

$$\begin{aligned}
 (\omega + i\eta - H_0 - V)G^R &= 1 \\
 (G_0^R - V)G^R &= 1 \\
 G_0^R G^R &= 1 + VG^R \\
 \boxed{G^R = G_0^R + G_0^R V G^R} & \text{ L.S.} \\
 \langle r | G^R | r' \rangle &= \langle r | G_0^R | r' \rangle + \int \langle r | G_0^R | r'' \rangle V(r'') \langle r'' | G^R | r' \rangle dr'' \\
 \boxed{G^R = G_0^R + G_0^R \sum^R G^R} & \\
 & \leftarrow \text{Self-énergie irréductible}
 \end{aligned}$$

$$G^R = \frac{1}{G_0^{-1R} - \Sigma^R} = \frac{1}{\omega + i\eta - \frac{\hbar^2}{2m} - \underbrace{\left(\frac{b_0}{\omega + i\eta - a_0 - \frac{b_1}{\omega + i\eta - a_1 - \dots}} \right)}$$

↑ Pôles déplacés.

$$\langle r | G^R | r' \rangle = \sum_n \frac{\varphi_n(r) \varphi_n^*(r')}{\omega + i\eta - E_n}$$

$$\Sigma^R = a - \frac{i}{\tau}$$

Moyenne sur impuretés:

$$N_c(q) = \sum_{i=1}^{N_i} e^{-iq \cdot R_i} N(q)$$

$$\rightarrow \overline{N_c(q)} = N(q) (2\pi)^3 \delta(q) \eta_i$$

$$\begin{aligned} \overline{N_c(q) N_c(q')} &= \eta_i N(q) N(-q) \delta^3(q+q') (2\pi)^3 \\ &\quad \uparrow \quad \uparrow \\ &\quad \leftarrow \quad \downarrow \\ &= \eta_i |N(q)|^2 \delta^3(q+q') (2\pi)^3 \end{aligned}$$

$$\Sigma^R = \begin{array}{c} \times \\ \vdots \\ \vdots \end{array} + \begin{array}{c} \times \\ \vdots \quad \vdots \\ \hline \xrightarrow{\quad} \\ \vdots \quad \vdots \\ \times \quad \times \end{array} + \dots$$

$h_1 \neq h_2$

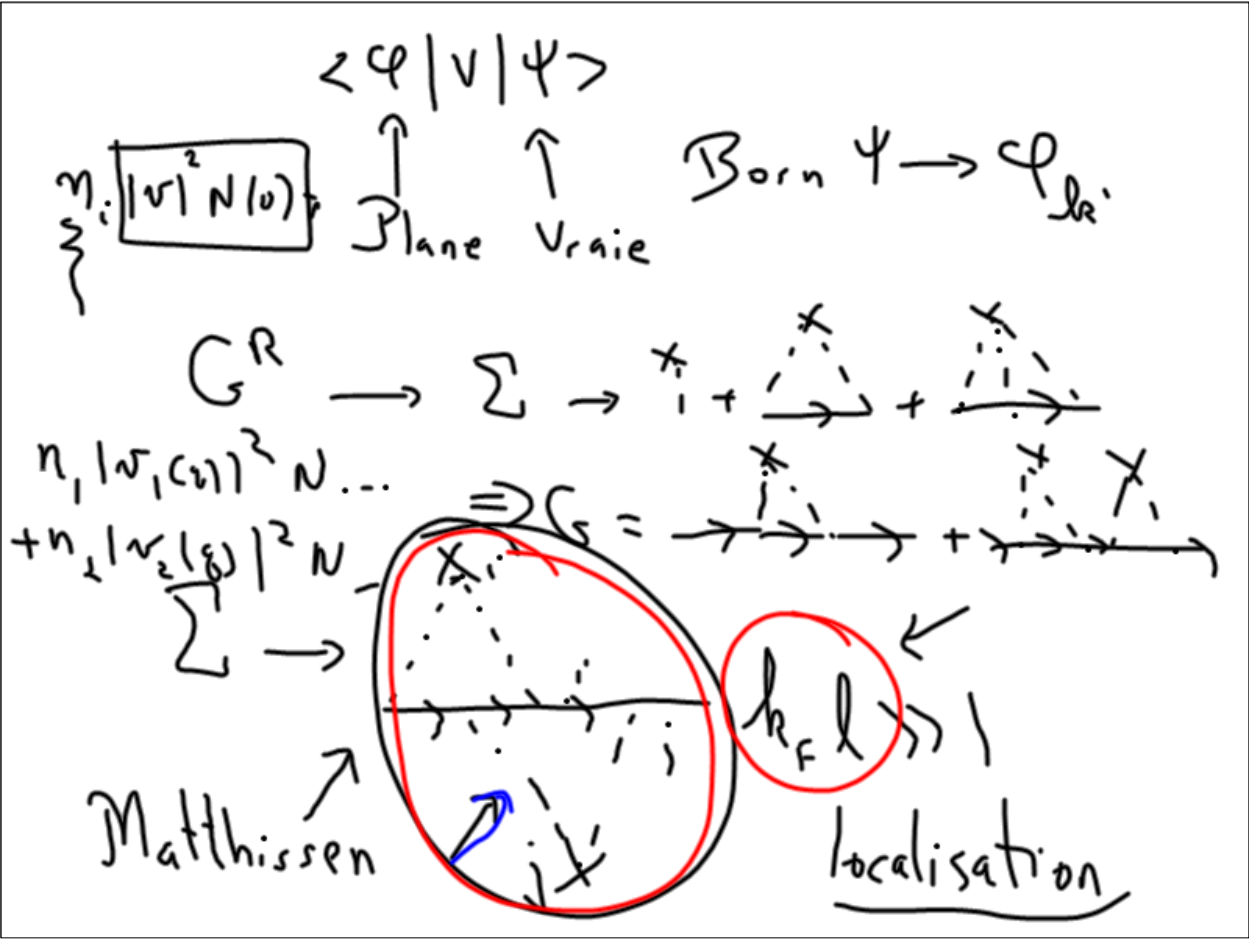
$$= n_i N(\Omega) + n_i \int_{\substack{d^3 h_i \\ (2\pi)^3}} |v(h_i)|^2 G_o^R(h_i, \omega)$$

