Phonons and Electron-Phonon Couplings

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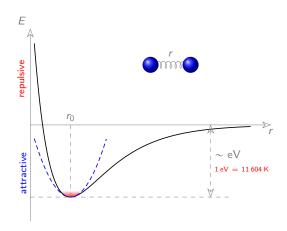
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Outline

- Lattice Dynamics
 - 1D Atomic Chain
 - 3D lattice
 - How to calculate the dynamical matrix

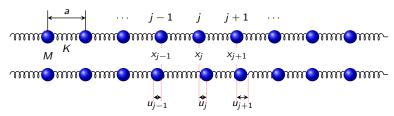
Density functional perturbation theory

Interatmoic Potential



1D Chain of Atoms — 1 Atom per Unit

A 1D chain of N equally spaced atoms at $R_j(t) = x_j + u_j(t)$



The Newton's Equation

$$M\frac{\mathrm{d}^2 u_j}{\mathrm{d}t^2} = K(u_{j+1} + u_{j-1} - 2u_j)$$
 $j = 1, ..., N$

Assume the solution has the form $u_j(t)=rac{A_q}{\sqrt{M}}e^{i(q\mathbf{x_j}-\omega t)}$, then 1

$$\omega^{2} = \frac{K}{M} (2 - e^{iqa} - e^{-iqa})$$
$$= \frac{2K}{M} (1 - \cos qa)$$
$$\Rightarrow \quad \omega = \sqrt{\frac{4K}{M}} \left| \sin \frac{qa}{2} \right|$$



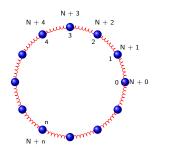
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 $^{^{1}}u_{j}(t)$ here is complex. In practice, take the real part, i.e. $\mathrm{Re}[u_{j}(t)]$.

Born-von Karman boundary condition

The Born-von Karman Periodic Boundary Condition:

$$u_n = u_{N+n}$$
 \Rightarrow $e^{iqx_j} = e^{iqx_{N+j}}$ \Rightarrow $e^{iqNa} = 1$
 \Rightarrow $q = \frac{2\pi}{a} \frac{I}{N};$ $(I = 0, 1, ..., N)$

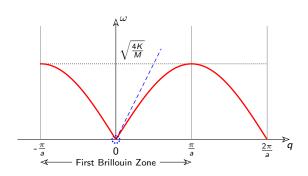


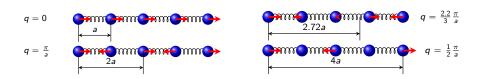
Usually, q is restricted within the first Brillouin Zone

$$-\frac{\pi}{a} \leqslant q \leqslant \frac{\pi}{a}$$



1D Chain of Atoms

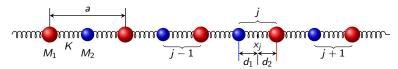






1D Chain of Atoms — 2 Atoms per Unit

A 1D chain with 2 atoms in each unit: $R_s^j(t) = x_j + d_s + u_s^j(t)$; s = 1, 2



The Newton's Equation

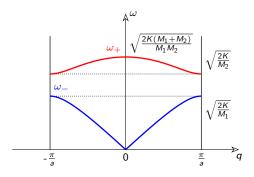
$$\begin{aligned} & M_1 \frac{\mathrm{d}^2 u_1^j}{\mathrm{d} t^2} = K(u_2^j + u_2^{j-1} - 2u_1^j) \\ & M_2 \frac{\mathrm{d}^2 u_2^j}{\mathrm{d} t^2} = K(u_1^j + u_1^{j+1} - 2u_2^j) \end{aligned} \implies \begin{cases} u_j^1(t) = \frac{A_q}{\sqrt{M_1}} \mathrm{e}^{i(q x_j - \omega t)} \\ u_j^2(t) = \frac{B_q}{\sqrt{M_2}} \mathrm{e}^{i(q x_j - \omega t)} \end{cases}$$

We then have

$$\begin{pmatrix} \frac{2K}{M_1} & \frac{-K}{\sqrt{M_1M_2}}(1+e^{-iqa}) \\ \frac{-K}{\sqrt{M_1M_2}}(1+e^{iqa}) & \frac{2K}{M_2} \end{pmatrix} \begin{pmatrix} A_q \\ B_q \end{pmatrix} = \omega^2 \begin{pmatrix} A_q \\ B_q \end{pmatrix}$$

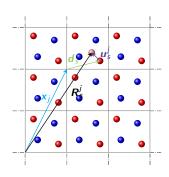
$$\implies \omega_{\pm}^2 = \frac{K}{M_1M_2} \left((M_1+M_2) \pm \sqrt{M_1^2 + M_2^2 + 2M_1M_2\cos qa} \right)$$

1D Chain of Atoms



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3D Lattice



- x_j : the position of unit cell j
- d_s: the equilibrum position of the atom s in the cell
- \mathbf{u}_s^j : displacement from the equilibrum positon for the atom s in the cell j
- R_s^j : the position of the atom s in the cell j

$$R_s^j(t) = x_j + d_s + u_s^j(t)$$

$$= r_s^j + u_s^j(t)$$

$$R_s^j(t) = r_s^j + u_s^j(t) \quad (c_s - x_s)$$

 $R_{s\alpha}^{j}(t) = r_{s\alpha}^{j} + u_{s\alpha}^{j}(t) \quad (\alpha = x, y, z)$

