

An Energy-Preserving Scheme for the Coupled Gross-Pitaevskii Equations

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Abstract. An energy-preserving scheme is proposed for the coupled Gross-Pitaevskii equations. The scheme is constructed by high order compact method in the spatial direction and average vector field method in the temporal direction, respectively. The scheme is energy-preserving, stable, and of sixth order in space and of second order in time. Numerical experiments verify the theoretical results. The dynamic behavior modeled by the coupled Gross-Pitaevskii equations is also numerically investigated.

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Key words: Coupled Gross-Pitaevskii equations, average vector field method, high order compact method, energy-preserving scheme.

1 Introduction

In Bose-Einstein condensates (BECs) community, using mean field approximation, the two-component time-dependent dimensionless coupled Gross-Pitaevskii (CGP) equations with external potentials and effective Rabi frequencies is as follows [11, 16, 22, 23, 25, 35]

$$iu_t = \left(-\frac{1}{2}\Delta + s_1(x, y) + \beta_{11}|u|^2 + \beta_{12}|v|^2 \right) u - L_z u - \lambda v, \quad (x, y) \in \mathbb{R}^2, \quad t > 0, \quad (1.1a)$$

$$iv_t = \left(-\frac{1}{2}\Delta + s_2(x, y) + \beta_{21}|u|^2 + \beta_{22}|v|^2 \right) v - L_z v - \lambda u, \quad (x, y) \in \mathbb{R}^2, \quad t > 0, \quad (1.1b)$$

where $i^2 = -1$, and Δ is the Laplace operator. We only consider two-dimensional case, i.e., $\Delta = \partial_{xx} + \partial_{yy}$ in this paper. $u(x, y, t)$, $v(x, y, t)$ are the complex-valued wave functions.

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$s_l(x, y)$ ($l=1,2$) are the real-valued external trapping potential functions. In the harmonic potentials, they are generally in the forms

$$s_l(x, y) = \frac{1}{2}(\omega_{x,l}^2 x^2 + \omega_{y,l}^2 y^2)$$

with dimensionless frequencies $\omega_{x,l}$ and $\omega_{y,l}$ ($l=1,2$) in the x and y directions, respectively. The action between intra-component and inter-component atomic is represented by

$$\beta_{lk} = \beta_{kl} = \frac{4\pi N_0 \alpha_{lk}}{\alpha_0}, \quad k, l = 1, 2,$$

with N_0 being the total numbers of the particles in BECs, α_0 being the dimensionless spatial unit, and $\alpha_{kl} = \alpha_{lk}$ being the s -wave scattering lengths between the l -th and k -th components (positive for repulsive interaction and negative for attractive interaction). L_z is the rotating angular momentum with rotating speed Ω , defined as

$$L_z = i\Omega(y\partial_x - x\partial_y).$$

λ is the effective Rabi frequency to realize the internal atomic Josephson junction by a Raman transition. If there is no effective Rabi frequency and no rotation, i.e., $\lambda = 0$ and $\Omega = 0$, the CGP equations collapse to the coupled nonlinear Schrödinger equations, they have been discussed by many authors [21].

Under appropriate assumption, to make the problem well-posed, with the following initial data

$$u(x, y, 0) = u_0(x, y), \quad v(x, y, 0) = v_0(x, y), \quad (1.2)$$

the problem (1.1a)-(1.2) satisfies the following laws [5].

- The component-masses

$$\mathcal{M}_u(t) = \int_{\mathbb{R}^2} |u(x, y, t)|^2 dx dy \quad \text{and} \quad \mathcal{M}_v(t) = \int_{\mathbb{R}^2} |v(x, y, t)|^2 dx dy$$

with respect to the u -component and v -component particles, respectively, satisfy the equalities

$$\frac{d}{dt} \mathcal{M}_u(t) = -\lambda \int_{\mathbb{R}^2} \left(v(x, y, t) \overline{u(x, y, t)} - u(x, y, t) \overline{v(x, y, t)} \right) dx dy, \quad (1.3a)$$

$$\frac{d}{dt} \mathcal{M}_v(t) = -\lambda \int_{\mathbb{R}^2} \left(u(x, y, t) \overline{v(x, y, t)} - v(x, y, t) \overline{u(x, y, t)} \right) dx dy. \quad (1.3b)$$

Therefore, the total mass of the particles of the system is an invariant, that is,

$$\mathcal{M}(t) = \mathcal{M}_u(t) + \mathcal{M}_v(t) = \mathcal{M}(0). \quad (1.4)$$

Furthermore, $\mathcal{M}_u(t) = \mathcal{M}_u(0)$ and $\mathcal{M}_v(t) = \mathcal{M}_v(0)$ in case of $\lambda = 0$. This indicates that the particle masses of each component are conserved.