi]$. Then for each $r_{i-\frac{1}{2}}$, $i = 1, 2, \dots, N$

$$-\Delta_\theta \phi \simeq \left[-\frac{1}{r_{i-\frac{1}{2}}^2} \frac{f(r_{i-\frac{1}{2}}, \theta_{j+1}) - 2f(r_{i-\frac{1}{2}}, \theta_j) + f(r_{i-\frac{1}{2}}, \theta_{j-1}))}{(\Delta\theta)^2} \right].$$

Let

$$\mathbf{B} = \text{diag} \left(\frac{2}{r_{\frac{1}{2}}^2 (\Delta\theta)^2}, \frac{2}{r_{\frac{3}{2}}^2 (\Delta\theta)^2}, \frac{2}{r_{\frac{5}{2}}^2 (\Delta\theta)^2}, \dots, \frac{2}{r_{N-\frac{1}{2}}^2 (\Delta\theta)^2} \right)$$

and

$$\mathbf{C} = \text{diag} \left(\frac{-1}{r_{\frac{1}{2}}^2 (\Delta\theta)^2}, \frac{-1}{r_{\frac{3}{2}}^2 (\Delta\theta)^2}, \frac{-1}{r_{\frac{5}{2}}^2 (\Delta\theta)^2}, \dots, \frac{-1}{r_{N-\frac{1}{2}}^2 (\Delta\theta)^2} \right),$$

$$\mathbf{f}_j = [f_{\frac{1}{2},j}, f_{\frac{3}{2},j}, \dots, f_{N-\frac{1}{2},j}]^\top \quad \text{for } j = 1, 2, \dots, M.$$

Then we can rewrite the Laplace operator as

$$\begin{bmatrix} \widehat{\mathbf{A}} + \mathbf{B} & \mathbf{C} & & & \mathbf{C} \\ \mathbf{C} & \widehat{\mathbf{A}} + \mathbf{B} & \mathbf{C} & & \\ & \ddots & \ddots & \ddots & \\ & & \mathbf{C} & \widehat{\mathbf{A}} + \mathbf{B} & \mathbf{C} \\ \mathbf{C} & & & \mathbf{C} & \widehat{\mathbf{A}} + \mathbf{B} \end{bmatrix} \begin{bmatrix} \mathbf{f}_1 \\ \mathbf{f}_2 \\ \vdots \\ \mathbf{f}_{M-1} \\ \mathbf{f}_M \end{bmatrix} \equiv \mathbf{A}\mathbf{F}.$$

Then the matrix \mathbf{A} represents the Laplace operator.

3.2 Algorithm

Now, we describe a numerical method to compute the spectrum of the linear operator \mathcal{L} defined by (7) for $\epsilon > 0$. There are two steps in our numerical method: First, we will solve the nonlinear problem (14) for ϕ . Second, compute the spectrum of the discretized linearized operator around ϕ .

Step I. Compute ϕ and it is energy minimizing among all solution of (14)–(17).

Step II. Compute the spectra of the discretized linearized operator \mathcal{L} . Since h is no longer real, so the linearized operator is a little different from \mathcal{L} in the equation (7).

Here, we denote some notation of the following. For $A \in \mathbb{R}^{M \times M}$, $q, m \in \mathbb{R}^M$, $m * q$ denote the Hadamard product of m and q , and $q^{\textcircled{p}}$ means the p -time Hadamard product of q . And $\text{diag}(q)$ represents the diagonal matrix of q .

In Step I. First, we represent $-\Delta q$ as

$$Aq = A[q_1, q_2, \dots, q_M]^T,$$

where q is an approximation of ϕ . As for the form of matrix A , it was shown in previous two subsections. Then the discretization of the equation (14) (and (16)) become

$$\epsilon Aq + (V + \lambda) * q - m * q^{\textcircled{3}} = 0,$$

where V and m are approximation of the function $V(x_1, x_2)$ and $m(x_1, x_2)$, respectively. We solve it by using an iteration method [7]:

$$[\epsilon A + \text{diag}(V + \lambda)]\tilde{q}_{new} = m * q_{old}^{\textcircled{3}}, \quad (18)$$

where \tilde{q}_{new} and q_{old} are unknown and known vector. The iteration step is shown in Algorithm 1 .

Algorithm 1 Iterative algorithm for solving $\phi(x)$

Step 1 Choose an initial guess of $q_{old} > 0$, and $\|q_{old}\|_2 = 1$.

Step 2 Solve (18), then obtain \tilde{q}_{new} .

Step 3 Let $\alpha_{new} = \frac{1}{\|\tilde{q}_{new}\|_2}$, $q_{new} = \alpha_{new}\tilde{q}_{new}$.

Step 4

if (converge) **then**

 Output the solution $(\alpha_{new})^{\frac{1}{2}}q_{new}$. Stop.

else

 Let $q_{old} = q_{new}$.

 Go to Step 2.

end if

When λ in a general case, λ is no longer a constant. λ varies from ϵ , so the numerical method has a little difference form. From (4), the discretization form becomes

$$[\epsilon A + \text{diag}(V - m * q^{\textcircled{2}})]q = -\lambda q. \quad (19)$$

It turns into an eigenvalue problem. In Algorithm 2, it shows the iteration step.

In Step II. Now we discretize \mathcal{L} of (8) into an eigenvalue problem.

$$L \begin{bmatrix} u \\ w \end{bmatrix} = \rho \begin{bmatrix} u \\ w \end{bmatrix}, \quad (20)$$

where

$$L = \begin{bmatrix} 0 & \epsilon A + \text{diag}(V + \lambda) - \text{diag}(m * q^2) \\ -\epsilon A - \text{diag}(V + \lambda) + \text{diag}(3m * q^2) & 0 \end{bmatrix}.$$

q and λ are obtain from **Step I**. We use ARPACK in MATLAB version R2007a to solve the linear algebraic eigenvalue problem and obtain eigenvalues ρ of L near origin.

4 Numerical Simulation

For each potential case, we summary the numerical results for three solution forms. Consider the solution form: $\psi(t, x) = e^{i\lambda t}\phi(x)$, $\lambda = 0, 1$, and general case.

Algorithm 2 Iterative algorithm for solving $\phi(x)$ in general solution

Step 1 Choose an initial guess of $q_{old} > 0$, and $\|q_{old}\|_2 = 1$.

Step 2 Solve

$$(\epsilon A + \text{diag}(V - m * q_{old}^{(2)}))\tilde{q}_{new} = -\lambda\tilde{q}_{new},$$

where $-\lambda$ is the smallest eigenvalue of $\epsilon A + \text{diag}(V - m * q_{old}^{(2)})$. Then obtain \tilde{q}_{new} .

Step 3 Let $\alpha_{new} = \frac{|\lambda|}{\|\tilde{q}_{new}\|_2}$, $q_{new} = \alpha_{new}\tilde{q}_{new}$.

Step 4

if (converge) **then**

 Output the solution q_{new} and λ . Stop.

else

 Let $q_{old} = q_{new}$.

 Go to Step 2.

end if

4.1 Setting

In ω case, the discretized matrix of Laplace operator has size NT by NT with $N = r/\Delta r$ and $T = 2\pi/\Delta\theta$. Here we use zero boundary condition and $r = 1$, $N = 32$, $T = 64$. And in μ case with square domain, the discretized matrix of Laplace operator has size N^2 by N^2 with $N = 2/h$ and here $N = 64$. And so the matrix size of the operator L are $2NT$ by $2NT$ and $2N^2$ by $2N^2$, respectively.

