

構造を仮定しない計算による 低密度原子核物質の非一様構造

Non-uniform structure of low-density nuclear matter
by three-dimensional calculation
without any assumption of the structures

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▪ *Menu*

I . Pasta structure

II . Relativistic mean field theory

III . Results

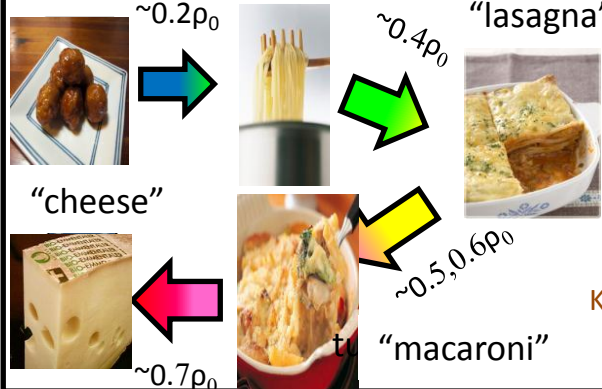
i) Fixed proton fraction

ii) β -equilibrium

IV . Conclusion

Pasta Structure : “Mixed phase with structure” in first phase transition

“meat ball” “spaghetti” Suggested by D.G.Ravenhall, C.J.Pethick, J.R.Wilson in 1983



Blue : nuclear matter
cyan : gas nuclei
 ρ_0 : normal nucleus density
 $\sim 0.16\text{fm}^{-3}$

Balance with Surface tension
and Coulomb repulsion

$$2E_C = E_S \quad (\text{Equilibrium condition})$$

K.Oyamatsu, NPA 561 (1993) 431

Pasta structure in Neutron star

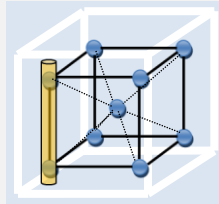
- Half mass of Crust region

→ Influence on EOS of Crust

- Glitch

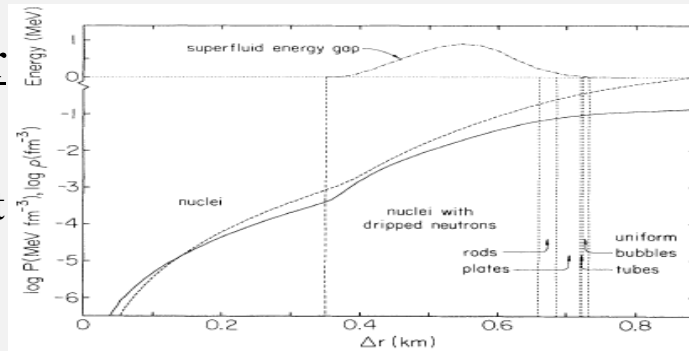
Rapid increasing of rotational speed of Neutron Star

Whirlpool of super fluidity of neutron pinned to nuclei



Nuclei

Super fluidity (Neutron)



C.P.Lorenz et al, PRL 70 (1993) 4

It may occur from the collapse of whirlpool of neutron super fluidity on lattice points?

→ It may depend on the crystalline of nuclei and more exotic structure like “pasta”

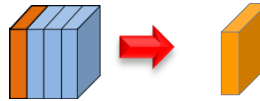
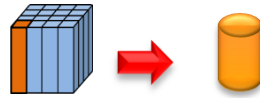
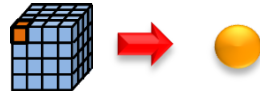
- Neutrino cross-section in supernova and collapsing stellar cores

Conventional Studies

Wigner-Seitz(WS) cell approximation

Whole space is divided into equivalent cells.
These are imposed geometrical symmetry as follow

- 3D → sphere
- 2D → cylinder
- 1D → plate



Merit : All calculations are 1D's

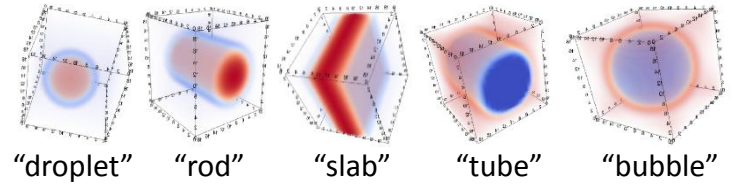
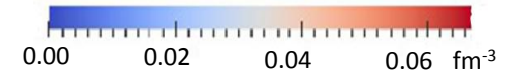


Fast calculation & low cost

But ...

