

# **Collective Beam-Beam: Theory and Simulation**

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# Contents

- Strong-Strong Beam-Beam Model and the results from paper "*A new model for the collective beam-beam interaction*" by J. A. Ellison and M. Vogt.
- Calculating eigen modes. Difficulties of numerical solution related to the type of the problem (3<sup>rd</sup> kind integral equation, not compactness of the operator.)
- Direct simulation of the density evolution. WMPT vs. PF approach. Parallel code. Numerical results.

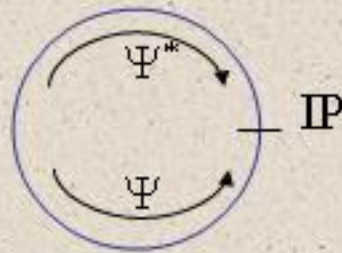


# Steps

- Rewrite in slow varying coordinates;
- Using averaging for the map, obtain continuous-time Vlasov equation for  $\Psi(t, v)$ ;
- Linearize this equation around equilibrium,
- Obtain integro-differential equation for  $\pi$  and  $\sigma$ -modes;
- Using Laplas transformation, obtain an integral equation;
- Make Fourier transformation in angle and obtain the system of decoupled integral equation.
- Study the stability by analyzing the spectrum of the integral operator.

# Model

- Strong-Strong Beam interaction
- Study density evolution in 2dof case
- One interaction point. Linear lattice.
- Two counter-rotating beams collide at IP.



The beams have densities

$$\psi_n(x, p_x, y, p_y), \quad \psi_n^*(x, p_x, y, p_y),$$



# Evolution Equation

The position of the particle in 4d phase space before the interaction point on  $n$ -th turn

$$u_n = (x, p_x, y, p_y)$$

changes according to

$$u_{n+1} = M \left( u_n + \zeta \begin{pmatrix} 0 \\ V_x(x_n, y_n) \\ 0 \\ V_y(x_n, y_n) \end{pmatrix} \right)$$

# Evolution Equation ★

The potential can be calculated using electrostatic formula in the frame of the moving beam.

$$V(\psi^*)(x, y) = \int_{\mathbb{R}^2} G(x - x', y - y') \rho^*(x', y') dx' dy'$$

$$G(x, y) = -\ln(\sqrt{x^2 + y^2}),$$

The potential in the frame of the kicked beam is calculated by multiplying by

$$\zeta \approx \frac{2N^*}{\gamma} \frac{qq^*}{4\pi\epsilon_0 mc^2},$$





# Equation in Slow-Varying Coordinates

Transformation to slow varying variables

$$u_n = R^n v_n$$

leads to

$$v_{n+1} = v_n + \zeta J_4 \nabla_v H(\Psi_n^*)(v_n, n),$$

where

$$H(\Phi)(v, n) = \int G(\sqrt{2\beta_x J_x} \sin(2\pi m v_x + \Theta_x) - \sqrt{2\beta_x' J_x'} \sin(2\pi m v_x^* + \Theta_x'), x \rightarrow y) \Phi(v') dv'.$$

Averaging leads to

$$v_{n+1} = v_n + \zeta J_4 \nabla_v \bar{H}(\Psi_n^*)(v_n) + O(\zeta^2).$$

This corresponds to the equations

$$\partial_t \Psi + \zeta \{\Psi, \bar{H}(\Psi^*)\} = 0, \quad \partial_t \Psi^* + \zeta^* \{\Psi^*, \bar{H}^*(\Psi)\} = 0.$$

- Linearize about equilibrium

$$\Psi(\mathbf{v}, t) = \Psi_e(\mathbf{v}) + \Psi_1(\mathbf{v}, t)$$

- Study  $\pi$  and  $\sigma$ -modes:

$$f = \Psi_1 \pm \Psi_1^*$$

- We are looking for the solutions in form.

$$g(J, t) = e^{i\omega t} \hat{g}(J, \omega)$$

- Azimuthal Fourier modes are the solutions of the uncoupled integral equation of the third kind.

$$(\omega - k \cdot \Omega_e(J)) \hat{g}(J, \omega) - \eta \int K_k(J, J') \hat{g}(J', \omega) dJ' = 0$$



## Difficulties of the numerical solution of the 3<sup>rd</sup> kind integral equation

- This kind of equation has been studied by Warnock (*“Linear Integral Equations of the third kind”*, G. R. Bart R.L. Warnock, SIAM, vol. 4, No. 4, Nov. 1973).
- This integral operator has continuous spectrum with eigenfunctions being delta functions. It makes it hard to present numerically. The operator is not compact therefore discretization of these equation doesn't converge.