In our numerical method, we use finite-difference method to solve NLS equation. The finite-difference method has truncation error $O(h^2)$, where h is the grid size. In our experiment, $O(h^2) \approx c \cdot 10^{-3}$.

The region of ϵ that we consider about is larger than 10^{-5} , that is because the value is smaller than 10^{-5} we treat it as zero. We should control the region of ϵ to satisfy the boundary condition, which means the region of ϵ would not be large. In the region that we considered, the boundary mean values of all cases are less than 10^{-5} . In Fig. 4, we plot the mean value of the boundary of ϕ for ϵ from 0 to 0.003. We can see the mean of boundary will be less than 10^{-5} as ϵ smaller than 0.0024. So in that case, we only try $\epsilon \in [10^{-5}, 0.0024]$ in our numerical experiment.

Besides the boundary condition, we also take care of the shape of ϕ , here we narrow down the focus on the solitary solution for this part only. As ϵ goes larger, sometimes the shape would change. The following statement like ϵ^* are all in the region of ϵ .

In the following table(Table 1), we write down the region of ϵ for each case. In each potential case, for our convenient, we choose the intersection of ϵ region for different parameters.

In this project, we focus on the parameter ϵ changing, and to study when the solution of the NLS would be unstable or stable. The notation ϵ^* denote as the bifurcation of stable and unstable.

We find that a pair of purely imaginary eigenvalues will collide at the origin and split into a pair of real eigenvalues for $\epsilon < \epsilon^*$. That is, when ϵ less than ϵ^* , the spectrum of the linearized operator are all pure imaginary. As ϵ larger than ϵ^* , the spectrum of the linearized operator has at least one pair of eigenvalue with nonzero real part. That means

Table 1: The region of ϵ .

	case μ	case ω
$\lambda = 0$	$[10^{-5}, 0.016]$	$[10^{-5}, 0.0024]$
$\lambda = 1$	$[10^{-5}, 0.025]$	$[10^{-5}, 0.005]$
general λ	$[10^{-5}, 0.001]$	$[10^{-5}, 0.001]$

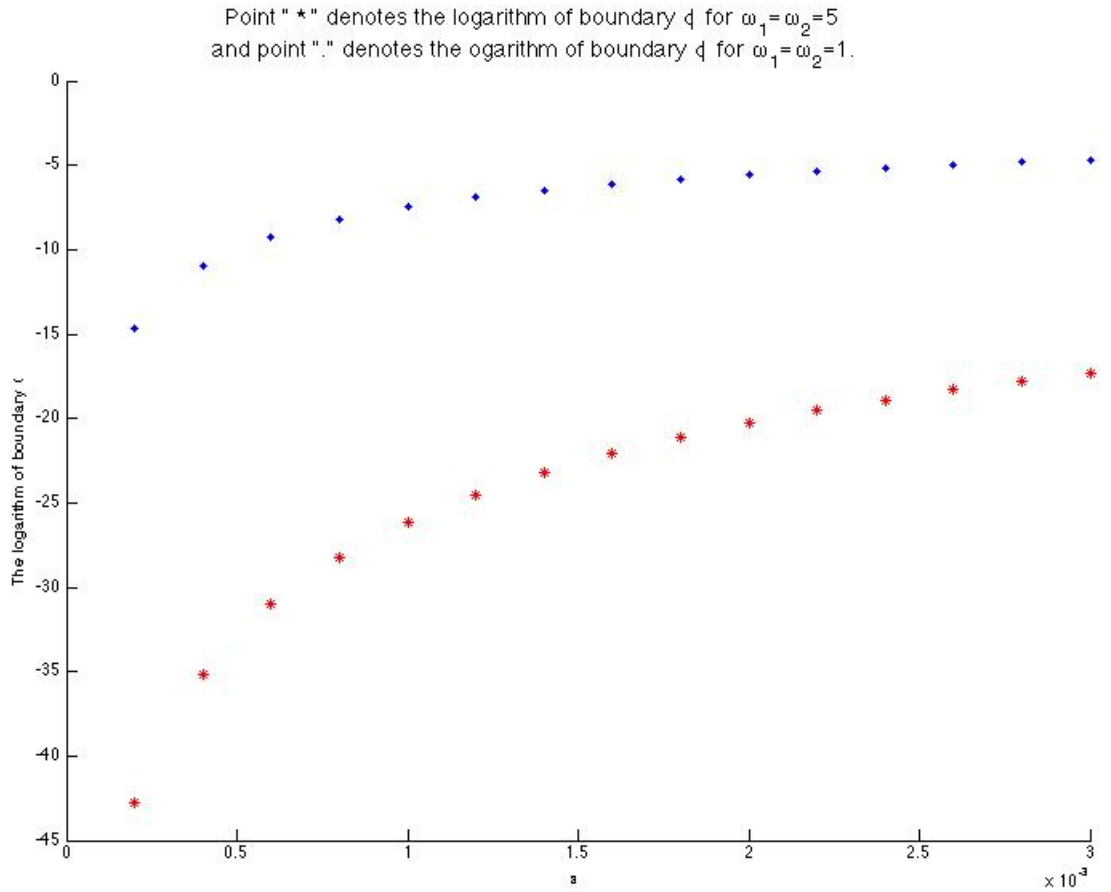


Figure 4: Boundary ϕ in ω case for $\lambda = 0$.