Only simple structure appear

Calculation in small cell



Proton density distribution (based on RMF)

- Fast calculation & low cost
- Several complex structures appear as meta stable state
- Restriction of periodic boundary condition

- Existence of Non-typical pasta structure

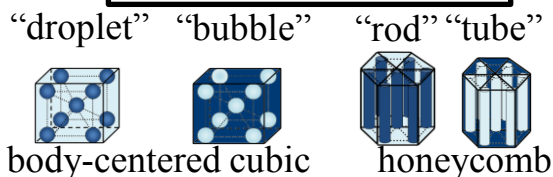
M.Matsuzaki PRC 73, 028801 (2006)

K.Nakazato,K.Oyamatsu PRL 103, 132501 (2009)

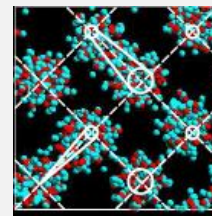
G.Watanabe et.al PRC 68, 035806 (2003)

- Crystalline of Pasta structure

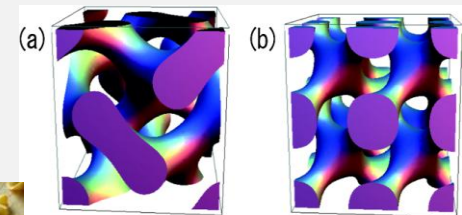
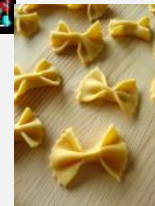
Conventional studies



K.Oyamatsu, NPA 561 (1993) 431



dumbbell (farfalle)



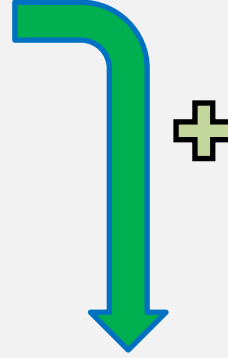
Gyroid double-diamond



Full real space
3D calculation

Relativistic mean field theory

$$\begin{aligned}
 L = & \bar{\psi} \left(i\gamma^\mu \partial_\mu - m - g_\sigma \sigma - g_\omega \gamma^\mu \partial_\mu - g_\rho \gamma^\mu \tau^a \rho_\mu^a \right) \psi \\
 & + \frac{1}{2} \partial_\mu \sigma \partial^\mu \sigma - \frac{1}{2} m_\sigma^2 \sigma^2 + \frac{1}{3} b m_\sigma (g_\sigma \sigma)^3 + \frac{1}{4} c (g_\sigma \sigma)^4 \\
 & + \frac{1}{2} m_\omega^2 \omega_\mu \omega^\mu + \frac{1}{2} \left[\frac{1}{2} (\partial^\mu \omega^\nu - \partial^\nu \omega^\mu) (\partial_\mu \omega_\nu - \partial_\nu \omega_\mu) \right] \\
 & + \frac{1}{2} m_\rho^2 R_\mu R^\mu - \frac{1}{4} R_{\mu\nu} R^{\mu\nu} \\
 & - \frac{1}{2} \left[\frac{1}{2} (\partial^\mu A^\nu - \partial^\nu A^\mu) (\partial_\mu A_\nu - \partial_\nu A_\mu) \right]
 \end{aligned}$$



Mean field approximation

$$\langle \sigma \rangle = \sigma$$

$$\langle \omega^\mu \rangle = \delta^{\mu,0} \omega^\mu$$

$$\langle R^{\mu\nu} \rangle = \delta^{\nu,0} R^{\mu\nu}$$

Thomas-Fermi approximation
&

Zero temperature

$$-\nabla^2 \sigma(\vec{r}) + m_\sigma^2 \sigma^2(\vec{r}) = -\frac{dU(\sigma)}{d\sigma} + g_{\sigma N} (\rho_n^s(\vec{r}) + \rho_p^s(\vec{r}))$$