The total energy can be written as

$$E_{\text{tot}}\left(\{\boldsymbol{R}_{s}^{j}(t)\}\right) = E_{\text{tot}}^{0}\left(\{\boldsymbol{r}_{s}^{j}\}\right) + \sum_{js\alpha} \frac{\partial E_{\text{tot}}^{0}}{\partial u_{s\alpha}^{j}} u_{s\alpha}^{j} + \frac{1}{2} \sum_{\substack{js\alpha \\ k \neq \beta}} \frac{\partial^{2} E_{\text{tot}}^{0}}{\partial u_{s\alpha}^{j} \partial u_{t\beta}^{k}} u_{s\alpha}^{j} u_{t\beta}^{k} + \dots$$

- The expression is exact if we take all the orders in the expansion.
- All the derivatives are taken at the equilibrium positions $\{r_s^j\}$, i.e. $\frac{\partial E_{\rm tot}^0}{\partial u_{\rm for}^j}=0$.
- Harmonic approximation: truncated at *second* order.



3D Lattice Dynamics

Within the harmonic approximation, the Newton's equation for the atom s in cell j

$$M_{s} \frac{\mathrm{d}^{2} u_{s\alpha}^{j}(t)}{\mathrm{d}t^{2}} = -\frac{\partial E_{\text{tot}}}{\partial u_{s\alpha}^{j}} = -\sum_{kt\beta} \frac{\partial^{2} E_{\text{tot}}^{0}}{\partial u_{s\alpha}^{j} \partial u_{t\beta}^{k}} u_{t\beta}^{k} = -\sum_{kt\beta} C_{s\alpha,t\beta}^{j,k} u_{t\beta}^{k}$$
(1)

The ansatz of the solution

$$u_{s\alpha}^{j}(t) = \frac{\eta_{s\alpha}^{\sigma}(\mathbf{q})}{\sqrt{M_{s}}} e^{i\mathbf{q}\mathbf{x}_{j}} e^{-i\omega_{\sigma}t}$$
(2)

Substitute Eq. 2 into Eq. 1

$$\omega_{\sigma}^{2}(\boldsymbol{q})\,\eta_{s\alpha}^{\sigma} = \sum_{t\beta} \left[\sum_{k} \frac{1}{\sqrt{M_{s}M_{t}}} \, \frac{\partial^{2} E_{\text{tot}}^{0}}{\partial u_{s\alpha}^{i} \partial u_{t\beta}^{k}} \, e^{i\boldsymbol{q}(\boldsymbol{x}_{k} - \boldsymbol{x}_{j})} \right] \eta_{t\beta}^{\sigma} = \sum_{t\beta} D_{s\alpha,t\beta}(\boldsymbol{q}) \, \eta_{t\beta}^{\sigma}$$

In matrix form

$$\begin{pmatrix} \ddots & & & \\ & D_{s\alpha,t\beta}(\mathbf{q}) & & \\ & & \ddots \end{pmatrix} \begin{pmatrix} \vdots \\ \eta_{t\beta}^{\sigma}(\mathbf{q}) \\ \vdots \end{pmatrix} = \omega^{2}(\mathbf{q}) \begin{pmatrix} \vdots \\ \eta_{t\beta}^{\sigma}(\mathbf{q}) \\ \vdots \end{pmatrix}$$

$$3N_{a} \times 3N_{a} \qquad 3N_{a} \qquad \text{polarization vector}$$

where $\sigma = 1, ..., 3N_a$ and N_a is the number of atoms in the *primitive cell*.

The Interatomic Force Constants

The Interatomic Force Constants (IFC)

$$\frac{\partial^2 E_{\text{tot}}^0}{\partial u_{s\alpha}^j \partial u_{t\beta}^k} = C_{s\alpha,t\beta}^{j,k}$$

• Symmetric because partial differentiation is commutative

$$\frac{\partial^2 E_{\text{tot}}^0}{\partial u_{s\alpha}^j \partial u_{t\beta}^k} = \frac{\partial^2 E_{\text{tot}}^0}{\partial u_{t\beta}^k \partial u_{s\alpha}^j} \quad \Rightarrow \quad C_{s\alpha,t\beta}^{j,k} = C_{t\beta,s\alpha}^{k,j} \tag{3}$$

ullet Translation invariance, only depend on the difference between j and k

$$\frac{\partial^2 E_{\text{tot}}^0}{\partial u_{s\alpha}^j \partial u_{t\beta}^k} = \frac{\partial^2 E_{\text{tot}}^0}{\partial u_{s\alpha}^0 \partial u_{t\beta}^{(k-j)}} \quad \Rightarrow \quad C_{s\alpha,t\beta}^{j,k} = C_{s\alpha,t\beta}^{0,k-j} \tag{4}$$

 Acoustic Sum Rule (ASR): if we displace the whole solid by an arbitrary uniform displacement, the forces acting on the atoms must be zero.

$$F_{s\alpha}^{j} = -\sum_{\beta} \left[\sum_{kt} \frac{\partial^{2} E_{\text{tot}}^{0}}{\partial u_{s\alpha}^{j} \partial u_{r\beta}^{k}} \right] \delta_{\beta} = 0 \qquad \Rightarrow \qquad \sum_{kt} \frac{\partial^{2} E_{\text{tot}}^{0}}{\partial u_{s\alpha}^{j} \partial u_{r\beta}^{k}} = 0$$
 (5)