# Regularizing the Integral Equation

- The equation must be transformed. For the 1dof case this was made in "*Linear Vlasov Analysis for stability of a bunched beam*" by R. Warnock, G. Stupakov, M. Venturini, Ellison and "*Nonsingular integral Equation for stability of bunched beam*" by R. Warnock, M. Venturini, and J. Ellison.

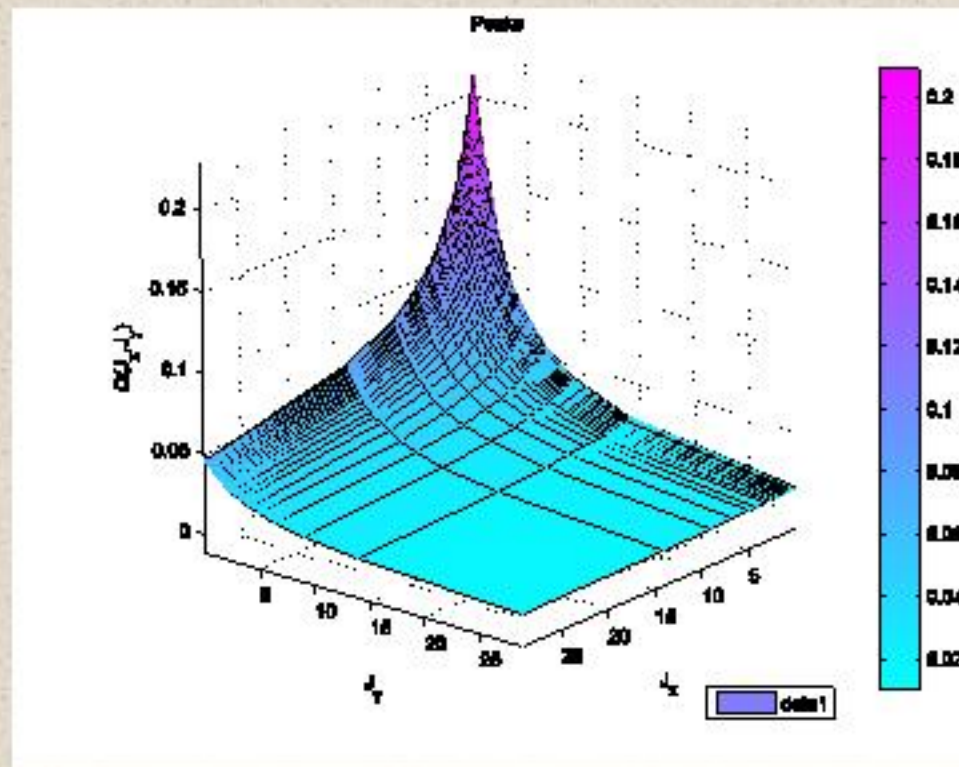
$$\hat{g}(J, \omega) - \eta \int \frac{\hat{g}(J', \omega)}{(\omega - k \cdot \Omega_e(J))} K_k(J, J') dJ' = 0$$