$$-\nabla^2 \omega_0(\vec{r}) + m_\omega^2 \omega_0(\vec{r}) = g_{\omega N} (\rho_p(\vec{r}) + \rho_n(\vec{r}))$$

$$-\nabla^2 R_0(\vec{r}) + m_\rho^2 R_0(\vec{r}) = g_{\rho N} (\rho_p(\vec{r}) - \rho_n(\vec{r}))$$

$$\nabla^2 V(\vec{r}) = 4\pi e^2 (\rho_p(\vec{r}) + \rho_n(\vec{r}))$$

$$\mu_p = \mu_B - \mu_e = \sqrt{k_{F,p}^2 + m_N^{*2}} + g_{\omega N} \omega_0 + g_{\rho N} R_0 - V$$

$$\mu_n = \mu_B = \sqrt{k_{F,n}^2 + m_N^{*2}} + g_{\omega N} \omega_0 - g_{\rho N} R_0$$

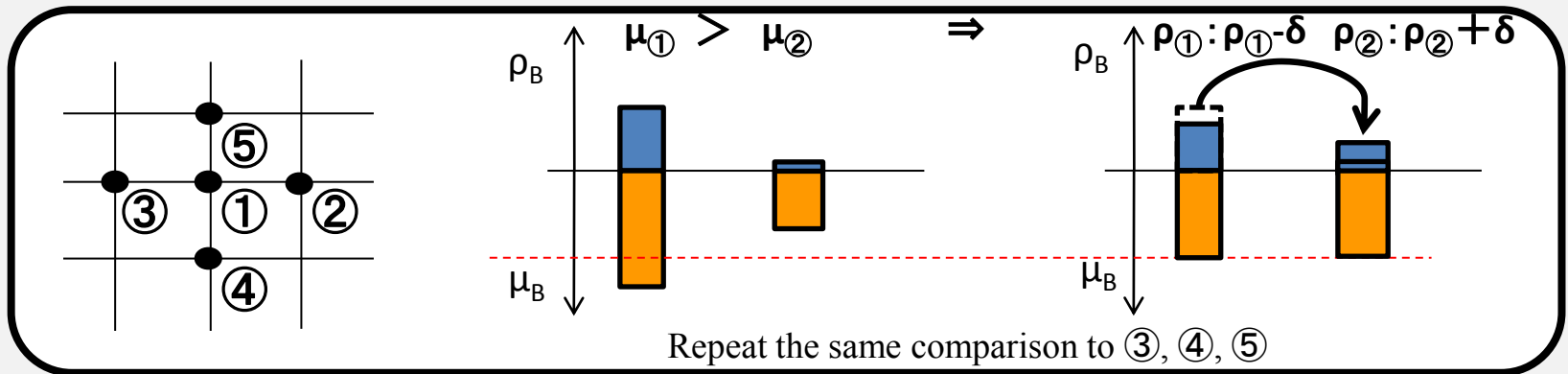
$$\rho_e(\vec{r}) = -\frac{(V(\vec{r}) - \mu_e)^3}{3\pi^2}$$

Solutions

- As an initial condition, randomly distribute fermions (n, p, e) over the grid
- We solve coupled differential equations, and simultaneously relax fermions density distributions to attain the uniformity of their chemical potential

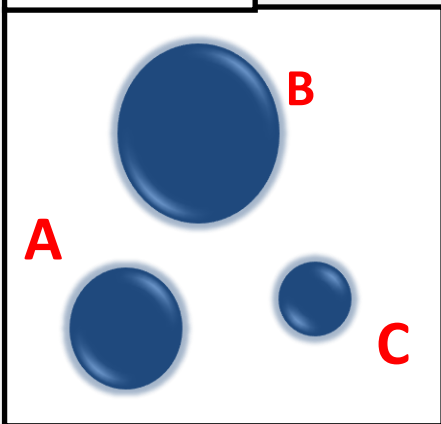
$$\mu_p(\vec{r}) = \sqrt{k_{F,p}^2(\vec{r}) + m_N^{*2}(\vec{r})} + g_{\omega N}\omega_0(\vec{r}) + g_{\rho N}\rho_0(\vec{r}) - V(\vec{r})$$

$$\mu_n(\vec{r}) = \sqrt{k_{F,n}^2(\vec{r}) + m_N^{*2}(\vec{r})} + g_{\omega N}\omega_0(\vec{r}) - g_{\rho N}\rho_0(\vec{r})$$



- Exchanging the baryon density, we solve coupled differential equation again.