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The Dynamical Matrix

The Dynamical Matrix

$$D_{s\alpha,t\beta}(\boldsymbol{q}) = \frac{1}{\sqrt{M_s M_t}} \sum_{l} \frac{\partial^2 E_{\text{tot}}^0}{\partial u_{s\alpha}^0 \partial u_{t\beta}^l} e^{i\boldsymbol{q} \boldsymbol{x}_l} = \frac{1}{\sqrt{M_s M_t}} \sum_{l} C_{s\alpha,t\beta}^{0,l} e^{i\boldsymbol{q} \boldsymbol{x}_l}$$

• If we define the distortion pattern $\boldsymbol{u}_s^I(\boldsymbol{q}) = \boldsymbol{v}_s(\boldsymbol{q}) e^{i \boldsymbol{q} \boldsymbol{x}_I}$

$$D_{s\alpha,t\beta}(\boldsymbol{q}) = \frac{1}{N} \frac{1}{\sqrt{M_s M_t}} \frac{\partial^2 E_{\text{tot}}^0}{\partial v_{s\alpha}^*(\boldsymbol{q}) \partial v_{t\beta}(\boldsymbol{q})}$$

ullet Dynamical matrix is Hermitian and admit real eigenvalues $\omega^2({m q})$

$$D_{s\alpha,t\beta}(\boldsymbol{q}) = D_{t\beta,s\alpha}^*(\boldsymbol{q})$$

Proof

$$\begin{split} D_{s\alpha,t\beta}(\boldsymbol{q}) &= \frac{1}{\sqrt{M_s M_t}} \sum_{l} C_{s\alpha,t\beta}^{0,l} e^{i\boldsymbol{q}\boldsymbol{x}_l} \\ &= \frac{1}{\sqrt{M_s M_t}} \sum_{l} C_{s\alpha,t\beta}^{l,0} e^{i\boldsymbol{q}\boldsymbol{x}_l} \\ &= \frac{1}{\sqrt{M_s M_t}} \sum_{l} C_{s\alpha,t\beta}^{l,0} e^{i\boldsymbol{q}\boldsymbol{x}_l} \\ &= \frac{1}{\sqrt{M_s M_t}} \sum_{l} C_{t\beta,s\alpha}^{0,l} e^{i\boldsymbol{q}\boldsymbol{x}_l} \\ &= D_{t\beta,s\alpha}(\boldsymbol{q}) \end{split}$$

Phonon Polarization Vectors

The eigenvectors $\eta^{\sigma}(\mathbf{q})$ of the Hermitian matrix $D_{s\alpha,t\beta}(\mathbf{q})$ are called the phonon polarization vector.

• The polarization vector is *cell-periodic*.

$$u_{s\alpha}^{j}(t) = \frac{\eta_{s\alpha}^{\sigma}(\mathbf{q})}{\sqrt{M_{s}}} e^{i\mathbf{q}\mathbf{x}_{j}} e^{-i\omega_{\sigma}t}$$

So the solution is a cell-periodic part multiply by e^{iqx} — Bloch's theorem.

Orthogonalization relation:

$$\sum_{s\alpha} \eta_{s\alpha}^{\sigma'*}(\boldsymbol{q}) \, \eta_{s\alpha}^{\sigma}(\boldsymbol{q}) = \delta_{\sigma\sigma'}; \qquad \sum_{\sigma} \eta_{s\alpha}^{\sigma*}(\boldsymbol{q}) \, \eta_{t\beta}^{\sigma}(\boldsymbol{q}) = \delta_{st} \, \delta_{\alpha\beta}$$

• Relation to phonon displacement — direction and amplitude of the vibration.

$$\xi_{s\alpha}^{\sigma} = \frac{1}{\sqrt{M_s}} \, \eta_{s\alpha}^{\sigma}(\mathbf{q})$$

• At some high-symmetry q-path

$$egin{cases} m{q} \parallel m{\eta}(m{q}) & ext{Longitudinal Wave} \ m{q} \perp m{\eta}(m{q}) & ext{Transverse Wave} \end{cases}$$

How to Calculate the Dynamical Matrix I

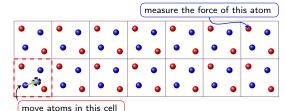
The definition of the dynamical matrix

$$D_{s\alpha,t\beta}(\mathbf{q}) = \frac{1}{\sqrt{M_s M_t}} \sum_{l=-\infty}^{\infty} C_{s\alpha,t\beta}^{0,l} e^{i\mathbf{q}\mathbf{x}_l} \approx \frac{1}{\sqrt{M_s M_t}} \sum_{|l| < l_{\text{cut}}} C_{s\alpha,t\beta}^{0,l} e^{i\mathbf{q}\mathbf{x}_l}$$
(6)

Finite-difference and supercell approach — Frozen phonon method

IFC by finite-difference:

$$\begin{split} &\frac{\partial^2 E_{tot}^0}{\partial u_{s\alpha}^0 \partial u_{t\beta}^l} = \frac{\partial F_{t\beta}^l}{\partial u_{s\alpha}^0} \\ &\approx \frac{F_{t\beta}^l(\Delta_{s\alpha}) - F_{t\beta}^l(-\Delta_{s\alpha})}{2\Delta_{s\alpha}} \end{split}$$

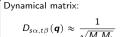


- Supercell must be large enough so that IFC is negligible at the cell boundary.
- Movements done only in one primitive cell.
- $3 \times N_a \times 2$ movements, i.e. move by $\pm \Delta$ in x/y/z directions for each atom in the primitive cell.
- Symmetry can be adopted to reduce the number of movements.
- The dynamical matrix can then be obtained at arbitrary q by Eq. 6.

