

Many fermions

$$\mathcal{H}_1 = \text{span} \{ |A\rangle \}$$

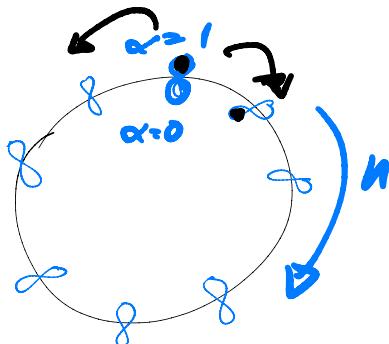
$$\Psi_F(A_1, A_2, \dots, A_n) = -\Psi_F(A_2, A_1, \dots, A_n)$$

\Rightarrow Pauli principle

$$\Psi_F(A, A, \dots) = 0.$$

\Rightarrow Band structure: metals vs insulators

Consider $\mathcal{H}_1 = \text{span} \{ |n\rangle \otimes |\alpha\rangle \}_{n=1..N, \alpha=0,1} \}$



$$|n+N\rangle = |n\rangle$$

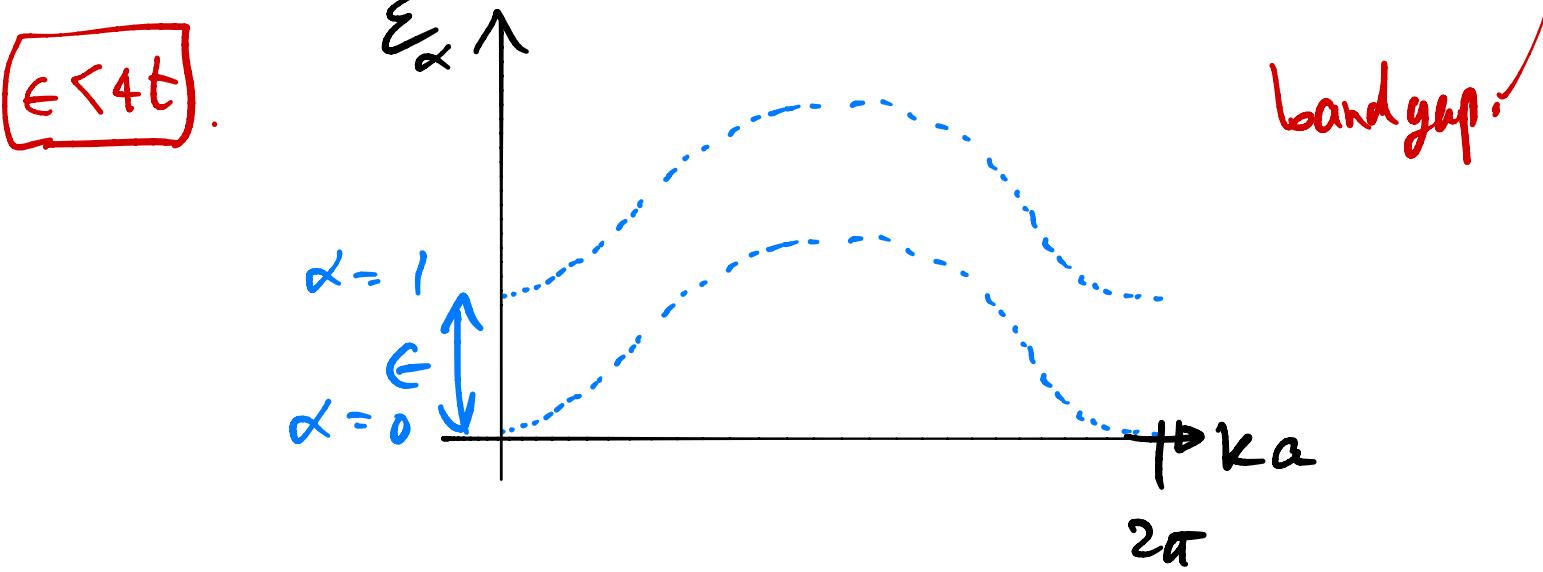