- We are calculating eigenvalues and eigenfunctions for 2dof case

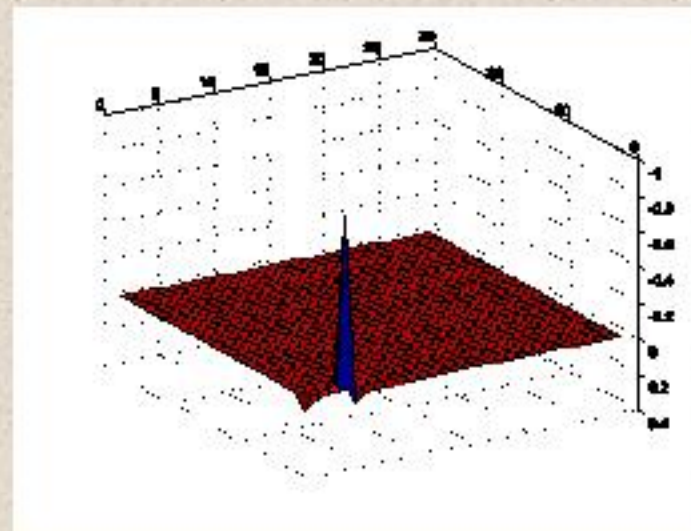
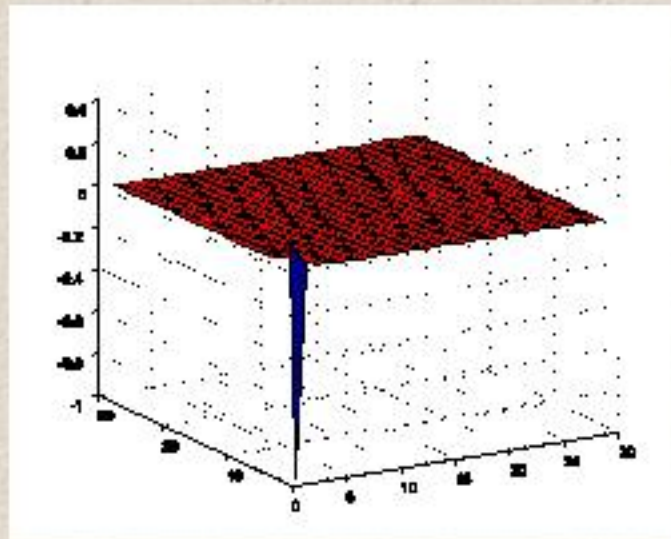
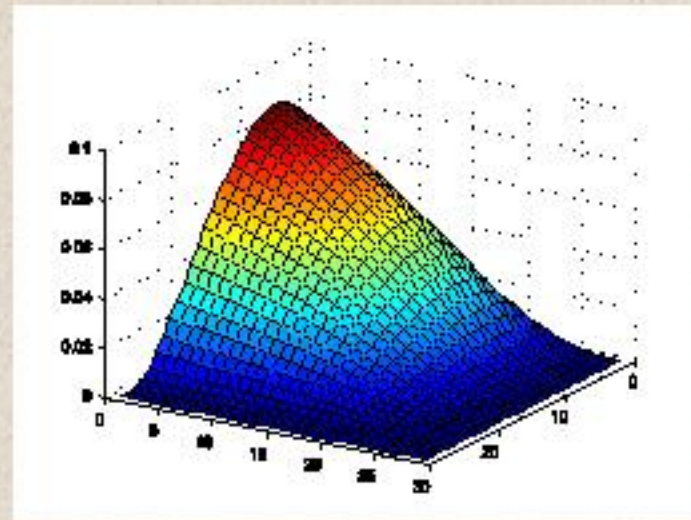
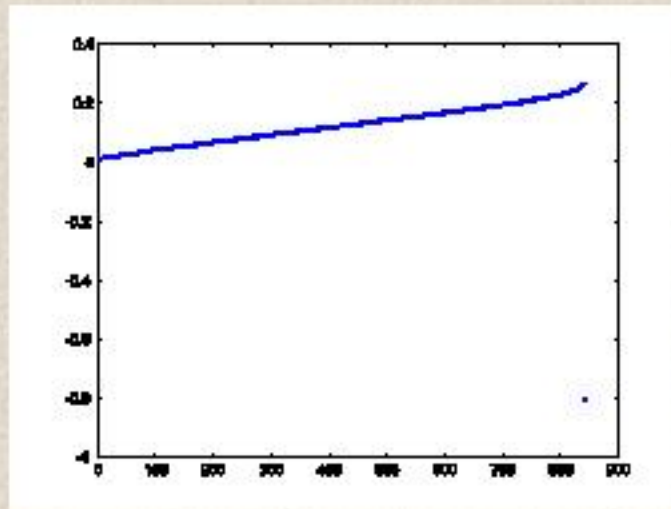


# 2dof case discretization

- To calculate the coefficients of this integral equation we have to deal with high dimensionality of the problem and the necessity of taking improper integrals.