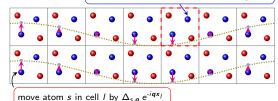
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How to Calculate the Dynamical Matrix II

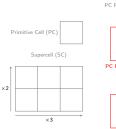
measure the force of atoms in arbitrary cell

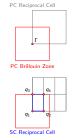


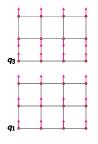
$$D_{s\alpha,t\beta}(\mathbf{q}) \approx \frac{1}{\sqrt{M_s M_t}} \times \frac{F_{t\beta}^{I}(\Delta_{s,\mathbf{q}}) - F_{t\beta}^{I}(-\Delta_{s,\mathbf{q}})}{2\Delta_{s,\mathbf{q}}}$$

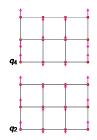


• Can only obtain dynamical matrix at certain q.





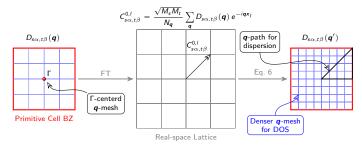




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$$D_{s\alpha,t\beta}(\mathbf{q}) \xrightarrow{\mathsf{FT}} C_{s\alpha,t\beta}^{0,l} \xrightarrow{\mathsf{Eq. 6}} D_{s\alpha,t\beta}(\mathbf{k})$$

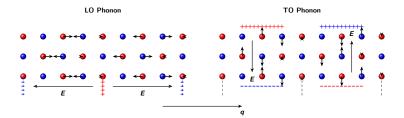
How to Calculate the Dynamical Matrix III

- Lineare response approach density functional perturbation theory
 - Can calculate $D_{s\alpha,t\beta}(\mathbf{q})$ at arbitrary \mathbf{q} using a primitive cell.
 - $D_{s\alpha,t\beta}({m q})$ is periodic in reciprocal space: $D_{s\alpha,t\beta}({m q}+{m G})=D_{s\alpha,t\beta}({m q})$



- In practice, first calculate $D_{s\alpha,t\beta}(q)$ with a small q-mesh. Then, perform FT to get the IFC in real space. Finally, dynamical matrix at arbitrary q can be obtained.
- Fails in metal with Kohn anomalies or in polar semiconductors where the dynamical matrix is non-analytic for $q \to 0$.
- Ocdes: Phonopy, PHON, YPHON, PhonTS, ShengBTE, ALM, ALAMODE, Quantum Espresso, Abinit. Siesta...

LO-TO Splitting



Long-wavelength limit in polar materials

The dynamical matrix in polar materials can be separated into two parts $^{\mathrm{2}}$

$$D_{s\alpha,t\beta}(\boldsymbol{q} \to 0) = D_{s\alpha,t\beta}^{\mathsf{an}}(\boldsymbol{q} = 0) + D_{s\alpha,t\beta}^{\mathsf{na}}(\boldsymbol{q} \to 0)$$

where the nonanalytic part is written as

$$D_{s\alpha,t\beta}^{\mathsf{na}}(\boldsymbol{q}\to 0) = \frac{1}{\sqrt{M_s M_t}} \frac{4\pi e^2}{\Omega} \frac{\left(\sum_{\gamma} q_{\gamma} Z_s^{*\gamma\alpha}\right) \left(\sum_{\mu} q_{\mu} Z_t^{*\mu\beta}\right)}{\sum_{\gamma\mu} q_{\gamma} \epsilon_{\infty}^{\gamma\mu} q_{\mu}} \tag{7}$$

• Z_s^* is the Born effective charge tensor for atom s

$$Z_s^{*\alpha\beta} = \Omega_0 \frac{\partial \mathcal{P}_\beta}{\partial u_{s\alpha}} = \frac{\partial \mathcal{F}_{s\alpha}}{\partial \mathcal{E}_\beta}$$

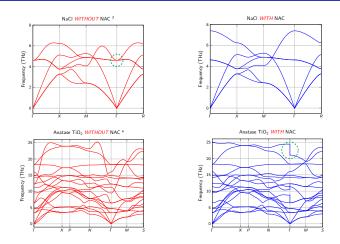
- **1** The response of the polarization per unit cell along the direction β induced by a displacement along the direction α of the atoms belonging to the sublattice s, under the condition of a zero electric field.
- ② The force on the atom s along α induced by the macroscopic field along β .
- **3** Sum rules: $\sum_s Z_s^{*\alpha\beta} = 0$
- $\bullet \ \, \text{ The infrared absorption: } I_{\text{IR}}(\omega_\sigma) \propto \sum_{\alpha=1}^3 \left| \sum_{s\beta} Z_s^{*\alpha\beta} \eta_{s\beta}^\sigma \right|^2$
- ϵ_{∞} is the *electronic* dielectric tensor of the crystal, i.e. the static dielectric constant with clamped nuclei.

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²X. Gonze and C. Lee, *Phys. Rev. B* 55, 10355(1997).

Phonon Band Structure with/without NAC



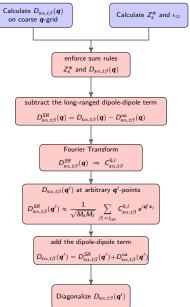
- The limiting value of Eq. 7 may be different for different q → 0 directions, as can be seen in the phonon dispersion of Anatase TiO₂.
- This strange behavior is physical and the frequency of the observed phonons depends on the polarizability
 in the different directions, with respect to the propagation of the incoming electric field.