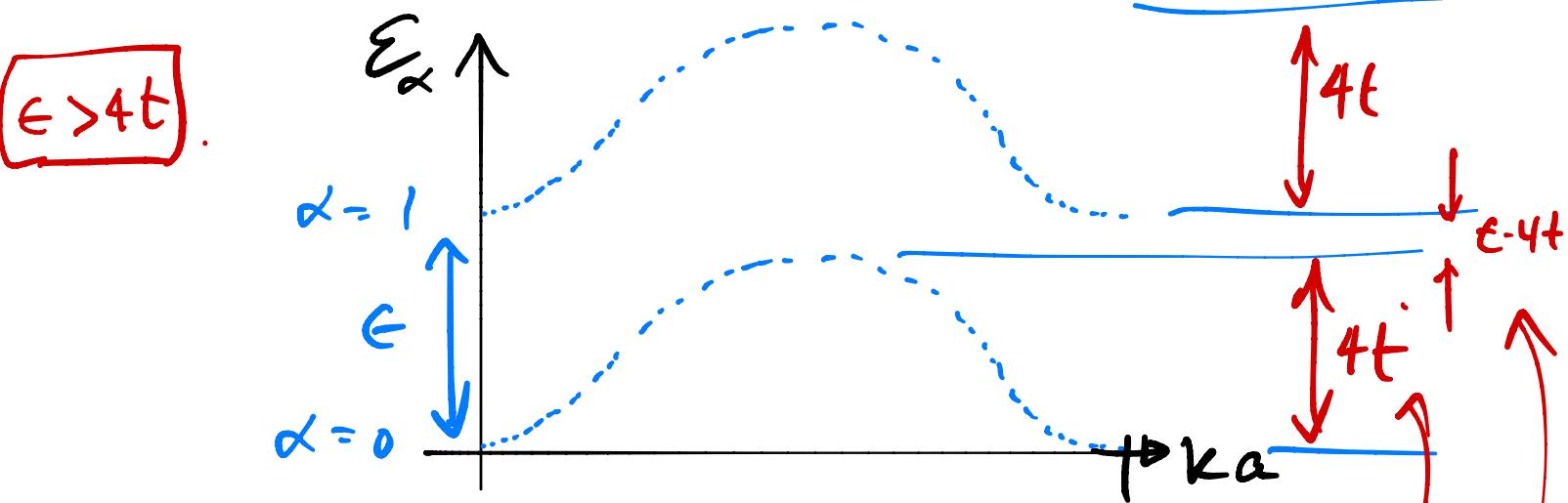
$$\begin{aligned} H_1 = & -t \sum_n (|n+1\rangle \langle n| + |n\rangle \langle n+1|) \otimes \mathbb{1} \\ & + \underbrace{\sum_n |n\rangle \langle n| \otimes (\epsilon |1\rangle \langle 1| + \delta |0\rangle \langle 0|)}_{=} \\ & = H_t + H_e. \end{aligned}$$

$$[H_t, H_\leftarrow] = 0.$$

eigenstates: $|k\rangle \otimes |\alpha\rangle = \frac{1}{\sqrt{n}} \sum_n e^{i k a n} |n\rangle \otimes |\alpha\rangle$