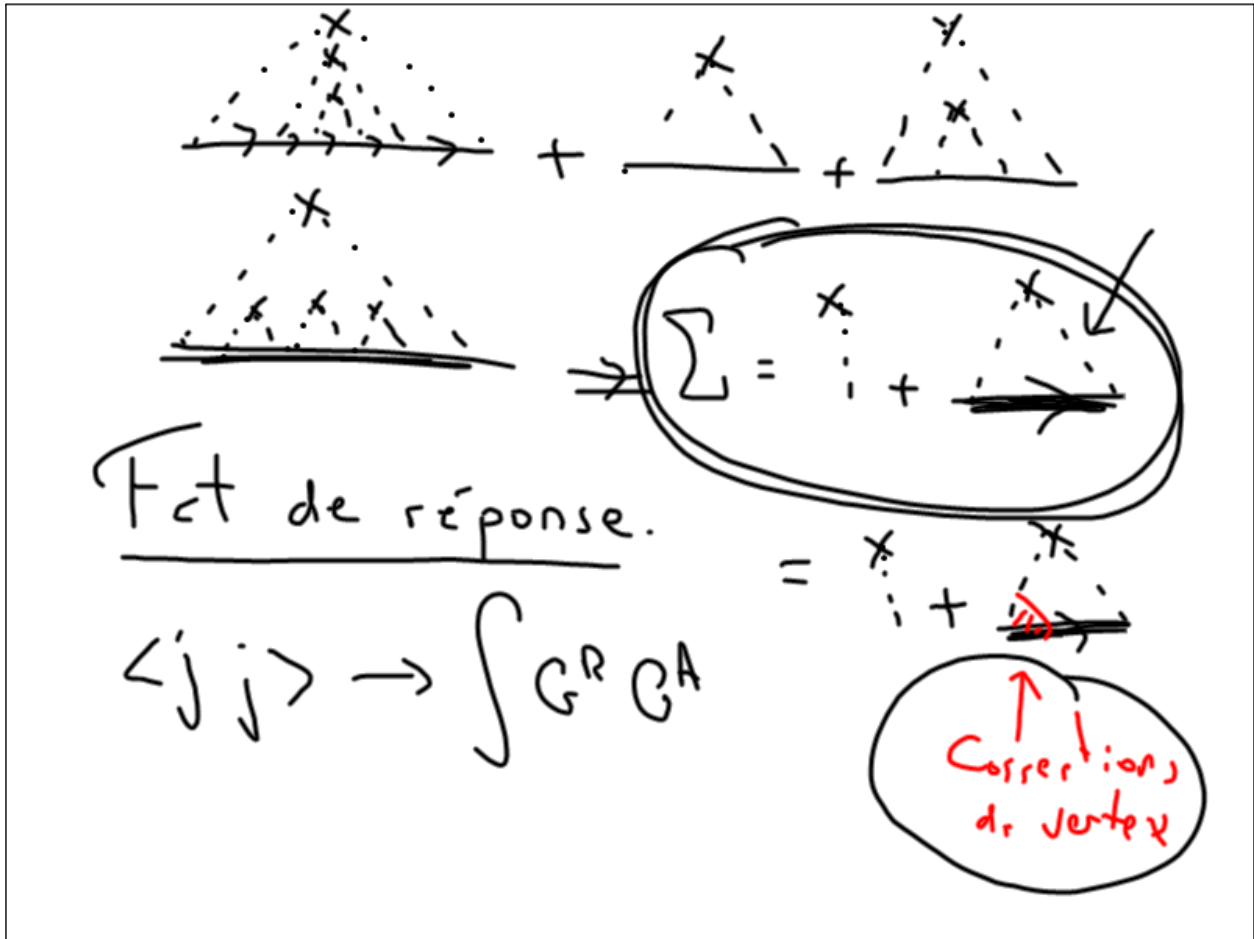
$h_1 \neq h_2$

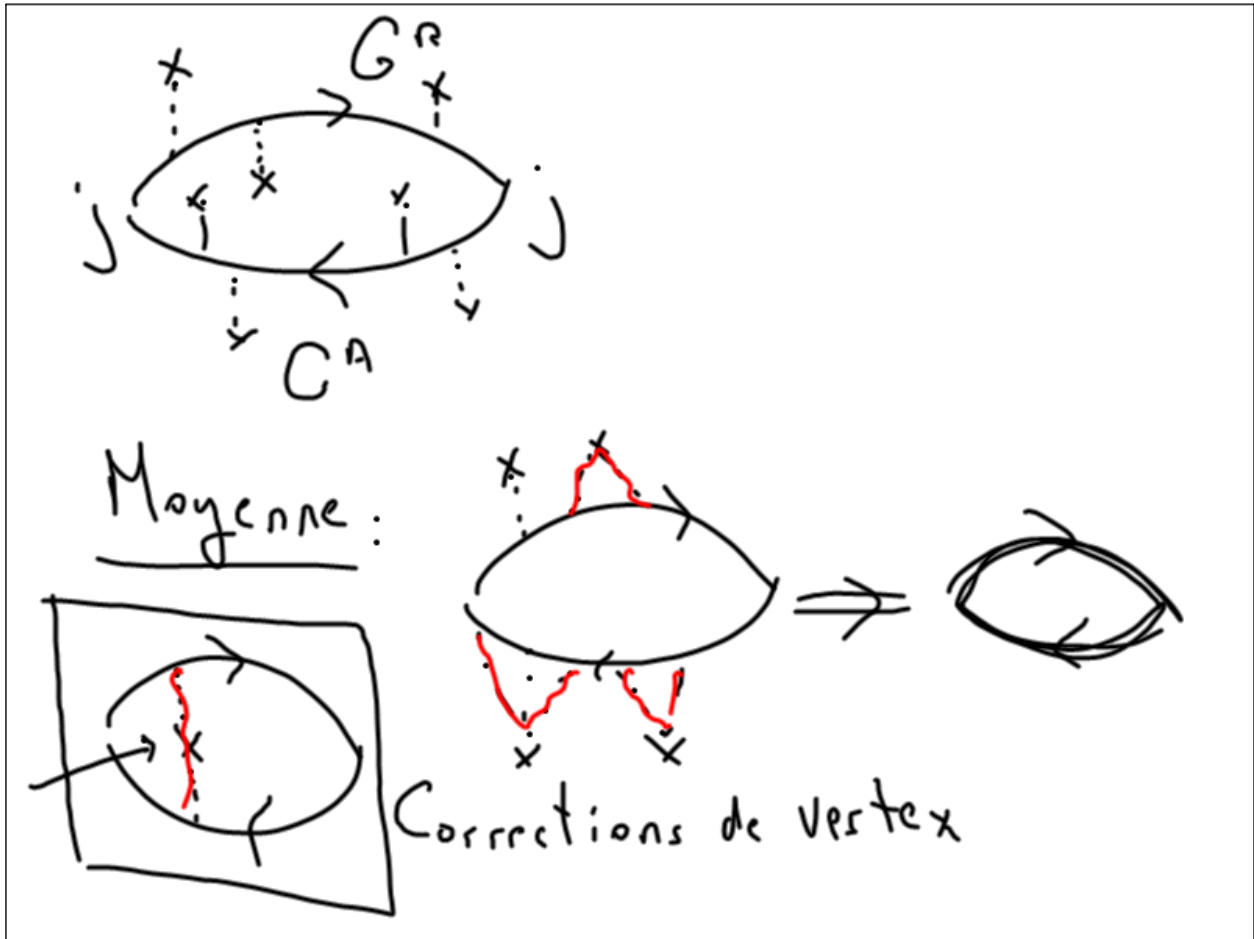
$$\Sigma = \begin{array}{c} \times \\ \vdots \\ \vdots \end{array} + \begin{array}{c} \times \\ \vdots \\ \vdots \\ \hline \xrightarrow{\quad} \\ \vdots \\ \times \end{array} + \begin{array}{c} \times \\ \vdots \\ \vdots \\ \hline \xrightarrow{\quad} \\ \vdots \\ \times \end{array} + \begin{array}{c} \times \\ \vdots \\ \vdots \\ \hline \xrightarrow{\quad} \\ \vdots \\ \times \end{array} + \dots$$

