

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

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

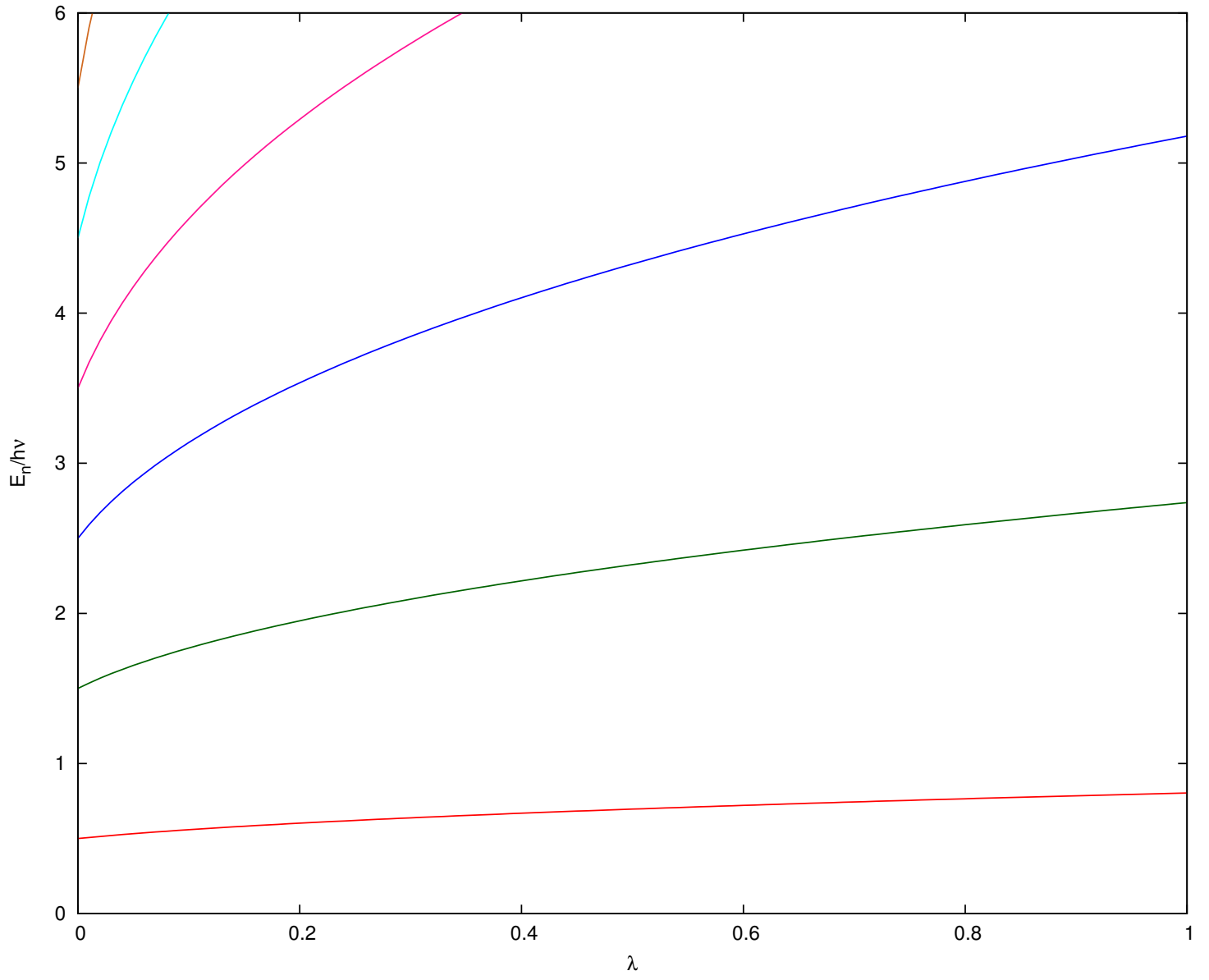
and the matrix for V is simply $\hbar\omega x^4$. We add this (multiplied by some λ) to

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

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 λ .

Energy spectrum of $H = H_0 + \lambda V$



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 ρ , 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:

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

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 $|0\rangle$, the timestep is $\Delta t = 1fs$ and that $\hbar\omega = \hbar\Omega = 1meV$.

Time evolution of the occupation of the five lowest states of $H = H_0 + H'(t)$