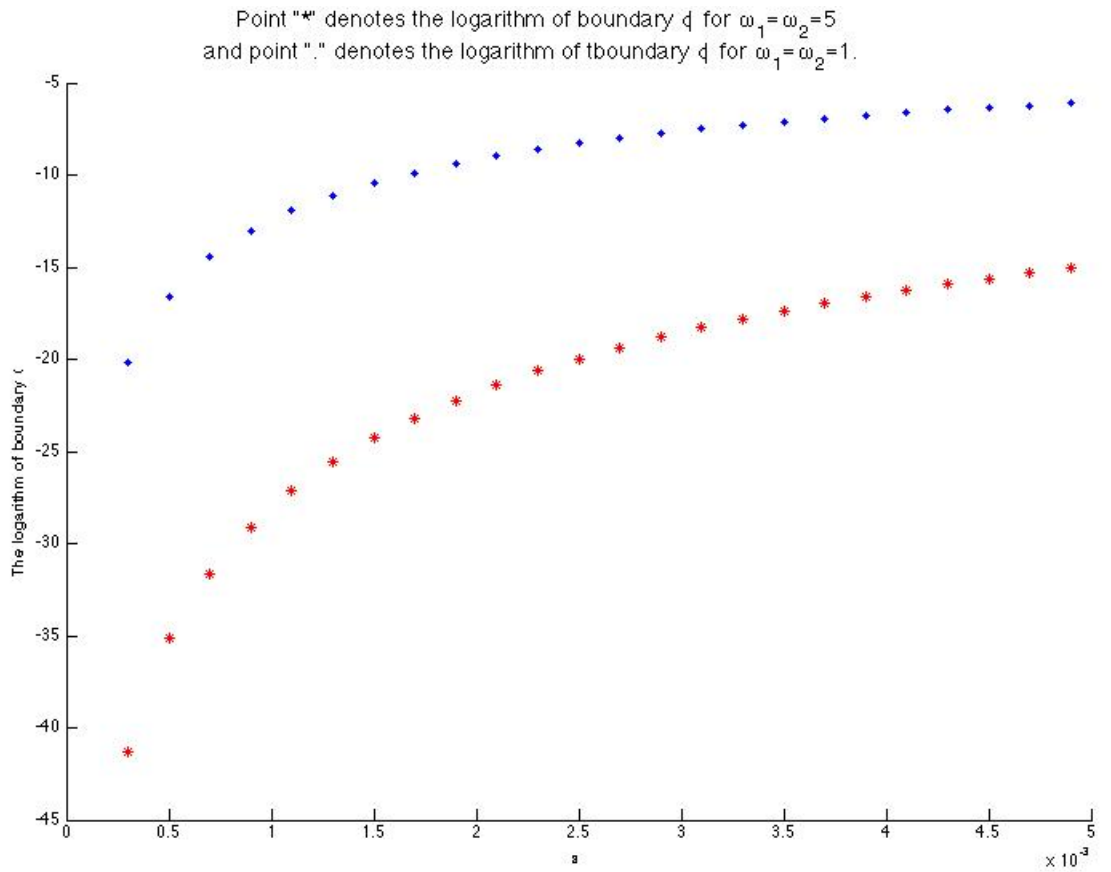


Figure 5: Boundary ϕ in ω case for $\lambda = 1$.

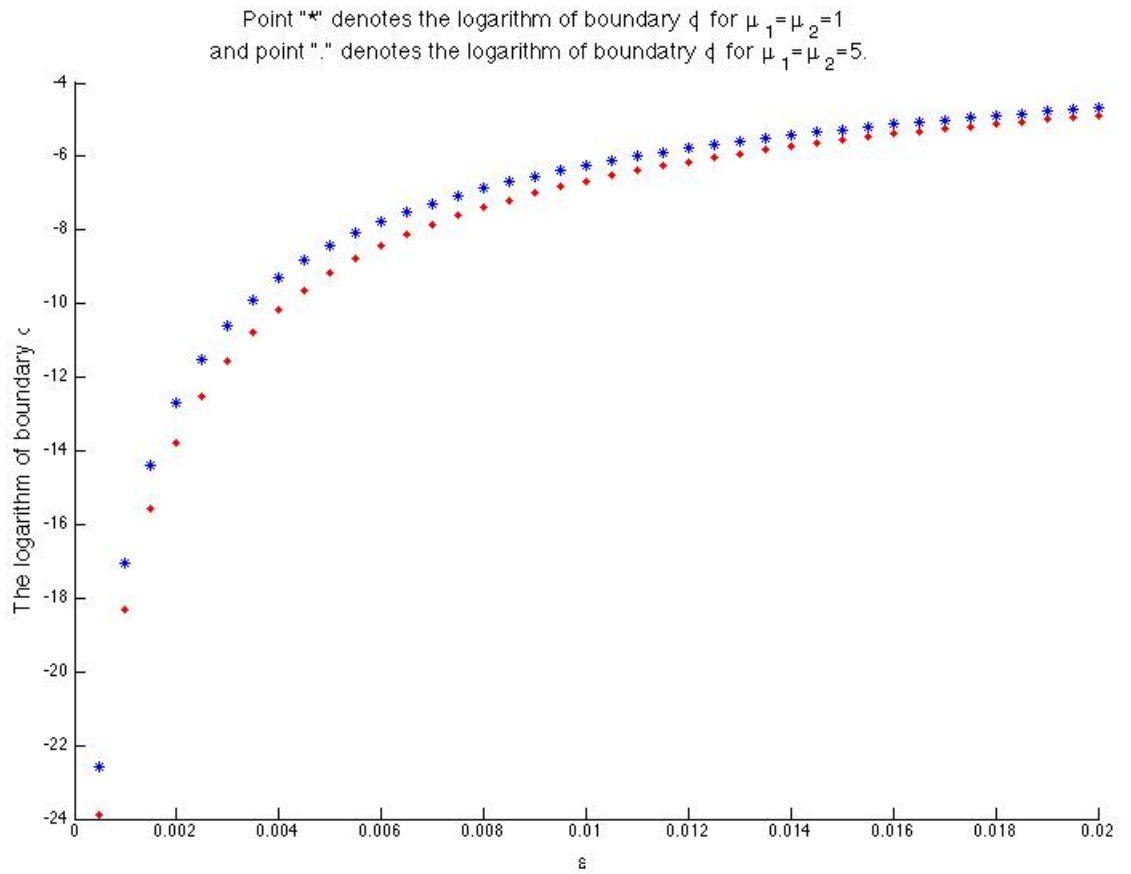


Figure 6: Boundary ϕ in μ case for $\lambda = 0$.

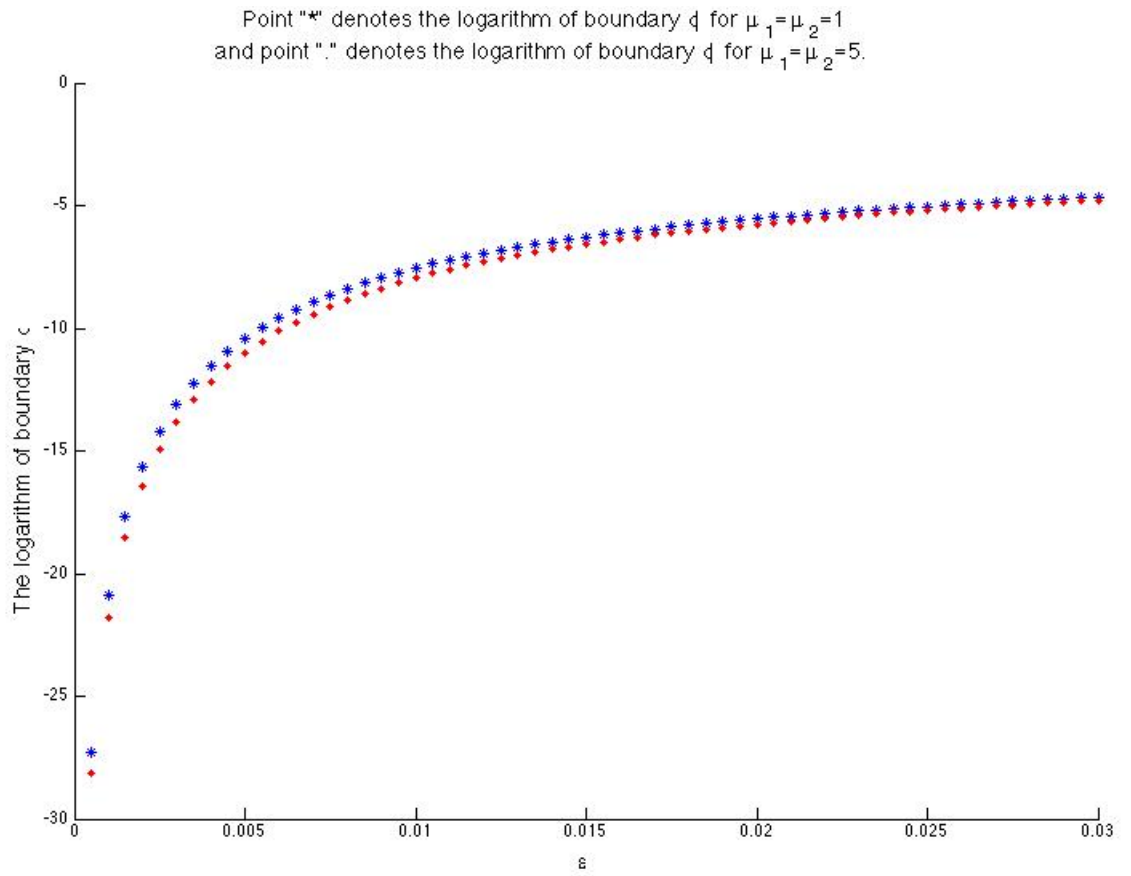


Figure 7: Boundary ϕ in μ case for $\lambda = 1$.

