

Tight Binding Model for Twist Bilayer Graphene*

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* A footnote to the article title

I. SYMBOL CONVENTION

II. THE TIGHT BINDING HAMILTONIAN FOR TWIST BILAYER GRAPHENE

The tight binding Hamiltonian written in the second quantization language is

$$\hat{H} = \sum_{\bar{\mathbf{k}}} \sum_{\alpha n, \beta m} H_{\alpha n, \beta m} C_{\alpha}^{\dagger}(\bar{\mathbf{k}} + \mathbf{G}_n) C_{\beta}(\bar{\mathbf{k}} + \mathbf{G}_m). \quad (1)$$

The basis wave function can be expressed as a summation of planewave:

$$\psi_{\alpha, n}(\bar{\mathbf{k}}) = \frac{1}{\sqrt{N_m N_a}} \sum_{\mathbf{I}, i} e^{i(\bar{\mathbf{k}} + \mathbf{G}_n) \mathbf{R}_{\mathbf{I}i\alpha}} |\phi_{p_z}\rangle, \quad (2)$$

$N_a = N_{\text{atom}}/4$, And

$$\mathbf{R}_{\mathbf{I}i\alpha} = \mathbf{L}_{\mathbf{I}} + \boldsymbol{\tau}_{i\alpha}. \quad (3)$$

The matrix element of the Hamiltonian should be written as:

$$\begin{aligned} H_{\alpha n, \beta m} &= \frac{1}{N_m N_a} \sum_{\mathbf{I}, \mathbf{J}, ij} t(\mathbf{R}_{\mathbf{I}i\alpha} - \mathbf{R}_{\mathbf{J}j\beta}) e^{-i(\bar{\mathbf{k}} + \mathbf{G}_n) \mathbf{R}_{\mathbf{I}i\alpha}} e^{i(\bar{\mathbf{k}} + \mathbf{G}_m) \mathbf{R}_{\mathbf{J}j\beta}} \\ &= \frac{1}{N_m N_a} \sum_{\mathbf{I}, \mathbf{J}, ij} t(\mathbf{R}_{\mathbf{I}i\alpha} - \mathbf{R}_{\mathbf{J}j\beta}) e^{-i(\bar{\mathbf{k}} + \mathbf{G}_n) \cdot (\mathbf{L}_{\mathbf{I}} + \boldsymbol{\tau}_{i\alpha})} e^{i(\bar{\mathbf{k}} + \mathbf{G}_m) \cdot (\mathbf{L}_{\mathbf{J}} + \boldsymbol{\tau}_{j\beta})} \\ &= \frac{1}{N_m N_a} \sum_{\mathbf{I}, \mathbf{J}, ij} e^{-i\mathbf{G}_n \boldsymbol{\tau}_{i\alpha}} e^{-i\bar{\mathbf{k}}(\mathbf{L}_{\mathbf{I}} - \mathbf{L}_{\mathbf{J}} + \boldsymbol{\tau}_{i\alpha} - \boldsymbol{\tau}_{j\beta})} t(\mathbf{L}_{\mathbf{I}} - \mathbf{L}_{\mathbf{J}} + \boldsymbol{\tau}_{i\alpha} - \boldsymbol{\tau}_{j\beta}) e^{i\mathbf{G}_m \boldsymbol{\tau}_{j\beta}} \\ \text{cutoff: } \langle \mathbf{I}, \mathbf{J} \rangle, &= \frac{1}{N_m N_a} \sum_{\mathbf{I}, ij} e^{-i\mathbf{G}_n \boldsymbol{\tau}_{i\alpha}} e^{-i\bar{\mathbf{k}}(\bar{\boldsymbol{\tau}}_{i\alpha, j\beta})} t(\bar{\boldsymbol{\tau}}_{i\alpha, j\beta}) e^{i\mathbf{G}_m \boldsymbol{\tau}_{j\beta}} \\ \frac{1}{N_m} \sum_{\mathbf{I}} &= 1, \quad = \frac{1}{N_a} \sum_{ij} e^{-i\mathbf{G}_n \boldsymbol{\tau}_{i\alpha}} e^{-i\bar{\mathbf{k}}(\bar{\boldsymbol{\tau}}_{i\alpha, j\beta})} t(\bar{\boldsymbol{\tau}}_{i\alpha, j\beta}) e^{i\mathbf{G}_m \boldsymbol{\tau}_{j\beta}}. \end{aligned} \quad (4)$$