levels: $\epsilon_\alpha(k) = (2 - 2 \cos ka)t + \epsilon_{\delta\alpha, 1}$

$N \gg \infty$: $k_j = \frac{2\pi j}{Na}, j=1..N$



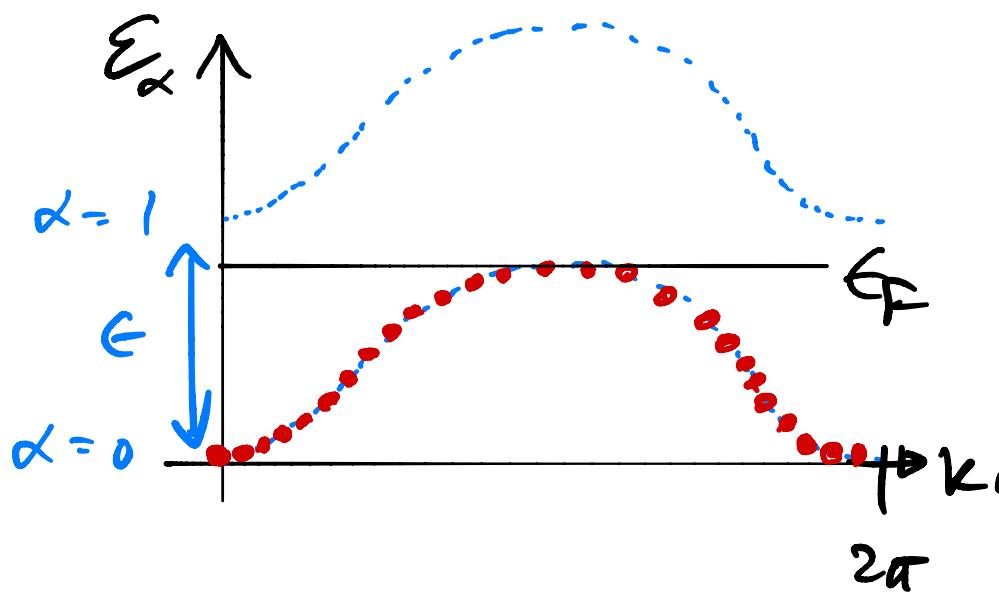
Many fermions :

$$H = \sum_i H_i(i)$$

eigenvalues are

$$\sum_{i=1}^n E_\alpha(k_i)$$

$$\{k_1\alpha_1, \dots, k_n\alpha_n\}$$



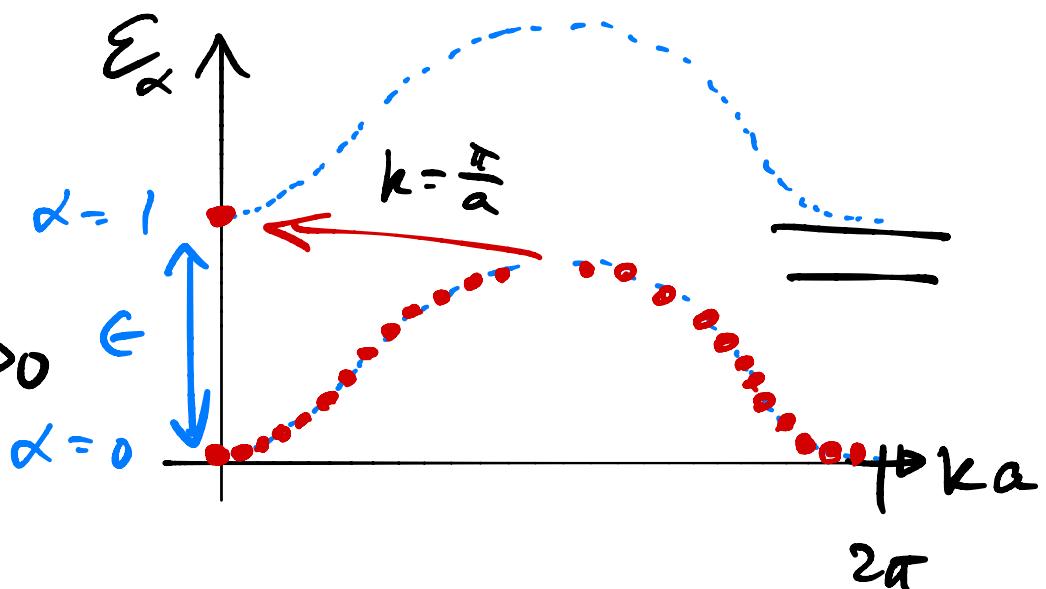
Half-filling: # of electrons = # of sites = N .