Calcul de σ
au delà de
Born

Diffuse plusieurs fois
sur même impulsion

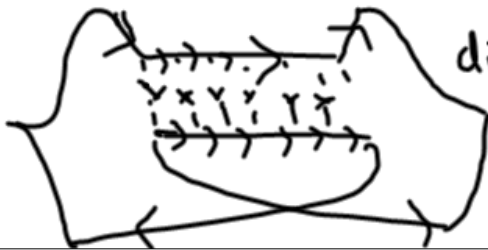






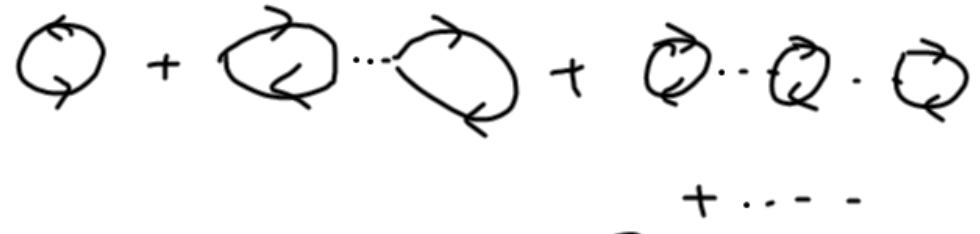
Pour satisfaire les lois de conservation il faut choisir les corrections de vertex de telle sorte qu'elles soient cohérentes avec la self-énergie retenue. Formel: Identités de Ward.

Corrections de vertex



diagrammes en échelle
(approx. de matière T)

