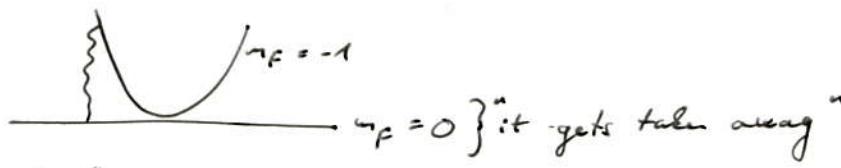


- using RF-resonance can put atoms to different m_F

↳ most energetic ones can be removed



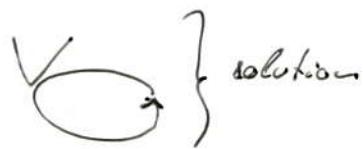
↳ "evaporation"

- modulus potential: spontaneous magnetization flip: $m_F = -1 \rightarrow m_F = 0$
→ problem...

→ using HO potential

→ apply rotating pot. on the modulus

⇒ the effective pot will be HO.



} solution

[2019.03.04.]

- $T = T_c$ (non-int., homogeneous sys)

$$N = V \underbrace{\left(\frac{m \varepsilon_0 T}{2\pi \hbar^2} \right)^{3/2}}_{\lambda_{dB}^{-3}(T_c) \text{ no thermal deBroglie wavelength}} \cdot \zeta(3/2)$$

$\lambda_{dB}^{-3}(T_c)$ no thermal deBroglie wavelength
 $\zeta(3/2)$ no characteristic length

$$\varepsilon_0 T = \frac{2\pi}{m} \frac{\hbar^2}{\lambda_{dB}^2(T)} \text{ no this defines it. (Kinetic energy)}$$

$$\boxed{\zeta(3/2) = n \lambda_{dB}^3(T_c)} \text{ critical combination}$$

phase-space density

this is the good check for BEC

- ζ func. has 0-s at even Θ numbers / Mathematica!

• explaining the deviation on the $N_0/N - T_c/T$ plot:

$$\boxed{T > T_c}$$

$$N = \left(\frac{\varepsilon_0 T}{\hbar \omega} \right)^3 F_-(3, \frac{\varepsilon_0 - \mu}{\varepsilon_0 T}) + \delta_1 \left(\frac{\varepsilon_0 T}{\hbar \omega} \right)^2 F_-(2, \frac{\varepsilon_0 - \mu}{\varepsilon_0 T})$$

$$\boxed{T < T_c}$$

(new stuff)

- big degeneracy on ground state $\rightarrow N_0$

- $M = E_0 \rightarrow 0\text{-s in } F_-!$

$$N = N_0 + \left(\frac{\epsilon_0 T}{\hbar \omega}\right)^3 \zeta(3) + \gamma \left(\frac{\epsilon_0 T}{\hbar \omega}\right)^2 \zeta(2)$$

$$1 - \frac{N_0}{N} = \left(\frac{\epsilon_0 T}{\hbar \omega}\right)^3 \frac{\zeta(3)}{N} + \gamma_1 \left(\frac{\epsilon_0 T}{\hbar \omega}\right)^2 \frac{\zeta(2)}{N}$$

$$\frac{N_0}{N} = 1 - \underbrace{\left(\frac{T}{T_0}\right) \left(\frac{\epsilon_0 T_0}{\hbar \omega}\right)^3 \frac{\zeta(3)}{N}}_1 - \gamma_1 \left(\frac{T}{T_0}\right)^2 \left(\frac{\epsilon_0 T_0}{\hbar \omega}\right)^2 \frac{\zeta(2)}{N}$$

$\left(\frac{N}{\zeta(3)}\right)^{2/3}$
 explains
why it is
underneath

$\sim N^{-1/3}$ behaviour
 \rightarrow finite size corr.
 \rightarrow gets smaller
 with $N \rightarrow \infty$

-
- depending on ω -s: $\omega_z < \omega_r ; \omega_z > \omega_r$
 - disc type cigar type condensate
 - oscillator length: $d = \sqrt{\frac{\hbar}{m\omega}} \sim$ depends on direction
 \sim usual condensate size
 - What is the half-width of C, if the non-int. modell is correct?