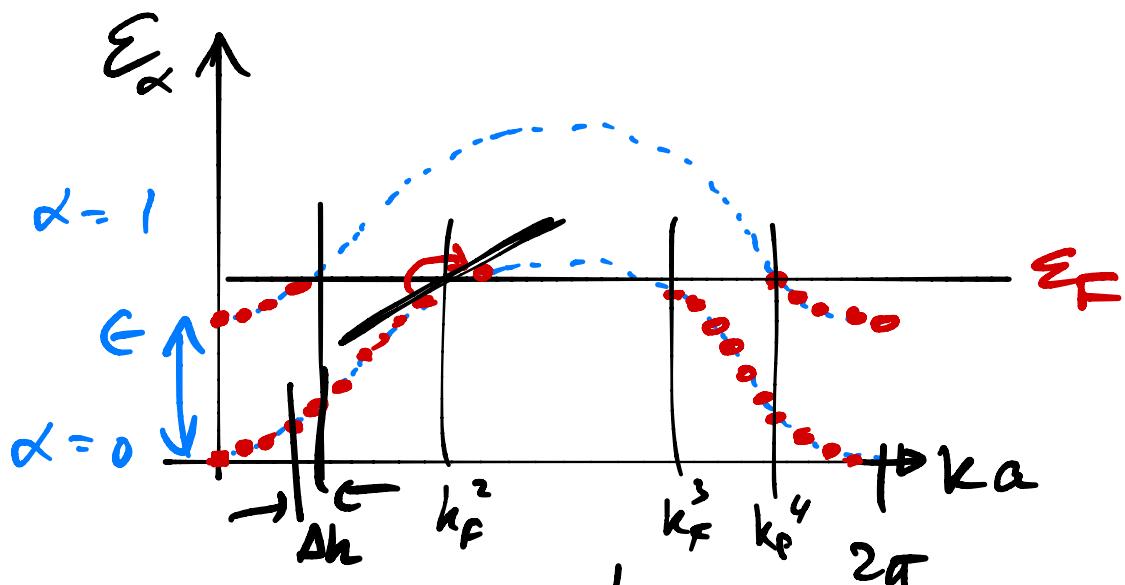
First excited state:

$$\Delta E \xrightarrow{N \rightarrow \infty} \underline{\underline{\epsilon - 4t}} > 0$$

ind. of N

ie energy gap.



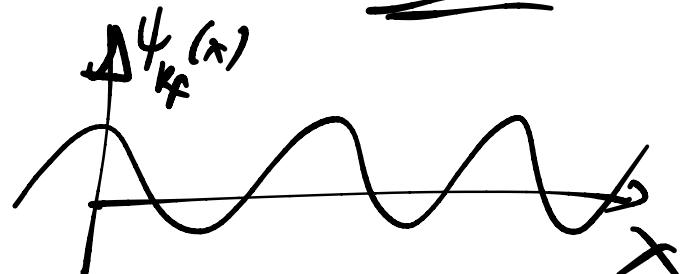


$$\Delta E = \frac{1}{N} \sum_{k \neq k_F} \frac{\partial_k E(k)}{E_F} \propto \frac{1}{N} \xrightarrow{N \rightarrow \infty} 0$$

$= \frac{2\pi}{Na} v_F$

gapless

Conduction requires that
the states near E_F → metal.
are extended.



$$j_x = \text{Im} \Psi_{k_F}^* \partial_x \Psi_{k_F} \neq 0$$

so far $\sigma_{DC} = \infty$.