Coulomb



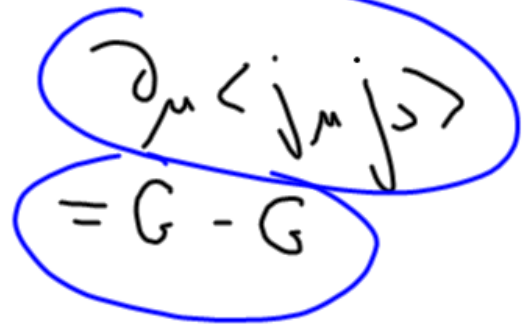
Bulles

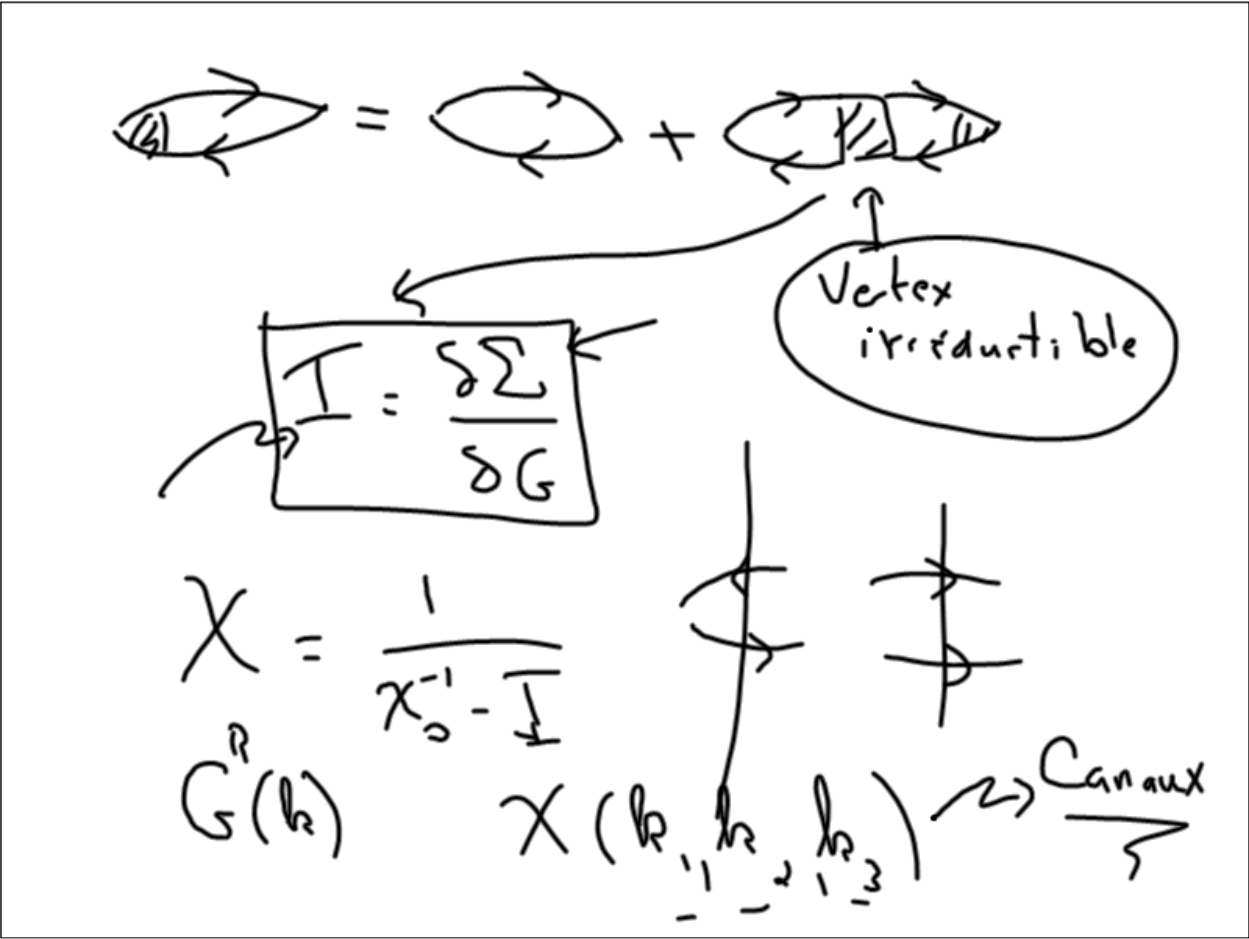
Échelles

Parquet

Bulles

Word





$$\langle \psi^\dagger \psi \psi^\dagger \psi \rangle$$

$$I_1, I_2, I_3$$

{ Relations de croisement
Crossing symmetry

Chapitre 4

4.1 - Seconde quantification

- 1 corps

- 2 corps

- Repr. Heisenberg

4.2. Motivation pour drif. de G^R

- Exemple H quadratique.

4.3 - Repr. int.

4.4 - K.B + K.S. (contours)

4.5 - Matsubara.

4.6 - Signif. physique. 4.7 Théorèmes.

$$\Psi(x_1, x_2, x_3) = \pm \Psi(x_2, x_1, x_3)$$

indiscernables

Symétriques: bosons

Antisymétriques: fermions.

Base à 1 particule.

$$A_1 |\alpha_1\rangle |\alpha_2\rangle |\alpha_3\rangle$$

$$+ A_2 |\beta_1\rangle |\beta_2\rangle |\beta_3\rangle$$

$$\Psi(x, y) = (x-y) N e^{-\frac{|x-y|}{a}}$$

✓

Seconde quantification (Neyele-Orland)

$|\alpha_1, \alpha_2, \dots, \alpha_n\rangle$ pas normalisé

$$|\alpha_1, \dots, \alpha_n\rangle = \frac{1}{\sqrt{\pi^{n!} \alpha_1! \dots \alpha_n!}} |\alpha_1, \alpha_2, \dots, \alpha_n\rangle$$

↙
↑
états à 1 corps.

Normalisé et (anti)symétrisé

$$|\alpha_1, \dots, \alpha_n\rangle = a_{\alpha_1}^+ a_{\alpha_2}^+ \dots a_{\alpha_n}^+ |0\rangle$$

$a_{\alpha_1}^+$

rajoute (crée)
une particule dans l'état α_1
et symétrise
↯
(ou antisymétrise) avec
les autres particules.