In process...

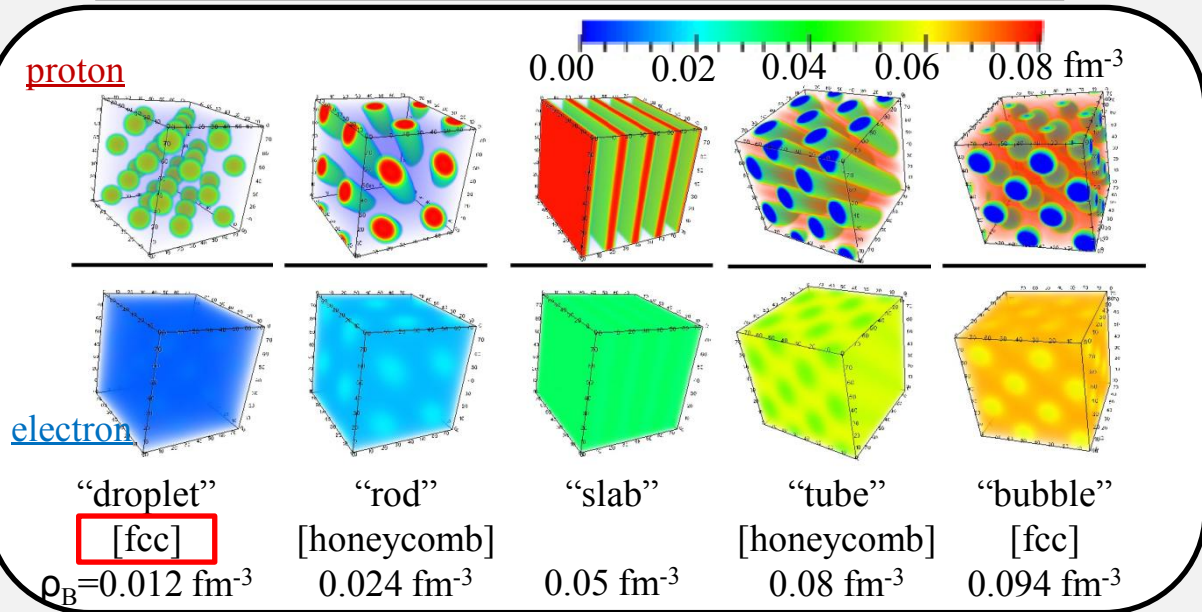


- We adjust the particle densities between distant grid points chosen randomly so as to avoid making separate droplets with different chemical potentials
- To save the calculation cost, we did not compare the chemical potential with all grids point and performed the comparison of the chemical potentials with the small regions frequently.

Results : Fixed proton ratio $Y_p=Z/A$

(A : total mass , Z : proton number)