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 $^{^3 {\}it https://github.com/phonopy/phonopy/tree/master/example/NaCl}$

 $^{^{4} {\}rm https://github.com/phonopy/phonopy/tree/master/example/TiO2-anatase}$

DFPT Interpolation Scheme for Polar Material



Outline

- Lattice Dynamics
 - 1D Atomic Chain
 - 3D lattice
 - How to calculate the dynamical matrix

Density functional perturbation theory



First and Second-order Derivatives of the Total Energy

Within DFT, the total energy is given by

$$E_{\rm tot} = \sum_{i}^{\rm occ} \langle \psi_i(\mathbf{r})| - \frac{\hbar^2}{2m} \nabla^2 |\psi_i(\mathbf{r})\rangle + \int_{\mathcal{V}} V_{\rm ext}(\mathbf{r}) \rho(\mathbf{r}) \, \mathrm{d}\mathbf{r} + E_H[\rho] + E_{xc}[\rho] + U_{II}$$
over the entire space

- $V_{\text{ext}}(\mathbf{r})$: electron-ion interaction
- U_{II}: ion-ion interaction
- $E_H[\rho]$: electron Hartree energy
- $E_{xc}[\rho]$: exchange-correlation energy

• $\rho(\mathbf{r})$: electron density

$$\rho(\mathbf{r}) = \sum_{i}^{occ} \psi_{i}^{*}(\mathbf{r}) \psi_{i}(\mathbf{r})$$

The first and second order derivative with respect to external parameter λ , μ

$$\frac{\partial E_{\text{tot}}}{\partial \lambda} = \int_{V} \frac{\partial V_{\text{ext}}(\mathbf{r})}{\partial \lambda} \rho(\mathbf{r}) \, d\mathbf{r} + \frac{\partial U_{II}}{\partial \lambda}$$
 (8)

$$\frac{\partial^{2} \mathsf{E}_{\mathsf{tot}}}{\partial \lambda \partial \mu} = \int_{V} \frac{\partial^{2} \mathsf{V}_{\mathsf{ext}}(\mathbf{r})}{\partial \lambda \partial \mu} \rho(\mathbf{r}) \, \mathrm{d}\mathbf{r} + \frac{\partial^{2} \mathsf{U}_{II}}{\partial \lambda \partial \mu} + \int_{V} \frac{\partial \mathsf{V}_{\mathsf{ext}}(\mathbf{r})}{\partial \lambda} \frac{\partial \rho(\mathbf{r})}{\partial \mu} \, \mathrm{d}\mathbf{r} \tag{9}$$

2-nd derivative requires the extra calculation of the *linear response* of the charge density to the external perturbation!

Density Functional Perturbation Theory I

The Kohn-Sham eqution

With a small perturbation μ

$$\begin{split} \left[-\frac{\hbar^2}{2m} \nabla^2 + V_{KS}(\boldsymbol{r}; \mu) \right] \psi_n(\boldsymbol{r}) &= \varepsilon_n \psi_n(\boldsymbol{r}) \\ \rho(\boldsymbol{r}) &= \sum_n^{\text{occ}} |\psi_n(\boldsymbol{r})|^2 \\ V_{KS}(\boldsymbol{r}) &= V_{\text{ext}}(\boldsymbol{r}) + V_H(\boldsymbol{r}) + V_{\text{xc}}(\boldsymbol{r}) \end{split}$$

$$V_{KS}(\mathbf{r}; \mu) = V_{KS}(\mathbf{r}; \mu = 0) + \mu \frac{\partial V_{KS}(\mathbf{r})}{\partial \mu}$$
$$\psi_n(\mathbf{r}; \mu) = \psi_n(\mathbf{r}; \mu = 0) + \mu \frac{\partial \psi_n(\mathbf{r})}{\partial \mu}$$
$$\varepsilon_n(\mu) = \varepsilon_n(\mu = 0) + \mu \frac{\partial \varepsilon_n}{\partial \mu}$$

Inserting these equations and keeps only the first-order terms

$$\left[-\frac{\hbar^2}{2m} \nabla^2 + V_{KS}(\mathbf{r}) - \varepsilon_n \right] \frac{\partial \psi_n(\mathbf{r})}{\partial \mu} = -\left[\frac{\partial V_{KS}(\mathbf{r})}{\partial \mu} - \frac{\partial \varepsilon_n}{\partial \mu} \right] \psi_n(\mathbf{r})$$
(10)

where

$$\frac{\partial V_{KS}(\mathbf{r})}{\partial \mu} = \frac{\partial V_{ext}(\mathbf{r})}{\partial \mu} + \frac{\partial V_{H}(\mathbf{r})}{\partial \mu} + \frac{\partial V_{xc}(\mathbf{r})}{\partial \mu}$$
(11)

$$\frac{\partial V_H(\mathbf{r})}{\partial \mu} = \int \frac{1}{|\mathbf{r} - \mathbf{r}'|} \frac{\partial \rho(\mathbf{r}')}{\partial \mu} d\mathbf{r}'; \qquad \frac{\partial V_{xc}(\mathbf{r})}{\partial \mu} = \int \frac{dV_{xc}}{d\rho(\mathbf{r}')} \frac{\partial \rho(\mathbf{r}')}{\partial \mu} d\mathbf{r}'$$
(12)

Density Functional Perturbation Theory II

Electron density response to the perturbation

$$\frac{\partial \rho(\mathbf{r})}{\partial \mu} = \sum_{n}^{\text{occ}} \left[\frac{\partial \psi_{n}^{*}(\mathbf{r})}{\partial \mu} \psi_{n}(\mathbf{r}) + \psi_{n}^{*}(\mathbf{r}) \frac{\partial \psi_{n}(\mathbf{r})}{\partial \mu} \right]$$
(13)

$$\frac{\partial \psi_n(\mathbf{r})}{\partial \mu} = \sum_{m \neq n} \frac{\langle \psi_m | \frac{\partial V_{KS}(\mathbf{r})}{\partial \mu} | \psi_n \rangle}{\varepsilon_n - \varepsilon_m} \psi_m(\mathbf{r})$$
(14)

Where n is the index for the occupied staets and m runs over all the states.