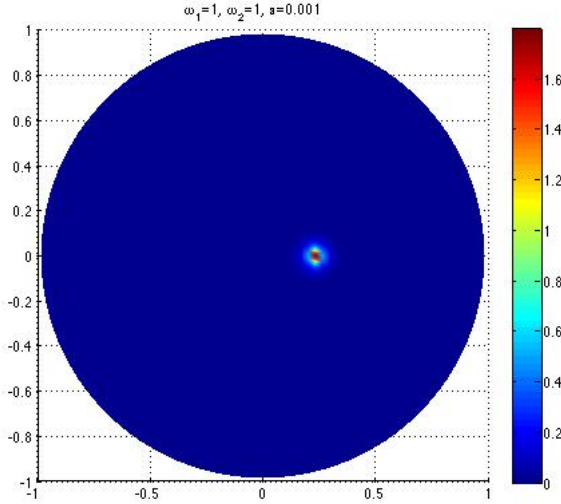


Figure 8: Solution of ϕ for $\omega_1 = 1, \omega_2 = 1, \epsilon = 0.001, \lambda = 1$.

the solution is stable while $\epsilon < \epsilon^*$, unstable when $\epsilon > \epsilon^*$.

Since the numerical results of eigenvalues are approximation of the exact mathematical values. There exist some error of the result. We determined that the eigenvalue has nonzero real part when the absolute value of the real part of eigenvalue is larger than 10^{-3} .

Since the regions of ϵ are small, and the spectrum of linearization operation changes so fast. We move ϵ in a small step for each case. $\delta\epsilon = 10^{-5}$. To prevent losing some pairs of pure complex eigenvalue from turning to real eigenvalues, we also move the target of the algorithm for finding eigenvalues.

This method is not an efficient method but an easy way to catch the changing of spectra.

4.2 ω case

In ω case, the maxima of ϕ is not at center of the disk; instead, it is on the right of the center. As ϵ go larger and the potential change, the maxima of ϕ seem to moving to the center of domain. In figure 10, we fix $\epsilon = 0.0002$ and changing the well depth of potential $V(x)$. We find that the location of maxima ϕ changes. It moves to the center of the domain.

We know trap potential $V(x)$ and $m(x)$ would effect the solution ϕ . The NLS equation in our model is an focusing case, that is the atoms would stay near the maxima of $m(x)$, ie, $x_1 = 1/4$.

1. $\lambda = 0$: Testing ϵ in the region that we mentioned before, and find that the eigenvalues of the linearized operation \mathcal{L} are all pure imaginary.
2. $\lambda = 1$: We find there is an ϵ^* in this solution case. The ϵ^* represent the bifurcation of pure imaginary to has a pair of eigenvalue with nonzero real part. In Fig. 11 shows the spectrum of \mathcal{L} . And in Fig. 12, it shows the ϵ^* of the parameters $\omega_1, \omega_2 = 1, 2, \dots, 5$

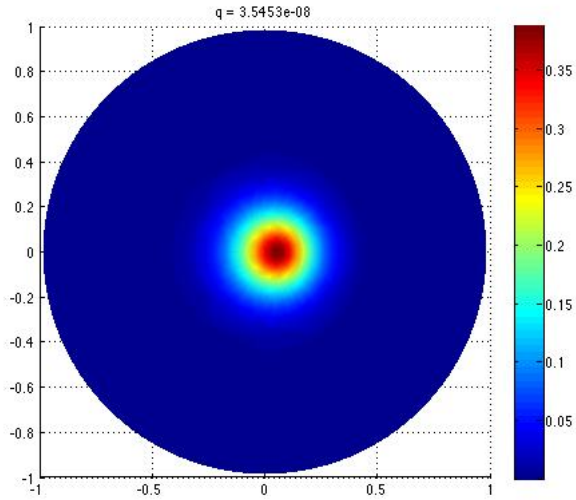


Figure 9: Solution of ϕ for $\omega_1 = 1, \omega_2 = 1, \epsilon = 0.001, \lambda = 0$.

3. In elliptic potential case, we can find ϵ^* in $e^{it}\phi(X)$ this solution form in our numerical method, but we can not find the ϵ^* in other solution form. At first we guess ϵ^* in $e^{i0t}\phi(x)$ form of solution can be found in smaller ϵ . But we do not find that in the region of ϵ as we mentioned before.
4. λ is general case: In this solution case, we can not find the ϵ^* . When ϵ is quite small, the value of λ is already negative. That means the trap potential is not positive at all, $V(x) < 0$ on the center of the disk. In Figure 13 it shows the value of λ as ϵ changes.