Tricky here, we write down the matrix form of the Hamiltonian,

$$\begin{aligned} \mathbf{H}_{\alpha\beta} &= \mathbf{X}_{\alpha}^{\dagger} \mathbf{T}_{\alpha\beta} \mathbf{X}_{\beta} \\ \mathbf{H} &= \oplus_{\alpha\beta} \mathbf{H}_{\alpha\beta} \end{aligned} \quad (5)$$

The above equation describes a block matrix multiplication. α, β runs in $[A_1, B_1, A_2, B_2]$

The matrix element:

$$\begin{aligned} (\mathbf{X}_{\alpha}^{\dagger})_{n,i} &= e^{i\mathbf{G}_n \boldsymbol{\tau}_{i\alpha}}, \\ (\mathbf{T}_{\alpha\beta})_{i,j} &= e^{-i\bar{\mathbf{k}}(\bar{\boldsymbol{\tau}}_{i\alpha, j\beta})} t(\bar{\boldsymbol{\tau}}_{i\alpha, j\beta}). \end{aligned} \quad (6)$$

III. TBG UNDER MAGNETIC FIELD

A. Magnetic Translation Operator

The original lattice translation symmetry is broken by the vector potential:

$$\hat{T}_{\mathbf{L}_I} \hat{T}_{\mathbf{L}_J} = \hat{T}_{\mathbf{L}_J} \hat{T}_{\mathbf{L}_I} \exp\left(i \frac{e}{\hbar} \oint \mathbf{A} d\mathbf{r}\right) \quad (7)$$

which gives us the form of AB effect in lattice:

$$\hat{T}_{\mathbf{L}_I}^{-1} \hat{T}_{\mathbf{L}_J}^{-1} \hat{T}_{\mathbf{L}_I} \hat{T}_{\mathbf{L}_J} = \exp \left(i \frac{e}{\hbar} \oint_S \mathbf{A} d\mathbf{r} \right) \quad (8)$$

S is the cell size spanned by $\mathbf{L}_I, \mathbf{L}_J$. Cosider a specific case:

$$\begin{aligned} \mathbf{L}_I &= \mathbf{L}_1 \\ \mathbf{L}_J &= q\mathbf{L}_2 \\ \Phi &= BS = \frac{p}{q} \Phi_0 \end{aligned} \quad (9)$$

where p, q are relatively prime, S is the moire unit lattice size. $\Phi_0 = h/e$. When B is uniform,

$$\begin{aligned} S &= qS \\ \exp \left(i \frac{e}{\hbar} \oint_S \mathbf{A} d\mathbf{r} \right) &= \exp \left(i \frac{e}{\hbar} \int_{\partial S} B dS \right) \\ &= \exp(2\pi i p) \\ &= 1 \end{aligned} \quad (10)$$

Subsequently, we have the commutation relationship:

$$\hat{T}_{\mathbf{L}_1} \hat{T}_{q\mathbf{L}_2} = \hat{T}_{q\mathbf{L}_2} \hat{T}_{\mathbf{L}_1} \quad (11)$$

It suggests that we should adopt a magnetic lattice unit cell q times larger than the original one.

B. Tight Binding Hopping term

By using Peierls substitution, the hopping term should be written as

$$\begin{aligned} \tilde{t}(\mathbf{R}_{Ii\alpha} - \mathbf{R}_{Jj\beta}) &= t(\mathbf{R}_{Ii\alpha} - \mathbf{R}_{Jj\beta}) \exp \left(-i \frac{e}{\hbar} \int_{\mathbf{R}_{Ii\alpha}}^{\mathbf{R}_{Jj\beta}} \mathbf{A} d\mathbf{r} \right) \\ &= t(\mathbf{R}_{Ii\alpha} - \mathbf{R}_{Jj\beta}) \exp \left(-2\pi i \frac{1}{\Phi_0} \int_{\mathbf{R}_{Ii\alpha}}^{\mathbf{R}_{Jj\beta}} \mathbf{A} d\mathbf{r} \right) \end{aligned} \quad (12)$$

Here we use the Landau gauge, $\mathbf{A} = (0, Bx, 0)$, the corresponding phase factor can be calculated in the following way:

$$\begin{aligned} \int_{\mathbf{R}_1}^{\mathbf{R}_2} \mathbf{A} d\mathbf{r} &= \int_{\mathbf{R}_1}^{\mathbf{R}_2} Bx dy \\ &= B \frac{y_2 - y_1}{x_2 - x_1} \int_{x_1}^{x_2} x dx \\ &= \frac{B(x_2 + x_1)(y_2 - y_1)}{2} \end{aligned} \quad (13)$$