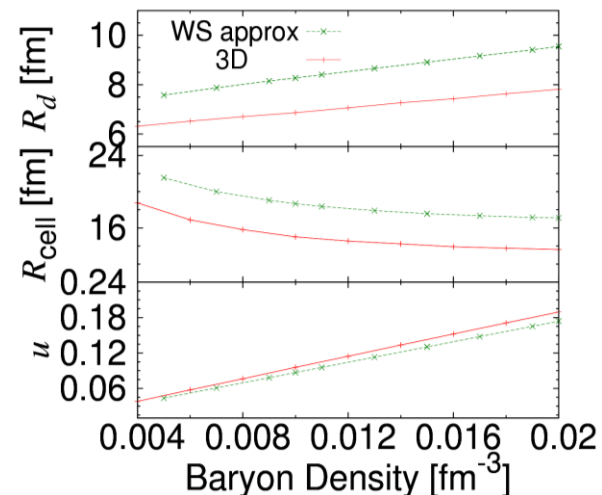
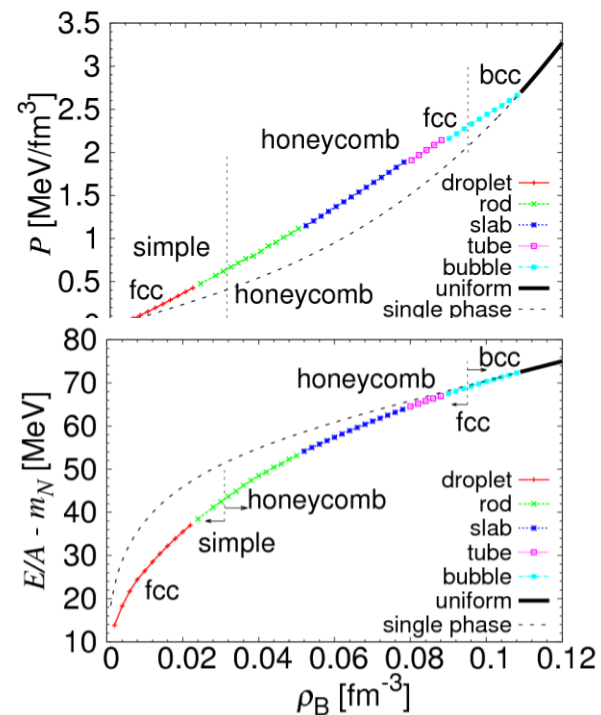
$$Y_p=N/A=0.5 \text{ (} A: \text{ total mass, } N: \text{ proton number)}$$



- We obtained the typical pasta structures as ground states.
- Crystalline of Pasta structure

New results

- Crystalline of droplet : “fcc” is energetically more favorable than “bcc”
- Size of droplet and lattice constant : Smaller than WS cell approximation

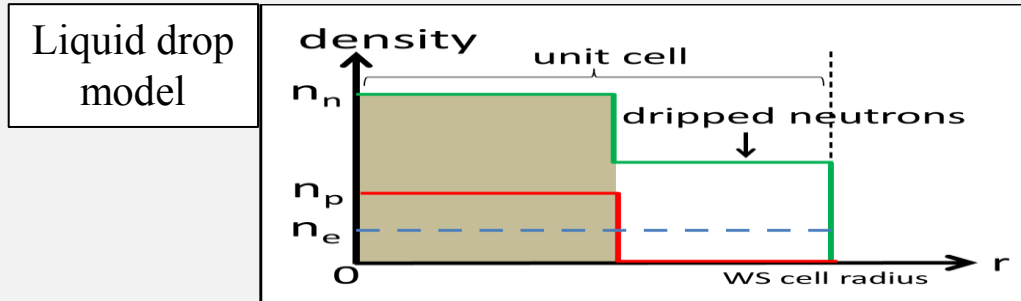


Effect of the different size of droplet

In the previous study of “Pasta structure” ...

D.G. Ravenhall et.al, PRL. 50, 2069 (1983)

K.Oyamatsu et.al, PTP. 72, 2 (1984)



$$\text{Total Energy} = (\text{bulk}) + (\text{Surface}) + (\text{Coulomb})$$

Assuming the same size of droplet for “bcc” and “fcc”



All the energy difference is the Coulomb Energy of “bcc” and “fcc”

In our calculation ...

ρ_B [fm^{-3}]	0.010	0.012	0.014	0.016	0.018
R(fcc) [fm]	6.33	6.67	6.79	7.02	7.31
R(bcc) [fm]	6.88	6.95	7.17	7.45	7.62

Radius of droplet

Energy difference between “fcc” and “bcc”:

0.2 ~ 0.8 MeV/A

The ratio of Coulomb energy difference is 20%



Need to calculate self-consistently including the size of “Pasta structure”