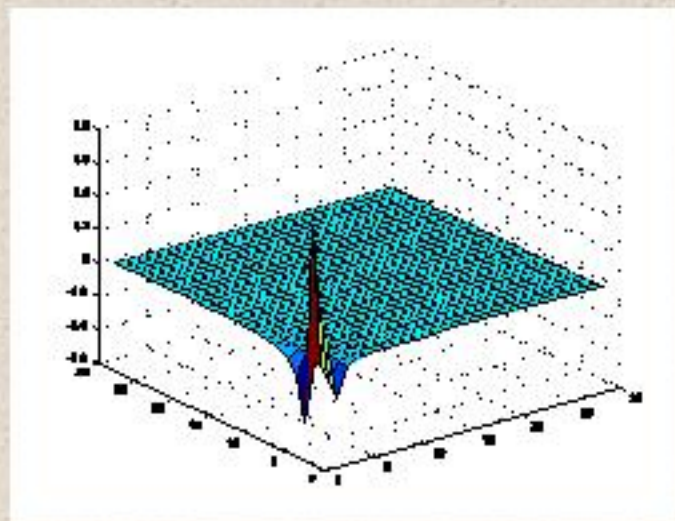
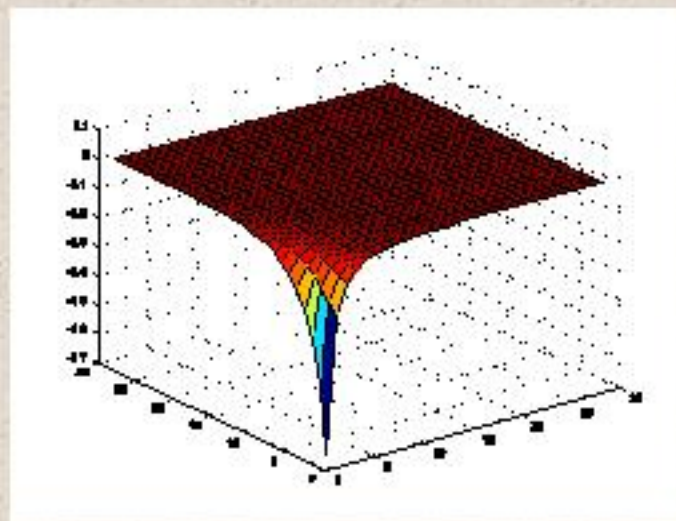
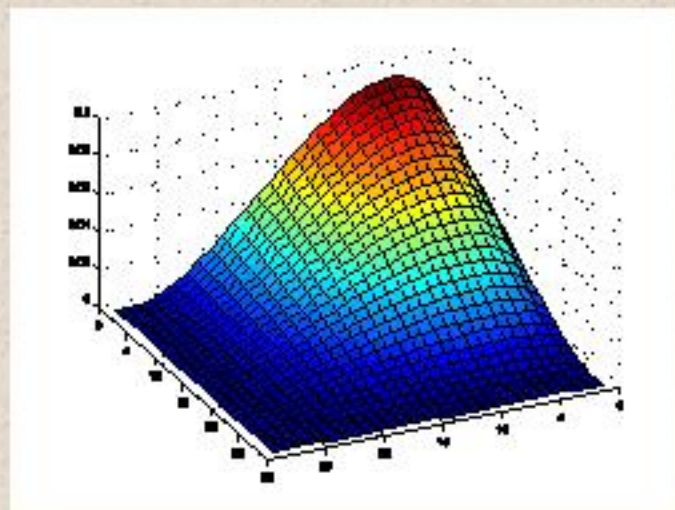
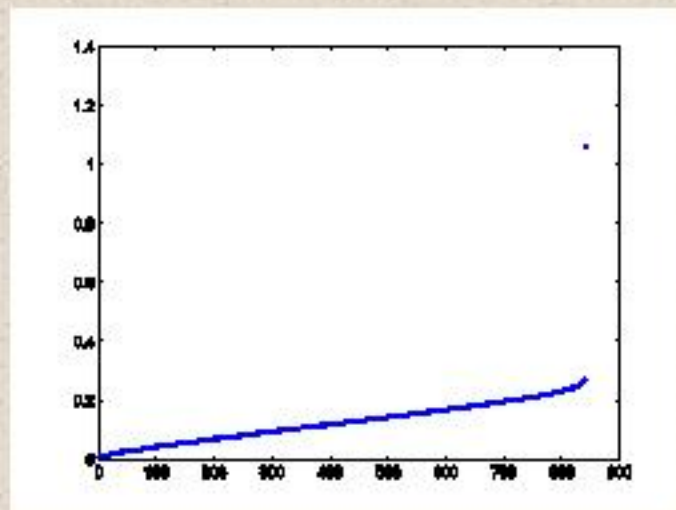


# $\sigma$ -mode





# $\pi$ -mode



# Analytical result

- Warnock suggested that the transformed equation is compact in some norm. I think I proved that it is indeed compact in

$$\|f\|_a = \sup \left| \frac{f(x) - f(a)}{x - a} \right| + |f(a)|$$

- This is an important step to find a convergent discretisation of the equation.