Define $P_v = \sum_n^{\rm occ} |\psi_n\rangle\langle\psi_n|$ as the projector on the valence bands, then $P_c = \mathbb{1} - P_v$ is the projector on the conduction bands

$$\begin{split} \frac{\partial \rho(\boldsymbol{r})}{\partial \mu} &= \sum_{n}^{\text{occ}} \left[P_{c} \frac{\partial \psi_{n}^{*}(\boldsymbol{r})}{\partial \mu} \psi_{n}(\boldsymbol{r}) + \psi_{n}^{*}(\boldsymbol{r}) P_{c} \frac{\partial \psi_{n}(\boldsymbol{r})}{\partial \mu} \right] + \sum_{n}^{\text{occ}} \left[P_{v} \frac{\partial \psi_{n}^{*}(\boldsymbol{r})}{\partial \mu} \psi_{n}(\boldsymbol{r}) + \psi_{n}^{*}(\boldsymbol{r}) P_{v} \frac{\partial \psi_{n}(\boldsymbol{r})}{\partial \mu} \right] \\ &= \sum_{n}^{\text{occ}} \left[P_{c} \frac{\partial \psi_{n}^{*}(\boldsymbol{r})}{\partial \mu} \psi_{n}(\boldsymbol{r}) + \psi_{n}^{*}(\boldsymbol{r}) P_{c} \frac{\partial \psi_{n}(\boldsymbol{r})}{\partial \mu} \right] + \sum_{mn}^{\text{occ}} \psi_{m}^{*}(\boldsymbol{r}) \psi_{n}(\boldsymbol{r}) \left(\left\langle \frac{\partial \psi_{n}}{\partial \mu} | \psi_{m} \right\rangle + \left\langle \psi_{n} | \frac{\partial \psi_{m}}{\partial \mu} \right\rangle \right) \end{split}$$

Due the orthonormality of $\psi_n(\mathbf{r})$, i.e. $\langle \psi_m | \psi_n \rangle = \delta_{mn}$

$$\frac{\partial \rho(\mathbf{r})}{\partial \mu} = \sum_{n}^{\text{occ}} \left[P_{c} \frac{\partial \psi_{n}^{*}(\mathbf{r})}{\partial \mu} \psi_{n}(\mathbf{r}) + \psi_{n}^{*}(\mathbf{r}) P_{c} \frac{\partial \psi_{n}(\mathbf{r})}{\partial \mu} \right]$$
(15)

The electron density response depends only on the component of the perturbation that *couples the occupied* states with the empty ones.

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Density Functional Perturbation Theory III

Apply P_c on the left and right-hand side of Eq. 10

$$\left[-\frac{\hbar^2}{2m} \nabla^2 + V_{KS}(\mathbf{r}) + \frac{\alpha P_v}{\rho} - \varepsilon_n \right] P_c \frac{\partial \psi_n(\mathbf{r})}{\partial \mu} = -(\mathbb{1} - P_v) \frac{\partial V_{KS}(\mathbf{r})}{\partial \mu} \psi_n(\mathbf{r})$$
(16)

where the αP_{ν} is added to make the left-hand-side nonsingular.

$$\frac{\partial V_{KS}(\mathbf{r})}{\partial \mu} = \frac{\partial V_{ext}(\mathbf{r})}{\partial \mu} + \int \frac{1}{|\mathbf{r} - \mathbf{r}'|} \frac{\partial \rho(\mathbf{r}')}{\partial \mu} d\mathbf{r}' + \int \frac{dV_{xc}}{d\rho(\mathbf{r}')} \frac{\partial \rho(\mathbf{r}')}{\partial \mu} d\mathbf{r}'$$
(17)

$$\frac{\partial \rho(\mathbf{r})}{\partial \mu} = \sum_{n}^{\text{occ}} \left[P_{c} \frac{\partial \psi_{n}^{*}(\mathbf{r})}{\partial \mu} \psi_{n}(\mathbf{r}) + \psi_{n}^{*}(\mathbf{r}) P_{c} \frac{\partial \psi_{n}(\mathbf{r})}{\partial \mu} \right]$$

Considering the time-reversal symmetry $\psi_{n,\mathbf{k}}^*(\mathbf{r}) = \psi_{n,\mathbf{k}}(\mathbf{r})$

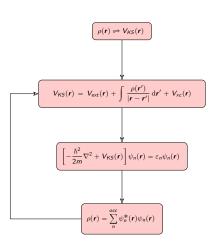
$$\frac{\frac{\partial \rho(\mathbf{r})}{\partial \mu}}{\frac{\partial \rho(\mathbf{r})}{\partial \mu}} = \sum_{n,-\mathbf{k}}^{\text{occ}} P_c \frac{\partial \psi_{n,-\mathbf{k}}^{*}(\mathbf{r})}{\partial \mu} \psi_{n,-\mathbf{k}}(\mathbf{r}) + \sum_{n,\mathbf{k}}^{\text{occ}} \psi_{n,\mathbf{k}}^{*}(\mathbf{r}) P_c \frac{\partial \psi_{n,\mathbf{k}}(\mathbf{r})}{\partial \mu}$$

$$= 2 \sum_{n,-\mathbf{k}}^{\text{occ}} \psi_{n,\mathbf{k}}^{*}(\mathbf{r}) P_c \frac{\partial \psi_{n,\mathbf{k}}(\mathbf{r})}{\partial \mu} \tag{18}$$

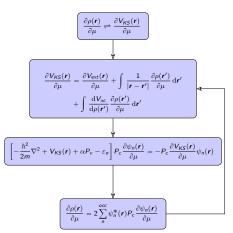
Eq. 16, Eq. 17 and Eq. 18 form a set of self-consistent eugations for the perturbed system.