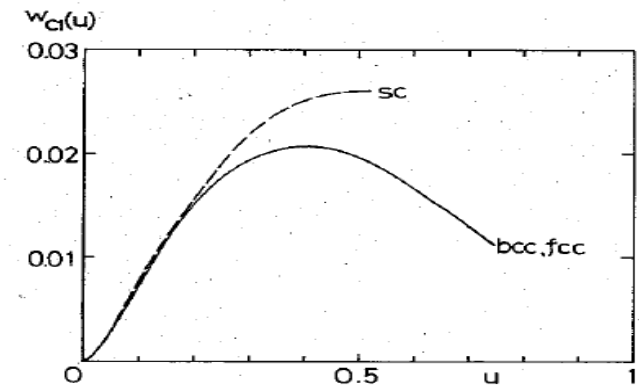
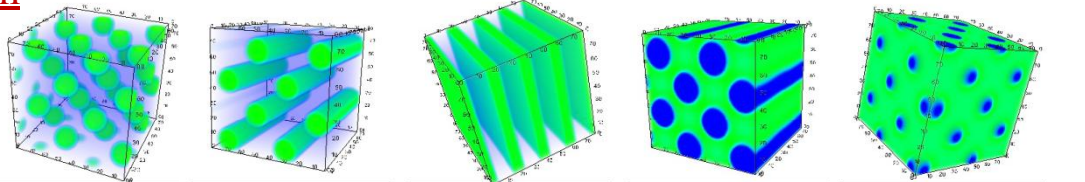


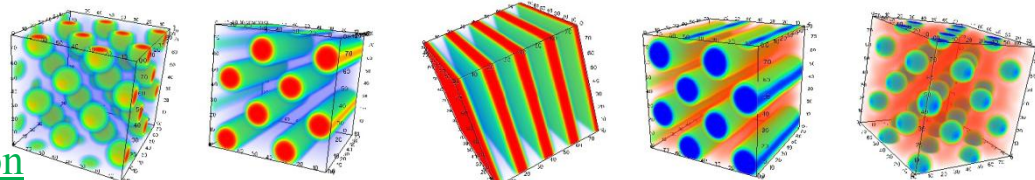
Fig. 1. Relative Coulomb energy $w_{C1}(u)$ for the matter with spherical nuclei. The curves correspond to three lattice types. The curve for bcc is actually lower than the one for fcc only by an indistinguishable amount. The range of u for bcc is $0 \leq u \leq 0.6802$, which is narrower than that for fcc.

$Y_p=0.3$

proton



neutron



“droplet”

“rod”

“slab”

“tube”

“bubble”

[fcc]

[simple]

[simple]

[fcc]

$\rho_B=0.016 \text{ fm}^{-3}$

0.030 fm^{-3}

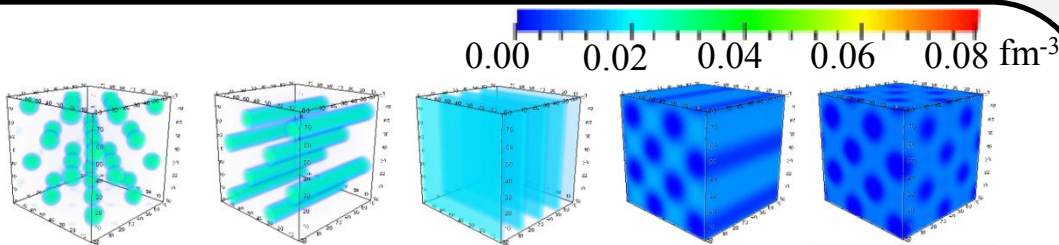
0.05 fm^{-3}

0.066 fm^{-3}

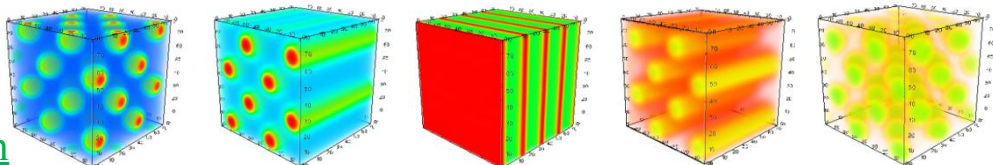
0.080 fm^{-3}

$Y_p=0.1$

proton



neutron



“droplet”

“rod”

“slab”

“tube”

“bubble”

[fcc]

[simple]

[simple]

[fcc]