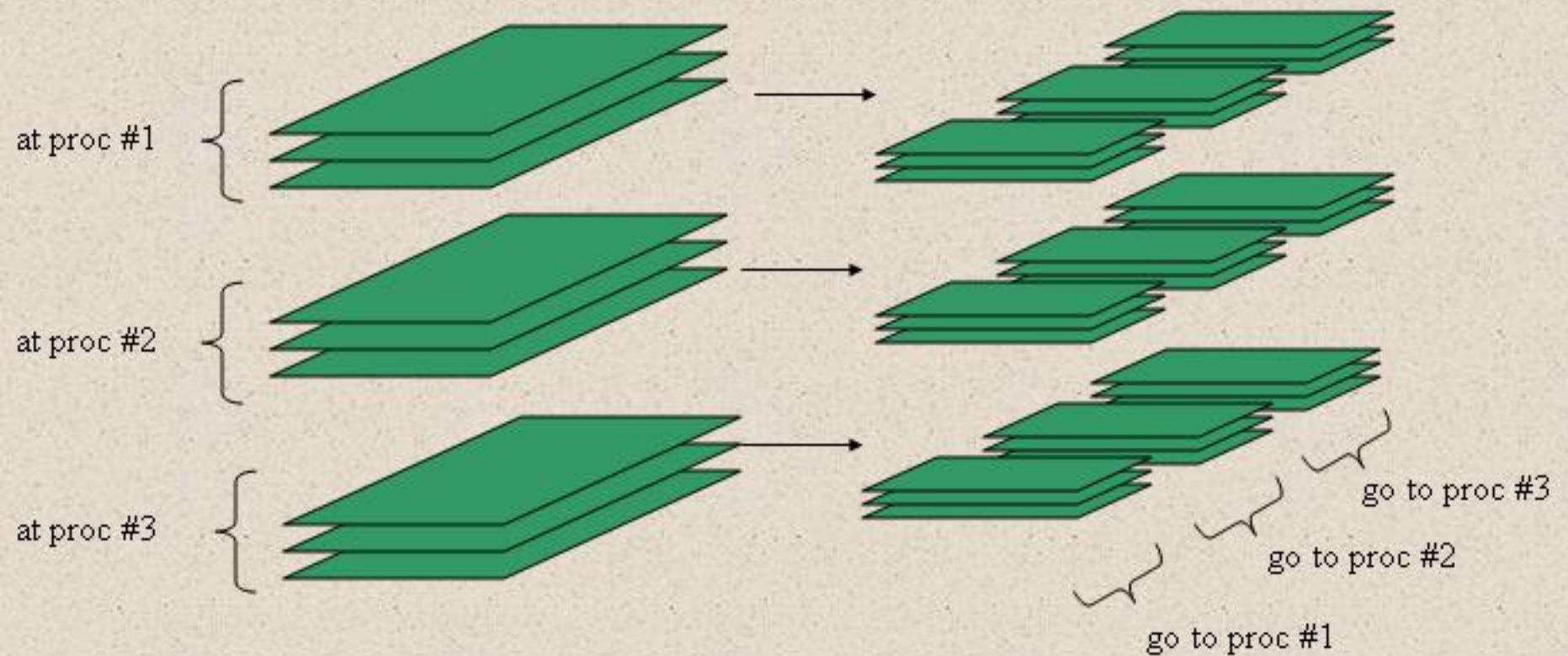


# Direct Simulations

- Macroparticle approach
- Weighted macroparticle approach. WMPT code by Mathias Vogt.
- PF method: 1 d.o.f. is done by Bob Warnock and Mathias Vogt.
- 2 dof. Simulation in 4 dimensional phase space. Need of parallel computations.

$$u_{n+1} = R \circ K[\Psi_n^*](u_n), \quad u_{n+1}^* = R \circ K[\Psi_n](u_n^*),$$
$$\Psi_{n+1} = \Psi_n \circ K^{-1}[\Psi_n^*] \circ R^{-1}, \quad \Psi_{n+1}^* = \Psi_n \circ K^{-1}[\Psi_n] \circ R^{-1},$$