To fix this:
 — phonons
 — impurities in the lattice.

$$\frac{d\vec{p}}{dt} = q\vec{E}$$

$$\vec{j} = \frac{e\vec{p}}{m} \stackrel{?}{=} \sigma \vec{E}$$

$$\overbrace{i\omega \vec{p}} = q\vec{E}$$

$$\Rightarrow \vec{j} = \frac{e}{m} \frac{q}{i\omega} \cdot \vec{E}$$

$$P = \frac{qE}{i\omega}$$

$$\sigma_{DC} = \lim_{\omega \rightarrow 0} \left(\frac{eq}{mw} \right) = \infty$$

Creation & annihilation

Fermionic operators

$$|0\rangle \equiv |n^0_{\text{electrons}}\rangle$$

$$c_{n\alpha}|0\rangle \stackrel{!}{=} 0. \quad c_{n\alpha}^+|0\rangle \stackrel{!}{=} |n\rangle \otimes |\alpha\rangle$$

Recall: $f_{AB} = [a_A, a_B^\dagger] = \cancel{a_A a_B^\dagger} - \cancel{a_B a_A^\dagger}$

$$[a_A^\dagger, a_B^\dagger] = 0 \quad a_A^\dagger a_B^\dagger |0\rangle = + a_B^\dagger a_A^\dagger |0\rangle.$$

Pauli: $(c_{n\alpha}^+)^2 = 0. \quad c_{n\alpha}^2 = 0. \quad \forall n, \alpha$

$$f_{AB} = c_{A\beta} c_{B\beta}^\dagger + c_{B\beta} c_{A\beta}^\dagger$$

$$\underbrace{\{c_A^+, c_B^+\}}_{\text{anti-commutator}} = 0 \implies c_A^+ c_B^+ |0\rangle = -c_B^+ c_A^+ |0\rangle$$

$$= c_A^+ c_B^+ + c_B^+ c_A^+.$$

$$\left\{ \begin{array}{l} \{c_{n\alpha}, c_{n'\alpha'}^+\} : f_{nn'} \cdot f_{\alpha\alpha'} \\ \{c_{n\alpha}, c_{n'\alpha'}^-\} = 0. \end{array} \right.$$

One Mode . $\mathcal{H}_1 = 1 \cdot >$

$$c^2 = 0, \{c, c^+\} = 1.$$

$$0 = c|0\rangle = c|\downarrow\rangle. \quad c^+|\downarrow\rangle = |\uparrow\rangle$$

$$\begin{aligned} \mathcal{H}_{\text{many}} &= \text{span} \{ |0\rangle, |1\rangle \} \quad (c^+)|\uparrow\rangle = (c^+)^2|\downarrow\rangle \\ &= \text{span} \{ |\downarrow\rangle, |\uparrow\rangle \} = \text{a single qubit}. \end{aligned}$$

$$[c^+ c, c] = -c, \quad [c^+ c, c^+] = +c^+.$$

$$N = c^+c$$

$$c^+c | \downarrow \rangle = 0 | \downarrow \rangle$$

$$c^+c | \uparrow \rangle = 1 | \uparrow \rangle$$

eigenstates of $N = c^+c$

ex:

Basis of hermitian ops on a qbit : $\mathbb{1}, \sigma^x, \sigma^y, \sigma^z$.

$$\sigma^x = c + c^\dagger, \sigma^y = \frac{c - c^\dagger}{i}, \sigma^z = 2c^\dagger c - 1 = 2N - 1$$

$$\sigma^i \sigma^j = \underbrace{i \epsilon^{ijk} \sigma^k}.$$

$$N = \sum_{i=1}^3 -\mathbb{1}.$$

$$\sigma^+ = c^+, \sigma^- = c.$$

$$\sigma^+ = \begin{pmatrix} 0 & 1 \\ 0 & 0 \end{pmatrix} \quad \sigma^- = \begin{pmatrix} 0 & 0 \\ 1 & 0 \end{pmatrix}.$$

More generally $D = \dim \mathcal{H}_1$.

$$\dim \mathcal{H}_{\text{many}} = 2^D = \sum_{N=0}^D \dim \mathcal{H}_F^{(N)}$$
$$= \sum_{N=0}^D \binom{N}{D}$$
$$= (1+1)^D.$$