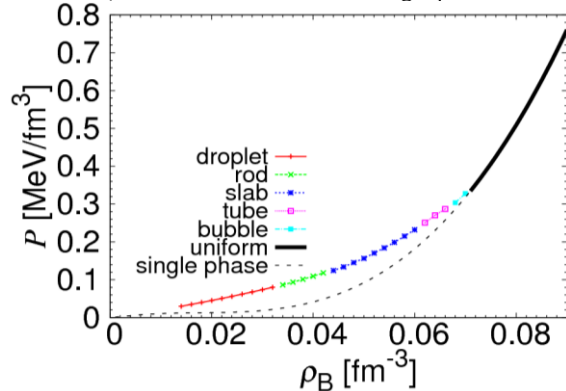
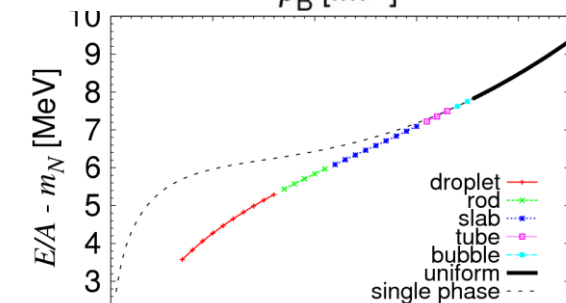
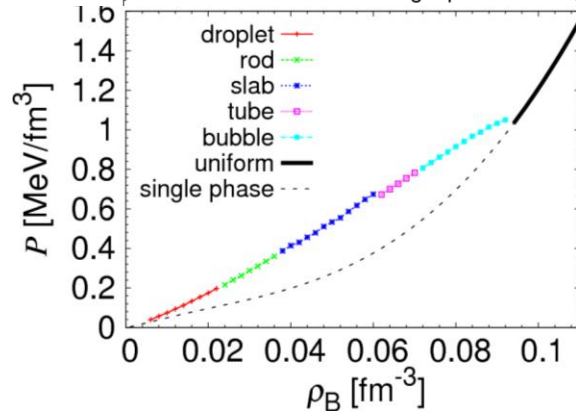
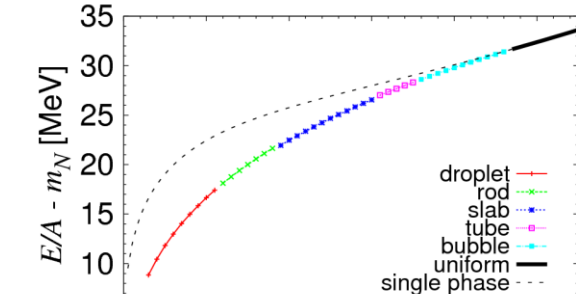
$\rho_B=0.020 \text{ fm}^{-3}$