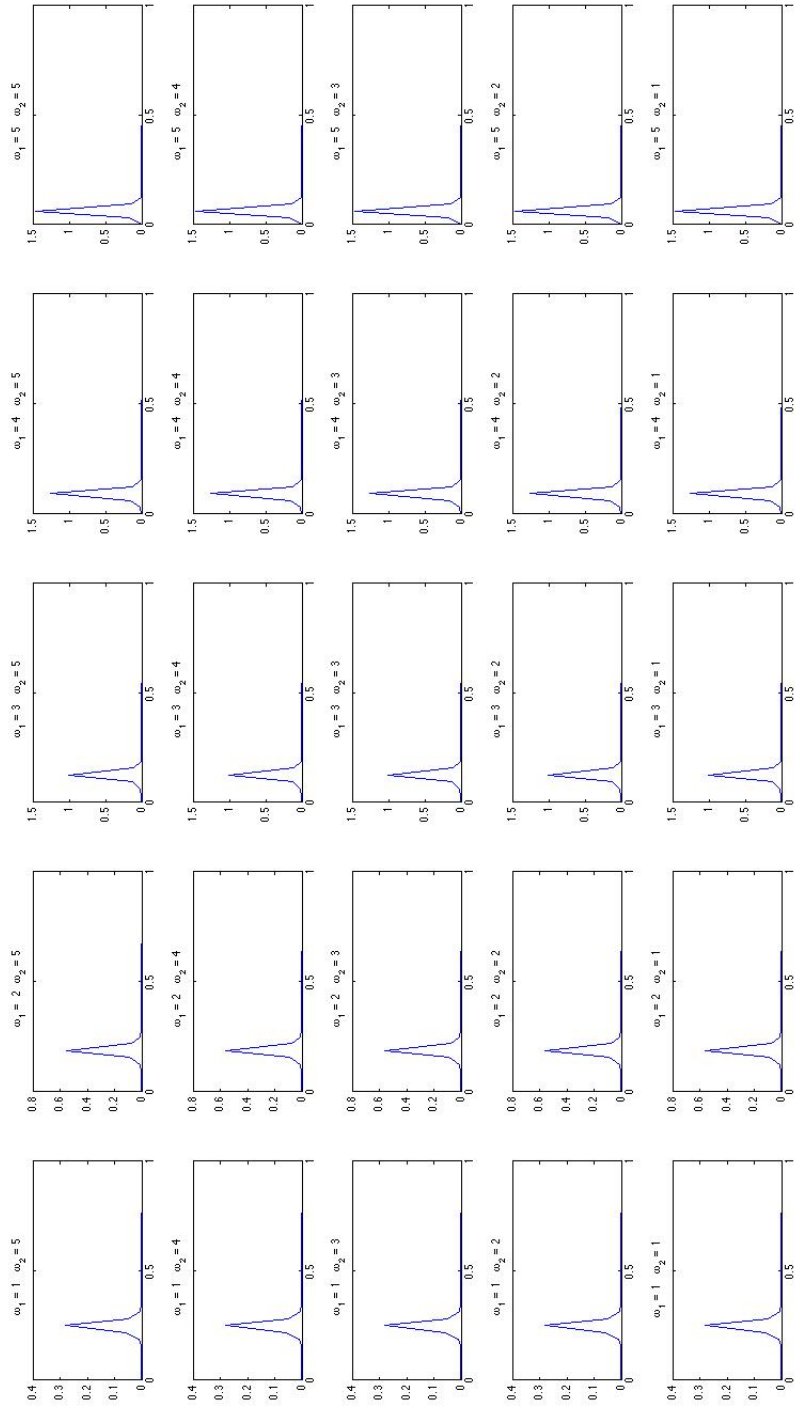


Figure 10: ϕ moves for $\lambda = 1$, $\epsilon = 0.0002$.

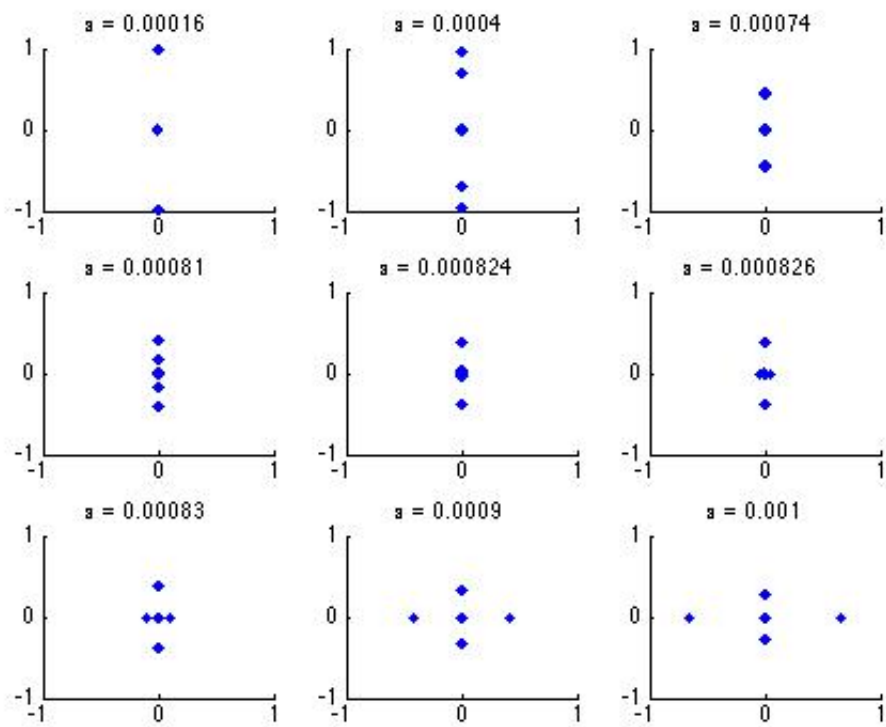


Figure 11: Spectrum of ω case for $\lambda = 1$.

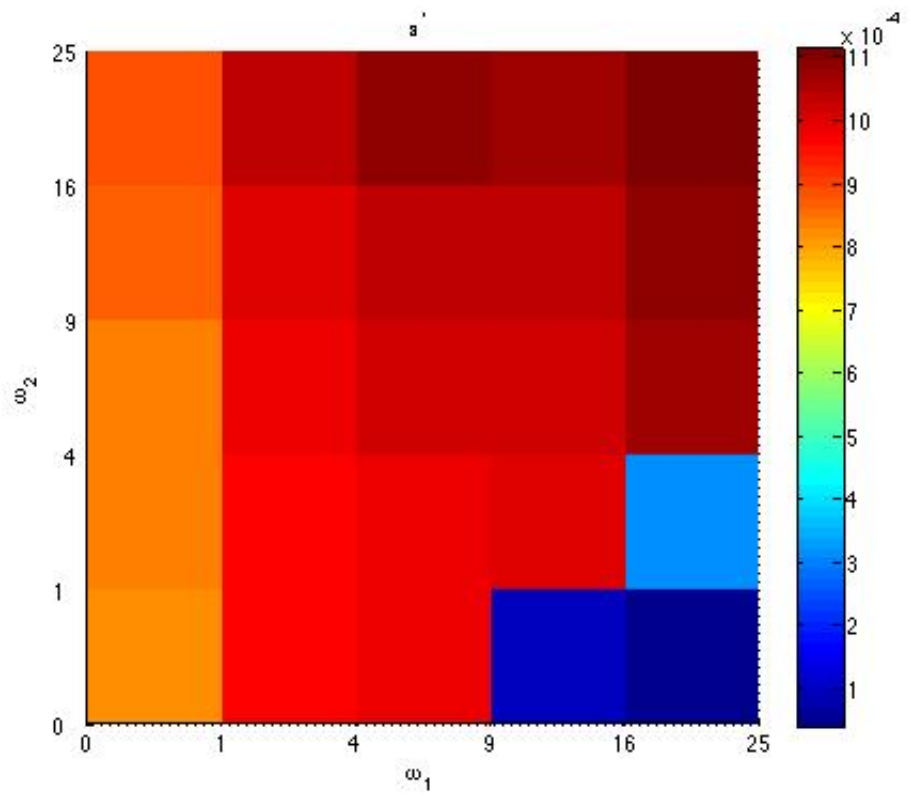


Figure 12: ϵ^* of ω case for $\lambda = 1$.

