# Computer Physics

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## Project 1

In this project I calculate the energy spectrum of  $H=H_0+\lambda V$  where  $H_0$  is the harmonic oscillator and  $V=\hbar\omega(\frac{x}{a})^4$  where  $a=\sqrt{\frac{\hbar}{m\omega}}$  is the natural length of the system.

This is achieved by adding the matrices of  $H_0$  and V in the basis of the eigenvectors of  $H_0$  and then truncating the basis. We know the matrix representation of  $\frac{x}{a}$  in this basis is

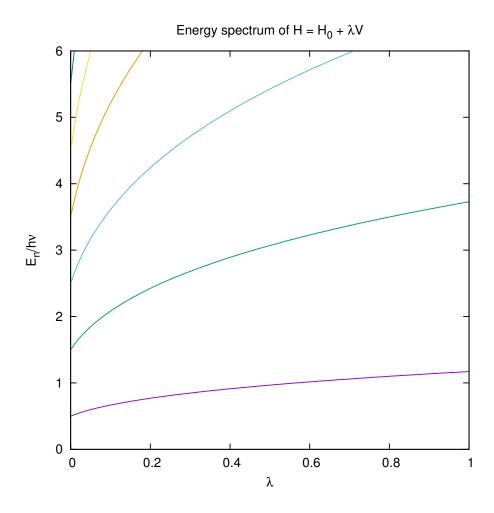
$$\mathbf{z} = \frac{1}{2} \begin{bmatrix} 0 & 2 & 0 & 0 & \dots \\ 2 & 0 & \sqrt{6} & 0 & \dots \\ 0 & \sqrt{6} & 0 & \sqrt{8} & \dots \\ 0 & 0 & \sqrt{8} & 0 & \dots \\ \vdots & \vdots & \vdots & \vdots & \ddots \end{bmatrix}$$

and the matrix for V is simply  $\hbar\omega x^4$ . We add this (multiplied by some  $\lambda$ ) to

$$H_0 = \hbar\omega \begin{bmatrix} 0.5 & 0 & 0 & 0 & \dots \\ 0 & 1.5 & 0 & 0 & \dots \\ 0 & 0 & 2.5 & 0 & \dots \\ 0 & 0 & 0 & 3.5 & \dots \\ \vdots & \vdots & \vdots & \vdots & \ddots \end{bmatrix}$$

and then truncate. In our case we used a basis of 128 eigenvectors of  $H_0$ .

The eigenvalues of the resulting matrix are of course the possible values of energy of the system. The following is a graph of the first six eigenvalues as a function of  $\lambda$ .



### Project 2

We continue with the harmonic oscillator  $H_0$  but this time we take a look at  $H = H_0 + H'(t)$  where  $H'(t) = \hbar\Omega(a^{\dagger} + a)\theta(t)$  where a is the ladder operator and  $\theta(t)$  is the Heaviside step function. We again use the energy basis of the harmonic oscillator. Then the matrix for a is

$$\begin{bmatrix} 0 & 1 & 0 & 0 & \dots \\ 0 & 0 & \sqrt{2} & 0 & \dots \\ 0 & 0 & 0 & \sqrt{3} & \dots \\ 0 & 0 & 0 & 0 & \dots \\ \vdots & \vdots & \vdots & \vdots & \ddots \end{bmatrix}$$

so the matrix for  $a^{\dagger} + a$  is

$$\begin{bmatrix} 0 & 1 & 0 & 0 & \dots \\ 1 & 0 & \sqrt{2} & 0 & \dots \\ 0 & \sqrt{2} & 0 & \sqrt{3} & \dots \\ 0 & 0 & \sqrt{3} & 0 & \dots \\ \vdots & \vdots & \vdots & \vdots & \ddots \end{bmatrix}$$

Now given some initial state of the density operator  $\rho$ , for example the ground state of the harmonic oscillator:

$$\rho = \begin{bmatrix} 1 & 0 & 0 & 0 & \dots \\ 0 & 0 & 0 & 0 & \dots \\ 0 & 0 & 0 & 0 & \dots \\ 0 & 0 & 0 & 0 & \dots \\ \vdots & \vdots & \vdots & \vdots & \ddots \end{bmatrix}$$

we can add these matrices in a truncated basis and then calculate the time evolution of the system using the Liouville-von Neumann equation:

$$i\hbar\dot\rho(t)=[H(t),\rho(t)]$$
 i.e.  $\dot\rho(t)=\frac{1}{\hbar}\Lambda[\rho(t)],\,\Lambda[\rho(t)]=-i[H(t),\rho(t)]$ 

or, more specifically, using iteration of the following Crank-Nicolson approximation of the Liouville-von Neumann equation:

$$\rho(t_{n+1}) = \rho(t_n) + \frac{\Delta t}{2\hbar} (\Lambda[\rho(t_n)] + \Lambda[\rho(t_{n+1})])$$

for some timegrid where  $t_0=0$  and  $t_{n+1}=t_n+\Delta t$ . In the graph below we show the time evolution of the occupation of the five lowest states assuming the initial state is |0>, the timestep is  $\Delta t=1fs$  and that  $\hbar\omega=\hbar\Omega=1meV$ .