- ground state wf:

$$\varphi_0(\vec{r}_1) \varphi_0(\vec{r}_2) \dots \varphi_0(\vec{r}_N) = \Psi(\vec{r}_1, \dots, \vec{r}_N)$$

$$\hat{n}(\vec{r}) = \sum_{i=1}^N \delta(\vec{r} - \vec{r}_i) \rightarrow \text{density } n \text{ in 1st quantization.}$$

↓

$$n(r) = \langle \Psi | \hat{n} | \Psi \rangle \rightarrow \text{density func.!}$$

$$\left(\int (\varphi_0(r))^2 d^3r = 1 \right)$$

$$n(r) = \prod_i \int d^3r_1 \int d^3r_2 \dots \int d^3r_N \varphi_0^*(r_1) \dots \varphi_0^*(r_N) \delta(r - r_i) \varphi_0(r_1) \dots \varphi_0(r_N)$$

$$= \prod_{i=1}^N |\varphi_0(r)|^2 = \underbrace{N |\varphi_0(r)|^2}_{\rightarrow} = n(r)$$

if we know the ground state non-int. wf.

- non-int. modell for harmonically trapped atoms:

$$\hat{H} = \left(-\frac{\hbar^2}{2m} \Delta + V(r) \right)$$

↓

$$\frac{1}{2} m (\omega_x^2 x_1^2 + \omega_y^2 x_2^2 + \omega_z^2 x_3^2)$$

$$\Rightarrow \varphi_0(r) = \left(\frac{m\omega}{\pi\hbar} \right)^{3/2} e^{-\frac{1}{2} \left(\frac{x_1}{d_x} \right)^2 - \frac{1}{2} \left(\frac{x_2}{d_y} \right)^2 - \frac{1}{2} \left(\frac{x_3}{d_z} \right)^2}$$

$$\text{and } d_i = \sqrt{\frac{\hbar}{m\omega_i}}$$

- total density is same φ_0 and $n(r)$



half width is in the order of d not few micrometers

- however the measured half width is in the order of 100 μm
 \Rightarrow INTERACTION is NON-NEGIGIBLE

Gross-Pitaevskii - eq. inclusion of the interaction

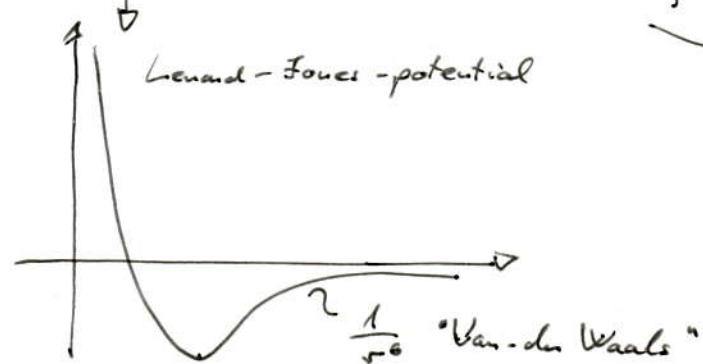
H_{ext} + H_{int}

spinless boson

- wf for rotating BEC

external, HO potential

$$\hat{H} = \sum_{i=1}^N \left(-\frac{\hbar^2}{2m} \Delta_i + V_{\text{ext}}(\vec{r}_i) \right) + \frac{1}{2} \sum_{i \neq j=1}^N v(\vec{r}_i - \vec{r}_j)$$



$$v \rightarrow v(|\vec{r}|)$$

• it acts weakly

- we can still use product ansatz for the whole wf.

$$\Psi(\vec{r}_1, \vec{r}_2, \dots, \vec{r}_N) = \varphi(\vec{r}_1) \cdot \dots \cdot \varphi(\vec{r}_N)$$

→ constraint: $1 = \int |\varphi|^2 d^3r$ → Ψ is normalized, too.

- Minimize:

$$\frac{\delta}{\delta \varphi^*(r)} \left(\langle \Psi | \hat{H} | \Psi \rangle - E \int d^3r |\varphi(r)|^2 \right) \stackrel{!}{=} 0$$

Lagrange-multiplicator...