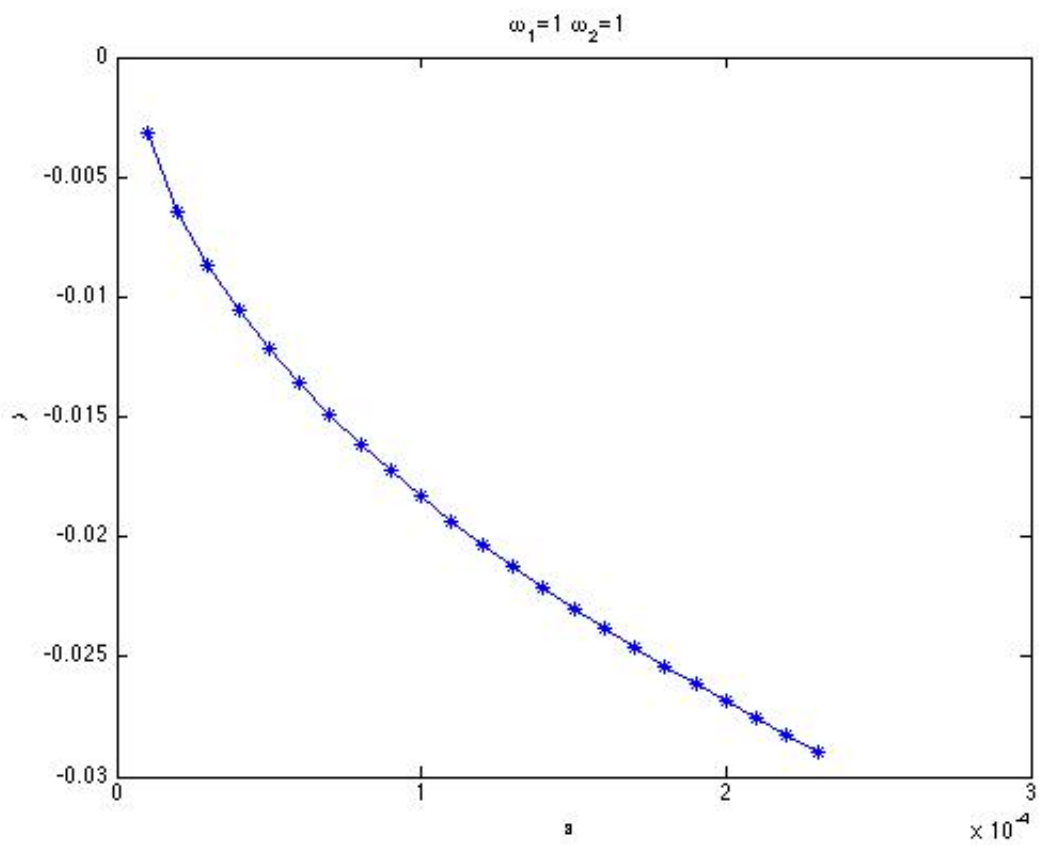


Figure 13: Value of λ in ω case.

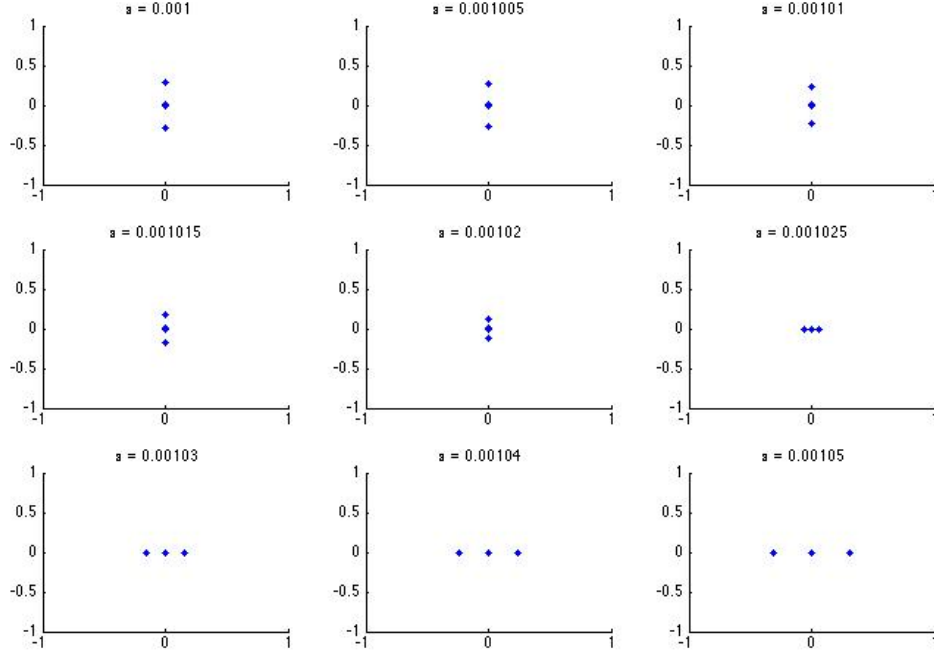


Figure 14: Spectrum of μ case for $\lambda = 0$.

4.3 μ case

1. $\lambda = 0$, we can find ϵ^* in this solution case. For $\mu_1, \mu_2 = 1, 2, \dots, 5$, we find that the eigenvalues are pure imaginary when $\epsilon < \epsilon^*$, and turn into a pair real value. Figure 14 plot the eigenvalues around original. And in Figure 15 show the ϵ^* for every μ_1, μ_2 .
2. $\lambda = 1$, we also find ϵ^* . Figure 16, 17 show the eigenvalues and values of ϵ^* for each μ_1, μ_2 .
3. Consider the same μ_1 and μ_2 , these two solution forms can be seen as changing the well depth of potential (from (5)). The ϵ^* are different. So we know that the potential well depth may change the stability of solution.

In optical lattices potential case, changing μ_1 or μ_2 for different periodic of potential. We can see that the value of ϵ^* increasing when μ_1, μ_2 go larger. To comparison $e^{it}\phi(x)$ and $e^{i0t}\phi(x)$ these two forms of solutions, we can see that the second kind of ϵ^* is less than the other ones.

4. λ is general case, there is no ϵ^* in this solution case. When ϵ is quit small(less than 1.2×10^{-4}), the value of λ is positive(In Figure 18). This result is consistent with the condition of Lin in [9](λ is positive for small enough ϵ). But λ become negative when ϵ gets larger, and the value of λ gets smaller as ϵ gets larger. When $\epsilon < 3.3 \times 10^{-4}$, λ is smaller than -1 . And the minimum of trap potential is 1, that means the trap potential is no longer positive value.

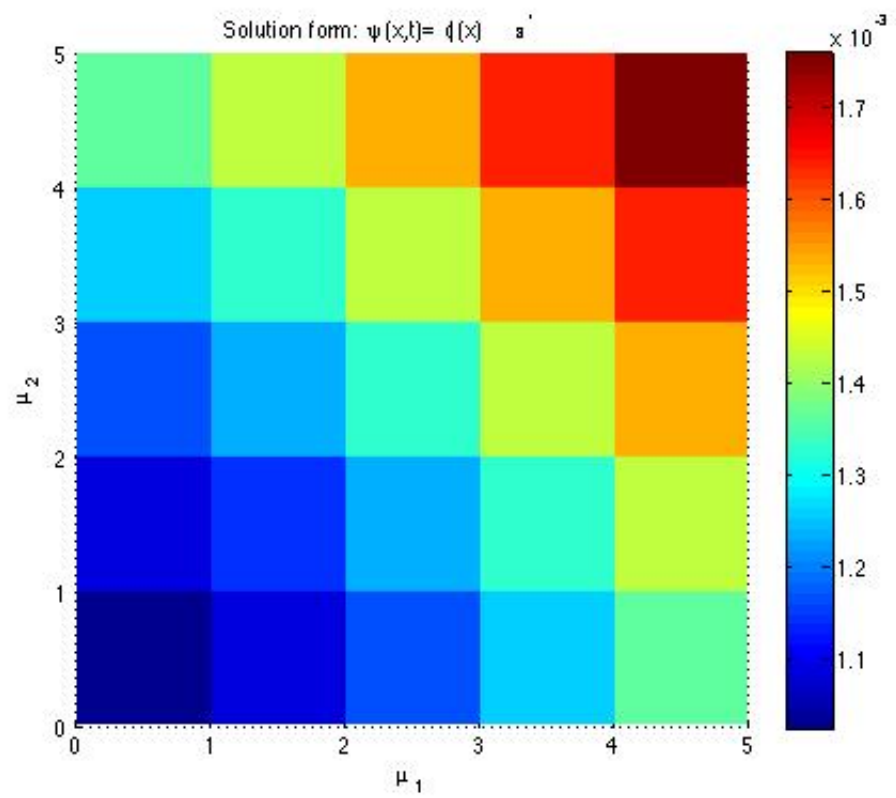


Figure 15: ϵ^* of μ case for $\lambda = 0$.