Back to tight-binding model:

$$H_e = \epsilon (\# \text{ of electrons w/ } \alpha = 1)$$
$$= \epsilon \cdot \sum_n c_{\alpha n}^+ c_{\alpha n} \delta_{\alpha=1}$$

$$\underline{(n+1 \times n)} = \underline{c_{n+1}^+} \underbrace{1_0 \times 0}_0 \underbrace{|}_{\sim} c_n$$

$$H_t = -t \sum_n \left(\underbrace{c_{n,\alpha}^+}_{\sim} c_{n+1,\alpha} + \underbrace{c_{n+1,\alpha}^+}_{\sim} c_{n,\alpha} \right)$$

$$H_{\text{many}} = -t \sum_{n,\alpha} (c_{n\alpha}^\dagger c_{n+1\alpha} + h.c.)$$

$$+ \sum_{n,\alpha\beta} \underline{\epsilon_{\alpha\beta}} c_{n\alpha}^\dagger c_{n\beta}$$

interaction

$$= \sum_{AB=1}^D c_A^\dagger h_{AB} c_B + \sum_{ABCD} \underline{\underline{\underline{\underline{\epsilon_{ABCD}}}}} c_A^\dagger c_B c_C^\dagger c_D$$

$$H_{\text{many}} = H_{\text{many}}^+ \Leftrightarrow h_{AB} = (h^+)_{AB} \underline{\underline{\underline{\underline{\quad}}}}$$

Diagonalize: $h_{AB} = U_{AK} \epsilon_K U_{KB}^+$

$$\text{i.e. } h_{AB} \underline{\underline{U_{BK}}} = \epsilon_K \underline{U_{AK}}.$$

$$\Rightarrow H_{\text{many}} = \sum_K \epsilon_K c_K^\dagger c_K$$

$$c_K = \sum_A (U^+)_K A c_A.$$

$$\text{ef: } \underline{\underline{C_{k\alpha}}} = \sum_n e^{\frac{ikn\alpha}{\sqrt{N}}} \underline{\underline{c_{n\alpha}}}$$

$$E_{\alpha\beta} = \begin{pmatrix} 0 & 0 \\ 0 & \epsilon \end{pmatrix} \Rightarrow H_{\text{many}} = \sum_{k\alpha} \underline{\underline{c_{k\alpha}^+ c_{k\alpha} \frac{\epsilon_{\alpha}(k)}{\epsilon_{\alpha}}}}$$

$$E(k_1, \dots, k_N) = \sum_{i=1}^N E_{\alpha_i}(k_i)$$

Ground state:

$$|g_s\rangle = \prod_{k\alpha} c_{k\alpha}^+ |0\rangle$$

$\{k\alpha \text{ with the smallest } \epsilon_{\alpha}(k)\}$

$$\Psi(n_1\alpha_1 \dots n_N\alpha_N) = \langle n_1\alpha_1 \dots n_N\alpha_N | g_s \rangle$$

$$= \langle 0 | c_{n_1\alpha_1} \dots c_{n_N\alpha_N} | g_s \rangle$$

$$\text{...} = \det \begin{pmatrix} u_{k_1\beta_1}(n_1\alpha_1) & u_{k_1\beta_1}(n_2\alpha_2) & \dots \\ u_{k_2\beta_2}(n_1\alpha_1) & u_{k_2\beta_2}(n_2\alpha_2) & \dots \\ \vdots & \vdots & \ddots \end{pmatrix}$$

$$= \det_{l,j} \left(e^{\frac{-ik_j l a}{\sqrt{N}}} \right).$$

ikna

$$u_{kp}(n\alpha) = \frac{e^{-ikna}}{\sqrt{N}} f_{\alpha p}$$

$$M_{lj} = \frac{e^{-ikja}}{\sqrt{N}}.$$

1.8 2d Quantiz'n, from scratch

creation ops for general 1-particle states

$$\psi(k) = \langle k | \psi \rangle \quad n_i = \sum_k |k X_k|$$

$$|\psi\rangle = \sum_k |k\rangle \langle k | \psi \rangle = \sum_k \psi(k) |k\rangle.$$