Exemple:

$$\langle \vec{r}_1 | \langle \vec{r}_2 | a_{\alpha_1}^+ a_{\alpha_2}^+ | 0 \rangle = \varphi_{\alpha_1, \alpha_2}(\vec{r}_1, \vec{r}_2)$$

Fermions:

$\xi = \text{zeta}$

$$a_\alpha^+ a_\beta^+ - \xi a_\beta^+ a_\alpha^+ = 0 \quad \xi \quad \xi$$

$$a_\alpha a_\beta - \xi a_\beta a_\alpha = 0$$

$$a_\alpha a_\beta^+ - \xi a_\beta^+ a_\alpha = \delta_{\alpha\beta}$$

$\underbrace{\hspace{1.5cm}} \quad \underbrace{\hspace{1.5cm}}$

on $\xi = -1$ fermions,
+1 bosons

Formule de changement de base

$$a_{\alpha}^{\dagger} = \sum_{\mu} c_{\mu}^{\dagger} \langle \mu | \alpha \rangle$$

$$|\alpha\rangle = \sum_{\mu} |\mu\rangle \langle \mu | \alpha \rangle$$

$$a_{\alpha} = \sum_{\mu} \langle \alpha | \mu \rangle c_{\mu}$$

Opérateur nombre

$$\hat{n}_\alpha = a_\alpha^\dagger a_\alpha$$

$$a_\alpha |0\rangle = 0$$

$$\hat{n}_\alpha (a_\alpha^{+2} |0\rangle)$$

$$= a_\alpha^\dagger a_\alpha a_\alpha^{+2} |0\rangle$$

$$= a_\alpha^\dagger (1 + a_\alpha^\dagger a_\alpha) a_\alpha |0\rangle = 2 (a_\alpha^{+2} |0\rangle)$$

$$\begin{aligned} n_\beta a_\alpha^\dagger |0\rangle &= 0 \\ a_\beta^\dagger a_\beta a_\alpha^\dagger |0\rangle &= 0 \end{aligned}$$

Opérateurs à l'op.

Dans la base propre

$$\hat{U}|\alpha\rangle = U_\alpha|\alpha\rangle = \langle\alpha|U|\alpha\rangle|\alpha\rangle$$

↑
Valeur propre.

$$\hat{U}|\alpha_1, \dots, \alpha_n\rangle = \sum_{i=1}^n U_{\alpha_i}|\alpha_1, \dots, \alpha_n\rangle$$

$$= \sum_{\alpha} U_{\alpha} \hat{n}_{\alpha} |\alpha_1, \dots, \alpha_n\rangle$$

$$\begin{aligned}
\hat{U} &= \sum_{\alpha} \hat{n}_{\alpha} U_{\alpha} \\
&= \sum_{\alpha \alpha'} \langle \alpha | U | \alpha' \rangle a_{\alpha}^{\dagger} a_{\alpha'} \\
&= \sum_{\alpha \alpha'} \sum_{\lambda \mu} C_{\mu}^{\dagger} \langle \mu | \alpha \rangle \langle \alpha | U | \alpha' \rangle \\
&= \sum_{\lambda} C_{\lambda}^{\dagger} \langle \lambda | U | \lambda \rangle C_{\lambda}
\end{aligned}$$

Base utilisée souvent

$$\sum_{\mathbf{k}} |\mathbf{k}\rangle \langle \mathbf{k}| = 1 = \int d\mathbf{r} |\mathbf{r}\rangle \langle \mathbf{r}|$$

$$- \langle \mathbf{r} | \mathbf{k} \rangle = \frac{1}{\sqrt{V}} e^{i\mathbf{k} \cdot \mathbf{r}}$$

$$- \langle \mathbf{k} | \mathbf{r} \rangle = \frac{1}{\sqrt{V}} e^{-i\mathbf{k} \cdot \mathbf{r}}$$

$$\begin{aligned} \langle \mathbf{r} | \mathbf{r}' \rangle &= \sum_{\mathbf{k}} \langle \mathbf{r} | \mathbf{k} \rangle \langle \mathbf{k} | \mathbf{r}' \rangle \\ &= \frac{1}{V} \sum_{\mathbf{k}} e^{i\mathbf{k}(\mathbf{r}-\mathbf{r}')} = \int \frac{d^3\mathbf{k}}{(2\pi)^3} e^{i\mathbf{k} \cdot (\mathbf{r}-\mathbf{r}')} = \delta(\mathbf{r}-\mathbf{r}') \end{aligned}$$

$$\begin{aligned} \langle \mathbf{k} | \mathbf{k}' \rangle &= \int d^3\mathbf{r} \langle \mathbf{k} | \mathbf{r} \rangle \langle \mathbf{r} | \mathbf{k}' \rangle \\ &= \frac{1}{V} \int d^3\mathbf{r} e^{-i(\mathbf{k}-\mathbf{k}') \cdot \mathbf{r}} = \delta_{\mathbf{k}\mathbf{k}'} \end{aligned}$$