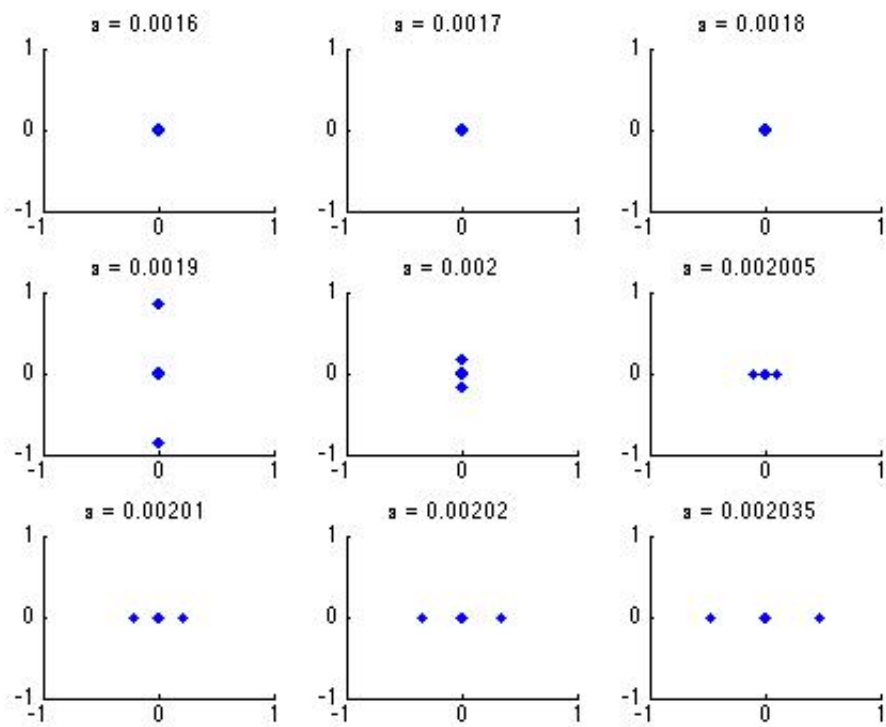


Figure 16: Spectrum of μ case for $\lambda = 1$.

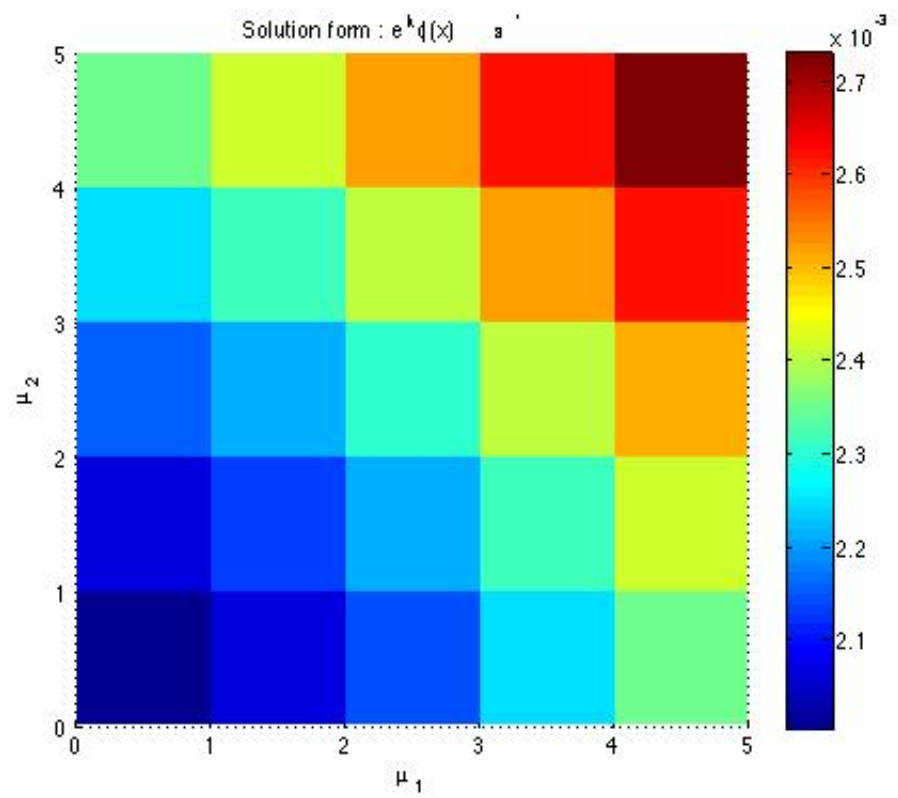


Figure 17: ϵ^* of μ case for $\lambda = 1$.

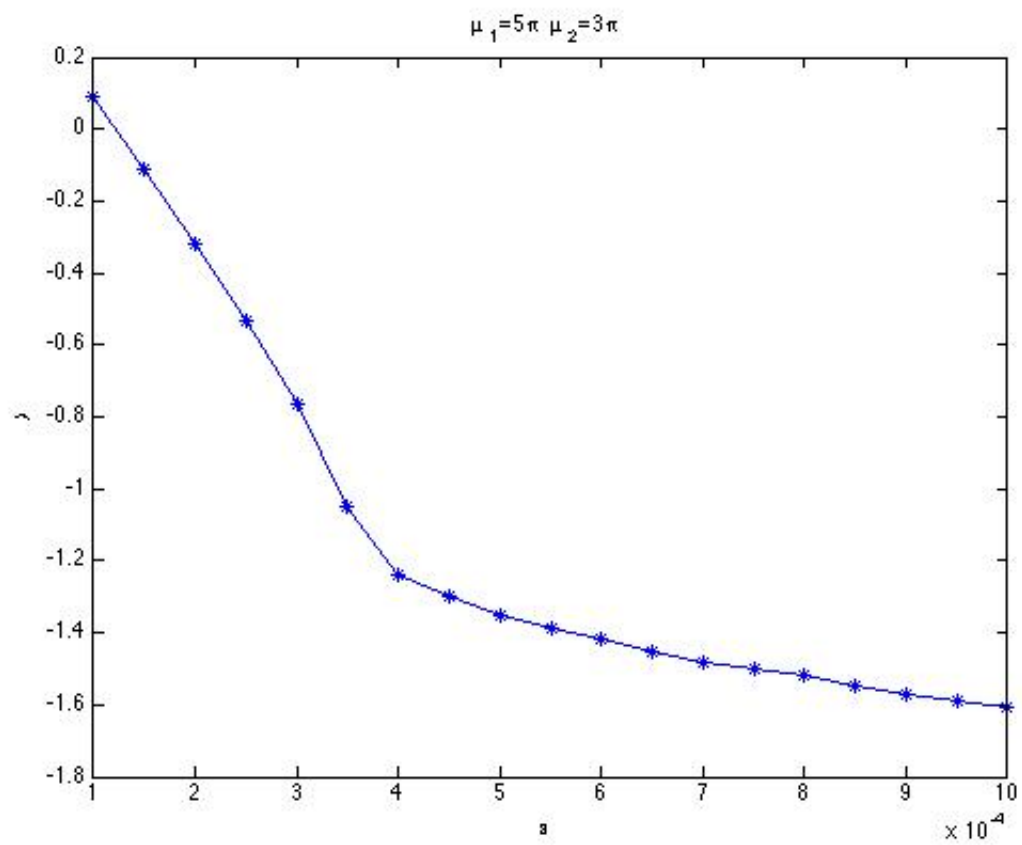


Figure 18: Value of λ in μ csae.