(nearly similar to Hartree-Fock, but we only have φ , not a Slater...)

$$\begin{aligned} \langle \Psi | \hat{H} | \Psi \rangle &= \sum_{i=1}^N \int d^3r_1 \dots d^3r_N \varphi^*(r_i) \dots \varphi^*(r_N) \left(-\frac{\hbar^2}{2m} \Delta_i + V(r_i) \right) \varphi(r_i) \dots \varphi(r_N) \\ &\quad + \frac{1}{2} \sum_{i \neq j} \int d^3r_1 \dots d^3r_N \varphi^*(r_i) \dots \varphi^*(r_N) v(\vec{r}_i - \vec{r}_j) \varphi(r_i) \dots \varphi(r_N) \end{aligned}$$

$$= \sum_{i=1}^N \int d^3 r_i \varphi^*(\vec{r}_i) \left(-\frac{\hbar^2}{2m} \Delta_i + V(\vec{r}_i) \right) \varphi(\vec{r}_i) + \frac{1}{2} \sum_{i,j=1, i \neq j}^N \int d^3 r_i d^3 r_j \varphi^*(\vec{r}_i) \varphi^*(\vec{r}_j).$$

19.

$$\cdot V(r_i - r_j) \varphi(r_i) \varphi(r_j) =$$

$r_i = r'$
 $r_j = r''$

change variable \Rightarrow summations are doable ...

$$= N \cdot \int d^3 r' \varphi^*(r') \left(-\frac{\hbar^2}{2m} \Delta' + V(r') \right) \varphi(r') + \frac{N(N-1)}{2} \int d^3 r' d^3 r'' \varphi^*(r') \varphi^*(r'').$$

$$\cdot V(r' - r'')$$

$$\frac{\delta}{\delta \varphi^*(r)} \left(\langle 4 | \hat{H} | 4 \rangle - E \int d^3 r' \varphi^*(r') \varphi(r') \right) =$$

$$\left[\frac{\delta \varphi(r)}{\delta \varphi^*(r)} = 0 ; \frac{\delta \varphi^*(r')}{\delta \varphi^*(r)} = \delta(r - r') \right]$$

\Rightarrow can be used for integration.

$$= N \int d^3 r' \delta(r - r') \left(-\frac{\hbar^2}{2m} \Delta' + V(r') \right) \varphi(r') +$$

$$+ \frac{N(N-1)}{2} \int d^3 r' d^3 r'' \left(\delta(r - r') \varphi^*(r'') + \delta(r - r'') \varphi^*(r') \right) V(r' - r'') \varphi(r') \varphi(r'')$$

$$- E \int d^3 r' \delta(r - r') \varphi(r') =$$

symmetric... ($V(r')$)

$$= N \left(-\frac{\hbar^2}{2m} \Delta + V(r) \right) \varphi(r) + N(N-1) \int d^3 r' \varphi^*(r') V(r' - r) \varphi(r') \varphi(r) -$$

$$- E \varphi(r)$$

$$\mathcal{O} = N \left(-\frac{\hbar^2}{2m} \Delta + V(r) \right) \varphi(r) + N(N-1) \left(\int d^3r' \varphi^*(r') V(r-r') \varphi(r') \right) \varphi(r) \sim E \varphi(r)$$

\rightsquigarrow non-linear - eq.

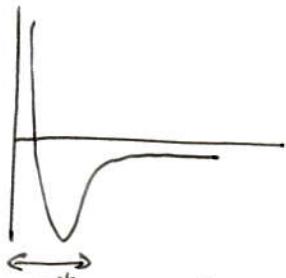
$$\Psi_0(r) = N^{1/2} \varphi(r) \rightsquigarrow \text{condensate - wf.}$$

$$\int |\Psi_0(r)|^2 d^3r = N \rightsquigarrow n_c(r) = |\Psi_0(r)|^2 \text{ condensate - density}$$

$$\mu = \frac{E}{N} \rightsquigarrow \text{chemical potential}$$

$$\mu \Psi_0(r) = \left(-\frac{\hbar^2}{2m} \Delta + V(r) \right) \Psi_0(r) + \underbrace{\frac{N-1}{N}}_{\rightarrow 1 \text{ if } N \rightarrow \infty} \varphi(r) \int d^3r' V(r-r') |\varphi(r')|^2$$

\rightsquigarrow non-local form of Gross-Pitaevski-equation



R^* \rightsquigarrow char. length of pot \rightsquigarrow if the $\sqrt{R} \gg R^*$ dist. of atoms.

and $T \rightarrow \infty$ and big density

$V(r-r')$ can be replaced with: $\frac{4\pi \hbar^2 a}{m} \delta(r-r')$

where a is the s-wave scattering length.

Then:

$$\boxed{\mu \Psi_0(r) = \left(-\frac{\hbar^2}{2m} \Delta + V(r) \right) + \frac{4\pi \hbar^2 a}{m} |\Psi_0(r)|^2 \Psi_0(r)}$$