DFT & DFPT Equations

DFT



DFPT



DFPT — Incommensurate Perturbations I

Add a monochromatic perturbation characterized by generic wavevector \boldsymbol{q} to the periodic potential

$$V_{\text{ext}}(\mathbf{r} + \mathbf{x}_j) = V_{\text{ext}}(\mathbf{r}); \qquad \Delta V_{\text{ext}}^{\mathbf{q}}(\mathbf{r} + \mathbf{x}_j) = e^{i\mathbf{q}\cdot\mathbf{x}_j} \, \Delta V_{\text{ext}}^{\mathbf{q}}(\mathbf{r})$$

It can be shown that all the linear responses are also characterized by wavevector q

$$\Delta V_{KS}^{\mathbf{q}}(\mathbf{r} + \mathbf{x}_j) = e^{i\mathbf{q} \cdot \mathbf{x}_j} \Delta V_{KS}^{\mathbf{q}}(\mathbf{r})$$
$$\Delta \rho^{\mathbf{q}}(\mathbf{r} + \mathbf{x}_j) = e^{i\mathbf{q} \cdot \mathbf{x}_j} \Delta \rho^{\mathbf{q}}(\mathbf{r})$$
$$\Delta \psi_{n,k}^{\mathbf{q}}(\mathbf{r} + \mathbf{x}_j) = e^{i\mathbf{q} \cdot \mathbf{x}_j} \Delta \psi_{n,k}^{\mathbf{q}}(\mathbf{r})$$

Let us define the related periodic quantities

$$\begin{split} \Delta V_{\text{ext}}^{\boldsymbol{q}}(\boldsymbol{r}) &= e^{i\boldsymbol{q}\cdot\boldsymbol{r}} \, \tilde{V}_{\text{ext}}^{\boldsymbol{q}}(\boldsymbol{r}); \\ \Delta V_{KS}^{\boldsymbol{q}}(\boldsymbol{r}) &= e^{i\boldsymbol{q}\cdot\boldsymbol{r}} \, \tilde{V}_{\text{ext}}^{\boldsymbol{q}}(\boldsymbol{r}); \\ \Delta V_{KS}^{\boldsymbol{q}}(\boldsymbol{r}) &= e^{i\boldsymbol{q}\cdot\boldsymbol{r}} \, \tilde{V}_{KS}^{\boldsymbol{q}}(\boldsymbol{r}); \\ \Delta \rho^{\boldsymbol{q}}(\boldsymbol{r}) &= e^{i\boldsymbol{q}\cdot\boldsymbol{r}} \, \tilde{\rho}^{\boldsymbol{q}}(\boldsymbol{r}); \\ \psi_{n,\boldsymbol{k}}(\boldsymbol{r}) &= e^{i\boldsymbol{k}\cdot\boldsymbol{r}} \, u_{n,\boldsymbol{k}}(\boldsymbol{r}); \\ \psi_{n,\boldsymbol{k}}(\boldsymbol{r}) &= e^{i(\boldsymbol{k}+\boldsymbol{q})\cdot\boldsymbol{r}} \, \Delta u_{n,\boldsymbol{k}}^{\boldsymbol{q}}(\boldsymbol{r}); \\ \Delta u_{n,\boldsymbol{k}}^{\boldsymbol{q}}(\boldsymbol{r}+\boldsymbol{x}_j) &= u_{n,\boldsymbol{k}}(\boldsymbol{r}); \\ \Delta u_{n,\boldsymbol{k}}^{\boldsymbol{q}}(\boldsymbol{r}+\boldsymbol{x}_j) &= u_{n,\boldsymbol{k}}(\boldsymbol{r}); \\ \Delta u_{n,\boldsymbol{k}}^{\boldsymbol{q}}(\boldsymbol{r}+\boldsymbol{x}_j) &= \Delta u_{n,\boldsymbol{k}}^{\boldsymbol{q}}(\boldsymbol{r}); \end{split}$$

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Substitute into Ea. 16

$$\left[H_{KS}(\mathbf{r}) + \frac{\alpha P_{\mathbf{v}}}{\epsilon} - \varepsilon_{n,\mathbf{k}}\right] P_{c} e^{i(\mathbf{k}+\mathbf{q}) \cdot \mathbf{r}} \frac{\partial u_{n,\mathbf{k}}^{\mathbf{q}}(\mathbf{r})}{\partial \mu} = -P_{c} e^{i(\mathbf{k}+\mathbf{q}) \cdot \mathbf{r}} \frac{\partial \tilde{V}_{KS}^{\mathbf{q}}(\mathbf{r})}{\partial \mu} u_{n,\mathbf{k}}(\mathbf{r})$$
(19)