0.040 fm^{-3}

0.05 fm^{-3}

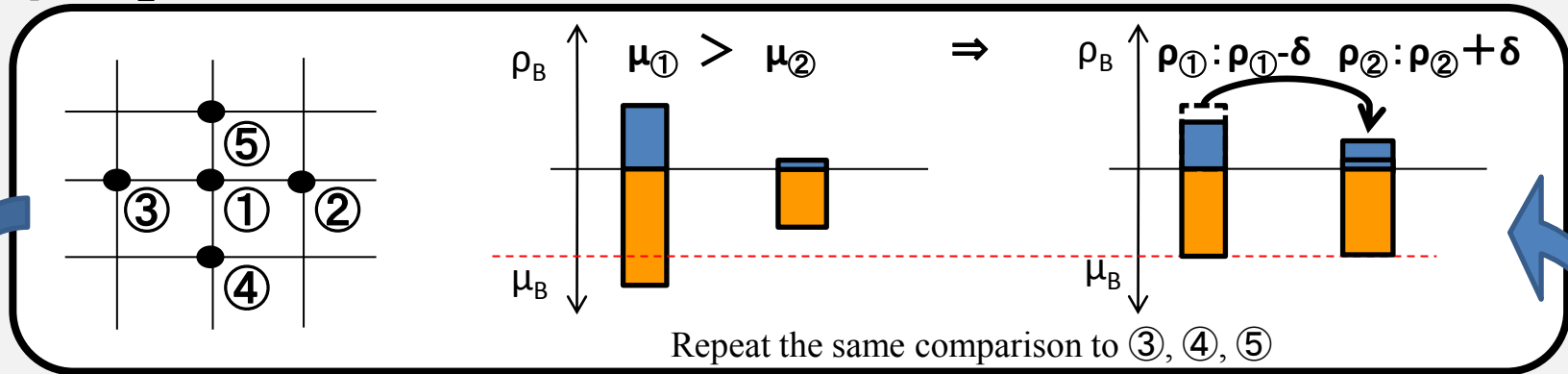
0.066 fm^{-3}

0.070 fm^{-3}

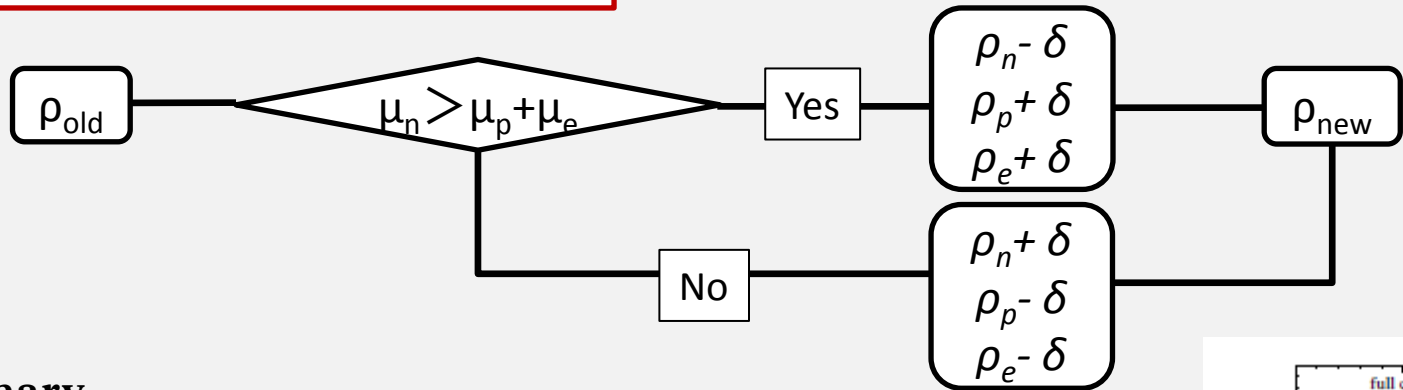


For any proton fraction, typical pasta structures appear as ground states

For β -equilibrium

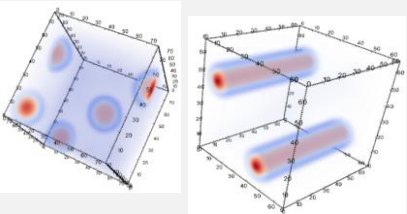


β -equilibrium : $\mu_n(\vec{r}) = \mu_p(\vec{r}) + \mu_e$

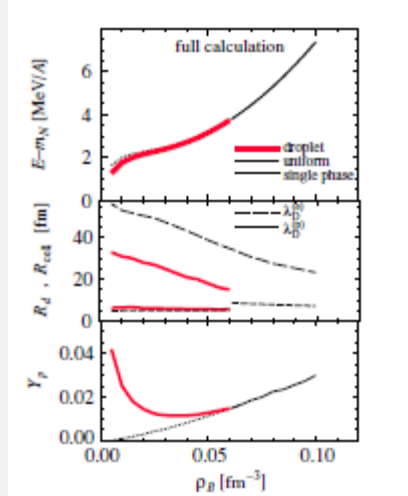


Preliminary

For WS cell approximation and small cell calculation, droplet only appear as ground state



- “droplet” and “rod” appear as ground states
- The comparison of energy between “droplet” and “rod” is not yet carried out.



▪ Summary

We explored the pasta structures of low-density nuclear matter based on RMF without any assumption for the structures.

- In 3D calculation, we obtained the typical pasta structures as ground state
- We explored the crystalline structure of pasta
- "fcc" is energetically more favorable than "bcc" for droplet.
⇒ Difference of droplet's size for "bcc" and "fcc"
- In β -equilibrium, we obtained "droplet" & "rod" structure as ground state.

▪ Future plans

- Explore of the EOS of β -equilibrium nuclear matter in detail
- Calculation of Shear modulus in the crust region of Neutron star
- Extension to the higher density and finite temperature nuclear matter
- Comparing with QMD calculation result for supernova compression process and local minimum states of our calculation