$$\text{let } a^\dagger(\psi) \equiv \sum_k a_k^\dagger \psi(k).$$

$$|k\rangle = a_k^\dagger |0\rangle \Rightarrow a^\dagger(\psi) |0\rangle = \sum_k \psi(k) \underbrace{a_k^\dagger |0\rangle}_{= |\psi\rangle} = |\psi\rangle.$$

$$\text{If } [a_k, a_{k'}^+]_s \equiv a_k a_{k'}^+ - s a_{k'}^+ a_k = \underline{\underline{f_{kk'}}}$$

$s = \pm 1$ for $\begin{array}{l} \text{Bosons} \\ \text{Fermions} \end{array}$

$$= \langle k' | k \rangle.$$

then $a(\varphi_1) a^+(\varphi_2) - s a^+(\varphi_2) a(\varphi_1)$

$$= \langle \varphi_2 | \varphi_1 \rangle.$$

and $[a(\varphi_1), a(\varphi_2)]_s = 0.$

One-body operators.

$$\mathcal{H}_1 = \text{Span} \left\{ |A\rangle \right\}$$

on basis.

$$O_{1A} = \sum_{AB} |A X_B| O_{AB}$$

$$1_A = \sum_A |A\rangle \langle A| \quad 1_B = \sum_B |B\rangle \langle B|$$

$$O_{AB} = \langle A | O_{1B} \rangle$$

$$|u_1 \dots u_N\rangle = \sum_{\pi} s^\pi |u_{\pi_1}\rangle \otimes |u_{\pi_2}\rangle \otimes \dots \otimes |u_{\pi_N}\rangle$$

$$\langle O | u_1 \dots u_n \rangle = \sum_{\pi} s^{\pi} \left[\underbrace{\langle O | u_{\pi_1} \rangle | u_{\pi_2} \rangle \dots | u_{\pi_N} \rangle}_{+ | u_{\pi_1} \rangle \langle O | u_{\pi_2} \rangle \dots + \dots + | u_{\pi_1} \rangle \otimes | u_{\pi_2} \rangle \dots \dots \langle O | u_{\pi_N} \rangle} \right]$$

$\langle O | u \rangle$

$$= \sum_{AB} \langle O_{AB} | A \times B | u \rangle$$

Replace B with A.

$$\overline{a^*(A)a(B)}$$

$$\Rightarrow \boxed{O = \sum_{AB} \overline{a^*(A)a(B)} O_{AB}}$$

eg: $H_i = \frac{P^2}{2m}$. $\mathcal{H}_i = \text{span} \{ | p \rangle \}$
 $P \in \mathbb{R}$

$$H_n = \sum_{i=1}^n \frac{P_i^2}{2m} = \Pi_n H \Pi_n$$

$$H = \sum_p \underbrace{a_p^* a_p}_{\sim} \frac{P^2}{2m}.$$

Π_n = projector onto \mathcal{H}_n .

and 316

$$|AB\rangle = \underline{c_A^+ c_B^+ |0\rangle} \quad \}$$

$$\begin{aligned} & \text{one particle, one particle} \\ & \text{in } A, \text{ in } B \end{aligned} \quad \underline{-c_B^+ c_A^+ |0\rangle} \\ & = -\underline{|BA\rangle}$$

$$\Leftarrow c_A^+ c_B^+ + c_B^+ c_A^+ = 0.$$

$$\underline{= \{c_A^+, c_B^+\}}$$

$$\mathcal{H}_1 = \text{span} \{ |A\rangle \}$$

$$(c^+)^2 = 0$$

$$N = c^+ c_-$$

$$\Rightarrow \left\{ \begin{array}{l} [c^+ c_-, c] = -c \\ [c^+ c_-, c^+] = +c^+ \end{array} \right. \quad \left\{ \begin{array}{l} c|0\rangle = 0 \\ c^+|0\rangle = |1\rangle. \end{array} \right.$$

$$A=1..D \rightarrow \mathcal{J}^D = \dim \mathcal{H}_1$$

$$\boxed{\{c, c^+ \} = 1} \implies cc^+ = 1 - \cancel{c^+c}.$$

$$\cancel{c^+c} + cc^+$$

$$\begin{matrix} |0\rangle \rightarrow |0\rangle & |\lambda|^2 \\ |1\rangle \rightarrow |1\rangle & |\lambda|^2 \\ |1\rangle \rightarrow 0 & \end{matrix}$$