$$\Psi(r) = \sum_{\mathbf{h}} \langle r | \mathbf{h} \rangle c_{\mathbf{h}}$$

$$\Rightarrow \{c_{\mathbf{h}}, c_{\mathbf{h}'}^\dagger\} = \delta_{\mathbf{h}\mathbf{h}'}$$

$$\{\Psi(r), \Psi^\dagger(r')\} = \delta^3(r-r')$$

$$\begin{aligned} \{\Psi(r), \Psi^\dagger(r')\} &= \sum_{\mathbf{h}\mathbf{h}'} \langle r | \mathbf{h} \rangle \underbrace{\{c_{\mathbf{h}}, c_{\mathbf{h}'}^\dagger\}}_{\delta_{\mathbf{h}\mathbf{h}'}} \langle \mathbf{h}' | r' \rangle \\ &= \sum_{\mathbf{h}} \langle r | \mathbf{h} \rangle \langle \mathbf{h} | r' \rangle = \langle r | r' \rangle = \delta^3(r-r') \end{aligned}$$

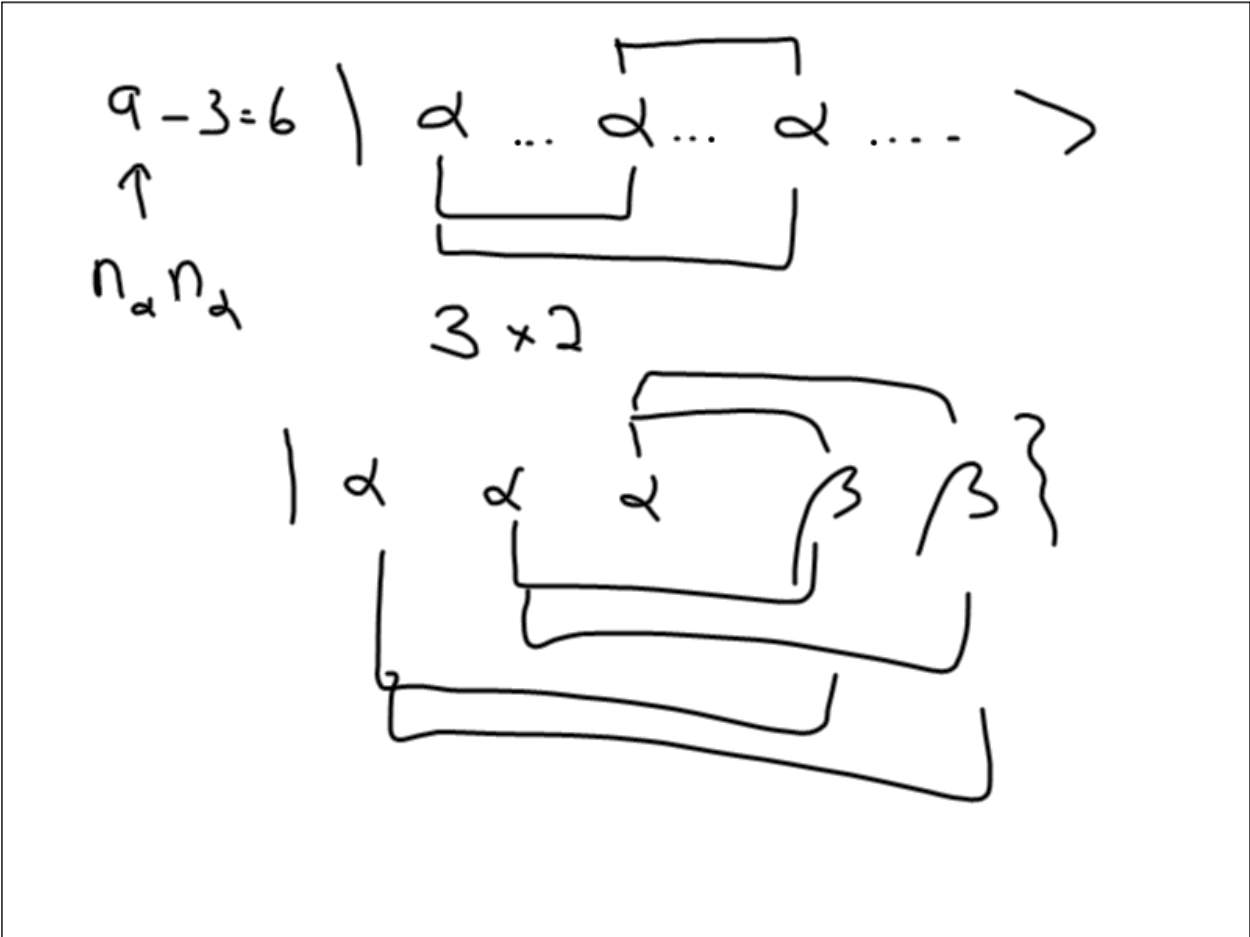
$$\begin{aligned}
 \hat{T} &= \sum_{\mathbf{k}} \langle \mathbf{k} | \frac{\hbar^2}{2m} | \mathbf{k} \rangle c_{\mathbf{k}}^{\dagger} c_{\mathbf{k}} \\
 &= -\frac{1}{2m} \int d^3r \psi^{\dagger}(\mathbf{r}) (\nabla^2 \psi(\mathbf{r})) \\
 &= \frac{1}{2m} \int d^3r (\nabla \psi^{\dagger}(\mathbf{r})) \cdot (\nabla \psi(\mathbf{r}))
 \end{aligned}$$