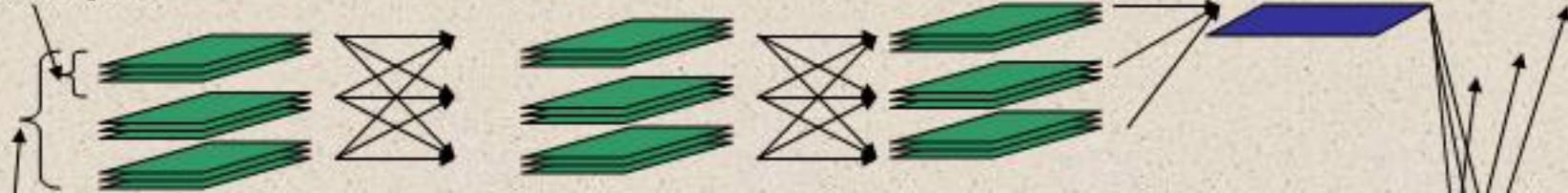
# Transformation of the Data





# Transformation of the Data

main proc



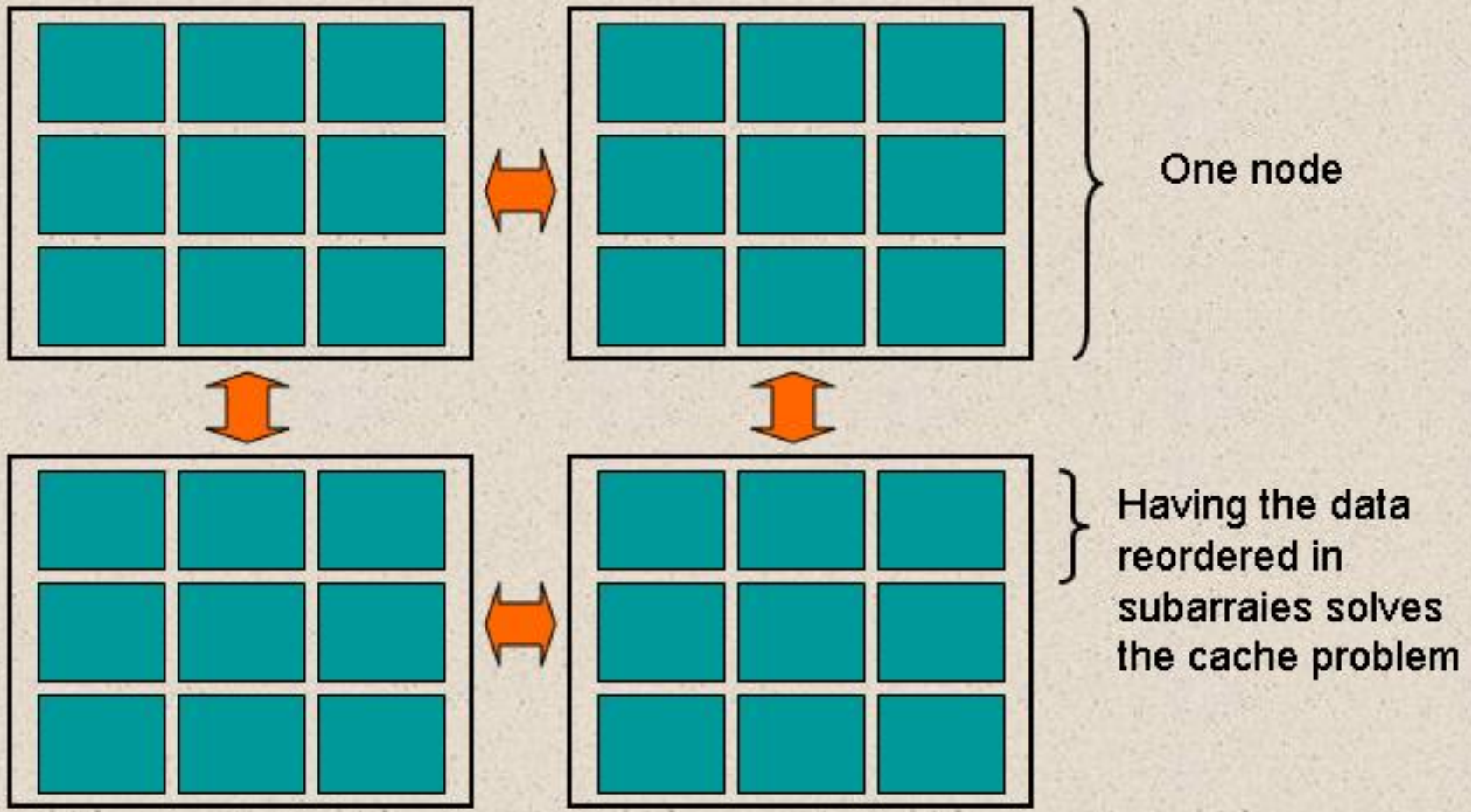
Beam #1

main proc



Beam #2

# Density evolution in slow varying coordinates in Slow Varying coordinates





# First results of parallel simulations

- Round Gaussian Source approximation
- Finding kick by solving Poisson equation.

