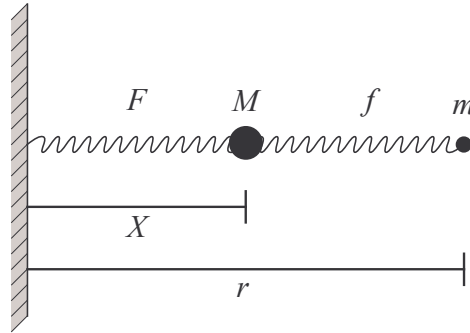


**Problem 1: Dynamics of two coupled masses**

**(4 points)**

Consider the Hamilton function of two particles in one dimension:



$$H = \frac{P^2}{2M} + \frac{p^2}{2m} + \frac{1}{2}F X^2 + \frac{1}{2}f (X - r)^2 .$$

- a) Calculate the frequencies of the normal modes of the classical system.
- b) Specify the Hamilton operator and calculate the eigenvalues of the corresponding quantum mechanical systems.
- c) Suppose that the mass  $m$  is much smaller than  $M$ . Calculate the eigenvalues of  $\hat{H}$  along the lines of the “Born-Oppenheimer approximation“: in a first step neglect the motion of the heavy mass  $M$  and calculate the eigenvalues of the corresponding Hamiltonian. Thereafter solve the problem of the motion of the heavy mass  $M$  in the “effective potential“, calculated in the previous step.
- d) Compare your results from b) and c). It is advantageous to use the abbreviations

$$\omega = \sqrt{\frac{f}{m}} , \quad \Omega = \sqrt{\frac{F}{M}} \quad \text{and} \quad \kappa = \sqrt[4]{\frac{m}{M}} .$$

**Problem 2: Fourier transformation of Coulomb potential**

**(2 points)**

- a) Calculate the Fourier transform

$$\tilde{v}(\vec{q}) = \frac{1}{\Omega} \int_{\Omega} v(\vec{r}) e^{-i\vec{q}\cdot\vec{r}} d^3 r \quad \text{of the Yukawa potential} \quad v(\vec{r}) = \frac{e^2}{4\pi\epsilon_0} \frac{1}{r} e^{-\gamma r}$$

in a Born-von Kármán cell with volume  $\Omega$ .  $\gamma$  is a real constant. Replace the integral  $\int_{\Omega} d^3 r$  by an integration over a sphere with radius  $R \rightarrow \infty$ .

- b) The Coulomb potential is the  $\gamma \rightarrow 0$  limit of the Yukawa potential. Determine the Fourier transform of the Coulomb potential.

**Problem 3: Ewald-Method****(4 points)**

The calculation of the electrostatic interaction energy  $U_{NN}$  between the nuclei in a solid requires the evaluation of a sum over all lattice vectors  $\vec{R}_j$ . A direct summation is numerically very demanding due to the long-range Coulomb potential. In the method of Ewald, the Coulomb potential is represented by error functions  $\text{erf}(x)$  and  $\text{erfc}(x)$ . One part of the sum is evaluated in Fourier space while the other part is carried out in real space.

a) Show that the sum  $S = \sum_j' \frac{1}{|\vec{a} - \vec{R}_j|}$  can be written as  $S = S_1 + S_2$  with

$$S_1 = \sum_j' \frac{2}{\sqrt{\pi}} \int_0^\eta e^{-|\vec{a} - \vec{R}_j|^2 x^2} dx, \quad S_2 = \sum_j' \frac{2}{\sqrt{\pi}} \int_\eta^\infty e^{-|\vec{a} - \vec{R}_j|^2 x^2} dx.$$

Thereby is  $\vec{a}$  the difference of two basis vectors  $\vec{\tau}_\nu - \vec{\tau}_{\nu'}$ ,  $\vec{R}_j$  is a lattice vector and  $\eta$  is a real number between 0 and  $\infty$  which is chosen in an appropriate way for the numerical evaluation. The prime at the sum indicates that the term  $\vec{R}_j = \vec{a}$  is excluded from the summation.

b) The case  $\vec{R}_j = \vec{a}$  can only occur for  $\vec{a} = 0$ . Show that  $S_1$  can be written as a sum without restriction in the form

$$S_1 = \sum_j \left( \frac{2}{\sqrt{\pi}} \int_0^\eta e^{-|\vec{a} - \vec{R}_j|^2 x^2} dx \right) - \delta_{\vec{a}, \vec{0}} \cdot \frac{2\eta}{\sqrt{\pi}}.$$

c) Represent in  $S_1$  the Gaussian function by its Fourier transform, evaluate the sum over  $j$  and integrate with respect to the variable  $x$ .

d) Write the sum  $S_2$  in terms of  $\text{erfc}(x)$ .

e) Why is the evaluation of  $S_1$  and  $S_2$  as outlined above more efficient than a direct summation in  $S$ ?

*Remark:* The divergent term in  $S_1$  with  $|\vec{G}| = 0$  is compensated by analogue terms in the electron-electron and electron-nucleus interaction.

Useful relations:

i) 
$$\int_0^\infty e^{-b^2 x^2} dx = \frac{\sqrt{\pi}}{2b}$$

ii) 
$$\text{erf}(y) = \frac{2}{\sqrt{\pi}} \int_0^y e^{-t^2} dt$$

iii) 
$$\text{erfc}(y) = 1 - \text{erf}(y)$$

iv) 
$$\sum_j e^{-|\vec{a} - \vec{R}_j|^2 x^2} = \sum_j \frac{1}{\Omega} \frac{\pi^{\frac{3}{2}}}{x^3} \sum_{\vec{G}} e^{i\vec{G} \cdot (\vec{a} - \vec{R}_j)} \cdot e^{-\frac{G^2}{4x^2}}.$$

$\Omega$  is the volume of the Born-von Kármán cell. There are  $N_{\text{cell}}$  lattice vectors.  $\Omega_0 = \Omega/N_{\text{cell}}$  is the volume of the unit cell.