Opérateurs à 2 corps

$$\hat{V} |\alpha\rangle \otimes |\beta\rangle = V_{\alpha\beta} |\alpha\rangle \otimes |\beta\rangle$$

Base propre.

$$\begin{aligned} \hat{V} |\alpha_1, \dots, \alpha_n\rangle &= \frac{1}{2} \sum_{i=1}^n \sum_{j \neq i}^n V_{\alpha_i, \alpha_j} |\alpha_1, \dots, \alpha_n\rangle \\ &= \frac{1}{2} \sum_{\alpha\beta} V_{\alpha\beta} (\hat{n}_\alpha \hat{n}_\beta - \delta_{\alpha\beta} \hat{n}_\alpha) |\alpha_1, \dots, \alpha_n\rangle \end{aligned}$$



$$\begin{array}{cc} n_\alpha & n_\beta \\ \uparrow & \uparrow \\ 3 & 2 \end{array}$$

$$\begin{aligned} & \hat{n}_\alpha \hat{n}_\beta - \delta_{\alpha\beta} \hat{n}_\alpha \\ &= a_\alpha^\dagger a_\alpha a_\beta^\dagger a_\beta - \delta_{\alpha\beta} a_\alpha^\dagger a_\alpha \\ &= a_\alpha^\dagger (\cancel{\delta_{\alpha\beta}} - \int a_\beta^\dagger a_\alpha) a_\beta - \delta_{\alpha\beta} \cancel{a_\alpha^\dagger} a_\alpha \\ &= - \int a_\alpha^\dagger a_\beta^\dagger a_\alpha a_\beta = a_\alpha^\dagger a_\beta^\dagger a_\beta a_\alpha \leftarrow \end{aligned}$$

Destruction à droite
de création par def. =

Ordre normal

$$\hat{V} = \frac{1}{2} \int d^3x d^3y v(x-y) \psi^\dagger(x) \psi^\dagger(y) \psi(y) \psi(x)$$

Heisenberg

$$\begin{aligned} \rightarrow c_h(t) &= e^{i\hat{H}t} c_h e^{-i\hat{H}t} \\ c_h^\dagger(t) &= e^{i\hat{H}t} c_h^\dagger e^{-i\hat{H}t} \end{aligned}$$

Cas particulier

$$\hat{H} = \sum_h \epsilon_h c_h^\dagger c_h$$

$$i \frac{\partial}{\partial t} c_h(t) = - [\hat{H}, c_h(t)] \leftarrow$$

$$\begin{aligned} \hat{H}(t) &= e^{i\hat{H}t} \hat{H} e^{-i\hat{H}t} \\ &= \sum_h \epsilon_h c_h^\dagger(t) c_h(t) \end{aligned}$$

$$\begin{aligned} \rightarrow i \frac{\partial}{\partial t} c_h(t) &= - \sum_{h'} \delta_{hh'} \epsilon_{h'} (-) c_{h'} \\ &= \epsilon_h c_h(t) \end{aligned}$$

$$\frac{\partial}{\partial t} c_h(t) = -i \epsilon_h c_h(t)$$

$$\rightarrow \boxed{c_h(t) = e^{-i\epsilon_h t} c_h} \leftarrow$$

$$\begin{aligned}
 [c_{h'}^+, c_{h'}] &= -c_{h'} \delta_{hh'} \\
 [AB, c] &= \overbrace{ABC - ACB} + \overbrace{ACB - CAB} \\
 &= A[B, c] + [A, c]B \\
 &= A\{B, c\} - \{A, c\}B \\
 &= -\delta_{hh'} c_{h'}
 \end{aligned}$$

A curved arrow points from the circled $c_{h'}$ in the first equation to the circled B in the third equation.