Demand: $\begin{matrix} |0\rangle \rightarrow 0 \\ |1\rangle \rightarrow |1\rangle & |\lambda|^2 \end{matrix}$

$$= 1 \cdot |1| + 1 \cdot 0 = 1$$

$$(c)^+ = c^+.$$

$$\left\{ \begin{array}{l} c \rightarrow \lambda c \\ c^+ \rightarrow \lambda^* c \end{array} \right. \quad \left\{ \begin{array}{l} \{c, c^+\} = |\lambda|^2 1 \\ N = \frac{cc^+}{|\lambda|^2} \end{array} \right.$$

$$|0\rangle.$$

$$\frac{c|0\rangle = 0}{c^+|0\rangle = |1\rangle}.$$

$$(c^+)^2 = 0.$$

$$\Rightarrow \mathcal{H} = \text{span}\{|0\rangle, |1\rangle\}.$$

$$\langle 1|0\rangle ? \quad \langle 1|0\rangle = \langle 0|c|0\rangle = 0 \quad \checkmark$$

$$\underbrace{\langle 1|1\rangle = 1}_{\text{ }} \quad \langle 1|1\rangle = \langle 0|cc^+|0\rangle$$

$$= \langle 0|(1 - \cancel{c^+c})|0\rangle$$

$$= \cancel{\langle 0|0\rangle}.$$

$$\langle \underline{e^{ikq(0)}} \rangle_0 = \underbrace{\int [Dq]}_{= \int \prod_{i=-M}^M dq_i} e^{-S[q]} e^{-ikq(0)}$$

$$= \int \prod_{i=-M}^M dq_i e^{-\sum_i (q_i - q_{-i})^2 - \sum_i V(q_i)}$$

if

$$V(q) = \cancel{k_{ij}} q_i q_j.$$

$$Z[J] = \int \prod dq_i e^{-q_i M_{ij} q_j + q_i J_i}$$

$$J_i = f_{ij} \cancel{ik}.$$

$$M_{ij} = \begin{pmatrix} 2 & 0 & 0 \\ 0 & 2 & 0 \\ 0 & 0 & 1 \end{pmatrix} + K_{ij}$$

$$Z[J] = \frac{1}{\sqrt{\det M}} e^{-\frac{1}{2} J M^{-1} J}$$

$$\langle e^{ikq} \rangle \propto e^{-\frac{1}{2} k^2 (\tilde{M})_{00}}$$

$$\langle \underline{q_{(0)}^2} \rangle = \left(\frac{\partial}{\partial J_0} \right)^2 \ln Z[J] \Big|_{J=0} = (\tilde{M})_{00}$$

$$\begin{aligned}
 e^{ikq} &= e^{ikN(a + a^\dagger)} \\
 &\stackrel{\text{BCH}}{=} e^{\cancel{ik a^\dagger N} \cancel{i k a N} e^{-\hbar^2 [a, a^\dagger]}} \cancel{e^{\frac{-\hbar^2}{2} N^2}} \\
 \langle 0 | e^{ik a^\dagger} e^{ik a} | 0 \rangle &= e^{-k^2 N^2} \\
 \underbrace{\langle 0 |}_{\langle 0 |} &\quad \underbrace{| 0 \rangle}_{| 0 \rangle} \quad \frac{N^2 = \langle q^2 \rangle}{}.
 \end{aligned}$$