Apply the phase $e^{-i(k+q)\cdot r}$ to the both sides of Eq. 19

$$\left[H_{KS}^{k+q}(\mathbf{r}) + \alpha P_{v}^{k+q} - \varepsilon_{n,k}\right] P_{c}^{k+q} \frac{\partial u_{n,k}^{q}(\mathbf{r})}{\partial \mu} = -P_{c}^{k+q} \frac{\partial \tilde{V}_{KS}^{q}(\mathbf{r})}{\partial \mu} u_{n,k}(\mathbf{r})$$
(20)

where

$$\frac{\partial \tilde{V}_{KS}(\mathbf{r})}{\partial \mu} = \frac{\partial \tilde{V}_{ext}(\mathbf{r})}{\partial \mu} + \int \frac{1}{|\mathbf{r} - \mathbf{r}'|} \frac{\partial \tilde{\rho}^{\mathbf{q}}(\mathbf{r}')}{\partial \mu} \, d\mathbf{r}' + \int \frac{dV_{xc}}{d\rho(\mathbf{r}')} \frac{\partial \tilde{\rho}^{\mathbf{q}}(\mathbf{r}')}{\partial \mu} \, d\mathbf{r}'$$

$$\frac{\partial \tilde{\rho}^{\mathbf{q}}(\mathbf{r}')}{\partial \mu} = 2 \sum_{n,k}^{occ} u_{n,k}^{*}(\mathbf{r}) P_{c}^{k+q} \frac{\partial u_{n,k}^{q}(\mathbf{r})}{\partial \mu}$$

$$P_{v}^{k+q} = \sum_{n=0}^{occ} |u_{n,k+q}\rangle \langle u_{n,k+q}|; \quad P_{c}^{k+q} = 1 - P_{v}^{k+q}$$

The treatment of perturbations incommensurate with the unperturbed system periodicity is mapped onto the original periodic system.

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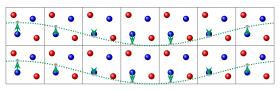
Periodic Displacement of Ions — Phonon Perturbation I

In a solid, the external potential can be expressed as the superposition of atomic potential $v_s(r)$

$$V_{\text{ext}}(\mathbf{r}) = \sum_{j} \sum_{s}^{N_a} v_s(\mathbf{r} - \mathbf{x}_j - \mathbf{d}_s - \mathbf{u}_s^j)$$

Consider the periodic displacement of the ions of the form

$$u_{s\alpha}^{j}(\mathbf{q}) = \mu_{s\alpha} e^{i\mathbf{q}\cdot\mathbf{x}_{j}} + \mu_{s\alpha}^{*} e^{-i\mathbf{q}\cdot\mathbf{x}_{j}} \qquad (\alpha = x, y, z)$$
 (21)



$$\frac{\partial V_{\text{ext}}^{q}(\mathbf{r})}{\partial \mu_{s\alpha}} = -\sum_{j} e^{i\mathbf{q}\cdot\mathbf{x}_{j}} \left. \frac{\partial v_{s}(\mathbf{r} - \mathbf{x}_{j} - \mathbf{d}_{s})}{\partial r_{\alpha}} \right|_{u_{s}^{j} = 0}
= -e^{i\mathbf{q}\cdot\mathbf{r}} \sum_{j} e^{i\mathbf{q}\cdot(\mathbf{x}_{j} - \mathbf{r})} \left. \frac{\partial v_{s}(\mathbf{r} - \mathbf{x}_{j} - \mathbf{d}_{s})}{\partial r_{\alpha}} \right|_{u_{s}^{j} = 0}$$

The colored quantity has the periodicity of the lattice.

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Periodic Displacement of Ions — Phonon Perturbation II

The dynamical matrix

$$D_{s\alpha,t\beta}(\boldsymbol{q}) = \frac{1}{\sqrt{M_s M_t}} \sum_k e^{-i\boldsymbol{q} \cdot \boldsymbol{x}_j} \frac{\partial^2 E_{\text{tot}}^0}{\partial u_{s\alpha}^j \partial u_{t\beta}^k} e^{i\boldsymbol{q} \cdot \boldsymbol{x}_k} = \frac{N}{\sqrt{M_s M_t}} \frac{\partial^2 E_{\text{tot}}^0}{\partial \mu_{s\alpha}^* \partial \mu_{t\beta}}$$

Inserting the equation for the second derivative Eq. 9

$$\begin{split} D_{s\alpha,t\beta}(\boldsymbol{q}) &= \frac{N}{\sqrt{M_s M_t}} \int_{V} \sum_{k} e^{-i\boldsymbol{q}\cdot\boldsymbol{x}_j} \frac{\partial^2 V_{\text{ext}}^{\boldsymbol{q}}(\boldsymbol{r})}{\partial u_{s\alpha}^i \partial u_{t\beta}^k} \, \rho(\boldsymbol{r}) \, e^{i\boldsymbol{q}\cdot\boldsymbol{x}_k} \, \mathrm{d}\boldsymbol{r} \\ &+ \frac{1}{\sqrt{M_s M_t}} \int_{V} \sum_{k} e^{-i\boldsymbol{q}\cdot\boldsymbol{x}_j} \frac{\partial V_{\text{ext}}^{\boldsymbol{q}}(\boldsymbol{r})}{\partial u_{s\alpha}^i \partial u_{t\beta}^k} \, \frac{\partial \rho(\boldsymbol{r})}{\partial u_{t\beta}^k} \, e^{i\boldsymbol{q}\cdot\boldsymbol{x}_k} \, \mathrm{d}\boldsymbol{r} \end{split}$$

Thank you!