5 Conclusions

In this project, we consider the NLS equation with the special solution $e^{i\lambda t}\phi(x)$. We mainly show the spectrum of the linearization operator \mathcal{L} . By numerical computation, we can conclude some results: when $\epsilon < \epsilon^*$ the eigenvalues are pure imaginary, and the eigenvalues turn to real as $\epsilon > \epsilon^*$. That is, there is a bifurcation of stable and unstable, when $\lambda = 0, 1$ in optical lattices potential case and $\lambda = 1$ in elliptic potential case.

Here, we use Matlab code to find the spectrum of the linearization operator. It is a convenient way to solve the eigenvalue problem. Since the method of finding eigenvalue in Matlab is to search some eigenvalues which are near the target. And the distribution of eigenvalues changes fast when ϵ near ϵ^* .

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出席國際學術會議心得報告

報告人姓名	張書銘	服務機構 及職稱	國立交通大學 助理教授
會議時間 地點	2010.06.27 - 2010.07.01 德國、柏林	本會核定 補助文號	NSC 98-2115-M-009-007
會議 名稱	(中文) 第八屆特徵值問題精確解國際會議 (英文) 8th International Workshop on Accurate Solution of Eigenvalue Problems (IWASEP 8)		

報告內容包括下列各項：

一、參加會議經過

本人於2010年2月開始規劃要參加2010年6月底的IWASEP 8-第八屆特徵值問題精確解國際會議，此國際會議是每兩年固定時間主辦的一個特徵值問題計算的重要聚會，此會議雖然僅為期四天，但是這四天的學術演講都是單場的，沒有平行的演講場次。因此，每個參與者都是同領域的學者，演講者更是在此方面的研究的翹楚者。

這次的參與是本人第一次以參與者身份出席國際會議，雖然沒有在外語演講的經驗上有所累積，但在與外國人的參與者之互動上，本人有更多的機會與其對談，並且進一步認識了荷蘭的學者Michiel Hochstenbach教授(Dept. of Math. And Computer Science, TU Eindhoven, Netherlands)。期望今年或是明年，能夠邀請Prof. Hochstenbach來台灣訪問。

二、與會心得

首先要感謝行政院國家科學委員會給予本人出席國際會議的差旅補助，讓本人有機會去見識國際會議，參與過程本人更增加不少國際觀與吸收到專業的研究知識。

當中本人與Prof. Hochstenbach有相當多次的交談，其研究領域相當廣泛(Numerical linear algebra, Eigenvalue problems, Model reduction, Ill-posed problems, inverse problems, regularization, Image analysis, Matrix functions, Linear systems, Scientific Computing, Nonlinear systems, Boundary element method, Numerical crack propagation, Industrial mathematics via LIME)。期望今年或是明年，能夠邀請Prof. Hochstenbach來台灣訪問。

三、考察參觀活動

無。

四、建議

無。

五、攜回資料名稱及內容

大會演講論文全文一冊及會議演講摘要手冊(內容包括此次會議的所有演講場次與講題)，收集完整的大會演講內容，相當豐富。

六、其他

無。

無研發成果推廣資料

98 年度專題研究計畫研究成果彙整表

計畫主持人：張書銘		計畫編號：98-2115-M-009-007-				計畫名稱：格若士-比塔烏斯基方程式的行進波解之穩定性研究	
成果項目		量化			單位	備註（質化說明：如數個計畫共同成果、成果列為該期刊之封面故事...等）	
		實際已達成數（被接受或已發表）	預期總達成數（含實際已達成數）	本計畫實際貢獻百分比			
國內	論文著作	期刊論文	0	0	0%	篇	
		研究報告/技術報告	0	0	0%		
		研討會論文	0	0	0%		
		專書	0	0	0%		
	專利	申請中件數	0	0	0%	件	
		已獲得件數	0	0	0%		
	技術移轉	件數	0	0	0%	件	
		權利金	0	0	0%	千元	
	參與計畫人力（本國籍）	碩士生	3	0	100%	人次	研究生兼任助理參與本計畫，在程式設計能力上均有成長。
		博士生	0	0	0%		
博士後研究員		0	0	0%			
專任助理		0	0	0%			
國外	論文著作	期刊論文	0	0	0%	篇	
		研究報告/技術報告	0	0	0%		
		研討會論文	0	0	0%		
		專書	0	0	0%		章/本
	專利	申請中件數	0	0	0%	件	
		已獲得件數	0	0	0%		
	技術移轉	件數	0	0	0%	件	
		權利金	0	0	0%	千元	
	參與計畫人力（外國籍）	碩士生	0	0	0%	人次	
		博士生	0	0	0%		
博士後研究員		0	0	0%			
專任助理		0	0	0%			

<p style="text-align: center;">其他成果</p> <p>(無法以量化表達之成果如辦理學術活動、獲得獎項、重要國際合作、研究成果國際影響力及其他協助產業技術發展之具體效益事項等，請以文字敘述填列。)</p>	<p style="text-align: center;">無</p>
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	成果項目	量化	名稱或內容性質簡述
科 教 處 計 畫 加 填 項 目	測驗工具(含質性與量性)	0	
	課程/模組	0	
	電腦及網路系統或工具	0	
	教材	0	
	舉辦之活動/競賽	0	
	研討會/工作坊	0	
	電子報、網站	0	
	計畫成果推廣之參與(閱聽)人數	0	

國科會補助專題研究計畫成果報告自評表

請就研究內容與原計畫相符程度、達成預期目標情況、研究成果之學術或應用價值（簡要敘述成果所代表之意義、價值、影響或進一步發展之可能性）、是否適合在學術期刊發表或申請專利、主要發現或其他有關價值等，作一綜合評估。

1. 請就研究內容與原計畫相符程度、達成預期目標情況作一綜合評估

達成目標

未達成目標（請說明，以 100 字為限）

實驗失敗

因故實驗中斷

其他原因

說明：

2. 研究成果在學術期刊發表或申請專利等情形：

論文： 已發表 未發表之文稿 撰寫中 無

專利： 已獲得 申請中 無

技轉： 已技轉 洽談中 無

其他：（以 100 字為限）

3. 請依學術成就、技術創新、社會影響等方面，評估研究成果之學術或應用價值（簡要敘述成果所代表之意義、價值、影響或進一步發展之可能性）（以 500 字為限）

本計畫原訂是規畫三年的，但僅獲得一年的補助。對於格若士-比塔烏斯基方程式的行進波解之穩定性研究，目前已達到原撰寫計畫書時所規劃的進度，在數值計算方法上設計出一個能夠求解格若士-比塔烏斯基方程式的行進波解，並期望能夠在不同參數下均可以用此數值計算方法穩定的求解行進波。若要能就有較有價值的研究成果，需要持續進行下去。

