

CHM452: Problem Set 2

Xinxian Chen*

March 24, 2020

Let $i = \sqrt{-1}$, $o(\cdot)$ and $O(\cdot)$ are little-o and big-O notations, respectively, and \mathbb{N} is the set of natural numbers ($\mathbb{N} = \{0, 1, 2, 3, \dots\}$) and $\mathbb{N}_+ = \mathbb{N} \setminus \{0\}$

1. (i) Since $\Psi(x, t)$ is a linear combination of $\Phi_n(x, t)$, where

$$\Phi_n(x, t) = \sqrt{\frac{2}{w}} \sin\left(\frac{n\pi}{w(t)}x\right) \exp\left(\frac{i(mvx^2 - 2E_n^i at)}{2\hbar w(t)}\right),$$



$w(t) = a + vt$, and $E_n^i = n^2\pi^2\hbar^2/2ma^2$, we only need to check that $\Phi_n(x, t)$ satisfies the boundary condition for all $n \in \mathbb{N}_+$

$$\Phi_n(x = 0, t) = 0, \quad \Phi_n(x = w(t), t) = 0.$$

the problem also ask to check if this wavefunction satisfies TDSE

On the other hand,

$$\begin{aligned} \Phi_n(x = 0, t) &= \sqrt{\frac{2}{w}} \sin\left(\frac{n\pi}{w(t)} \cdot 0\right) \exp\left(\frac{i(-2E_n^i at)}{2\hbar w(t)}\right) = 0, \\ \Phi_n(x = w(t), t) &= \sqrt{\frac{2}{w}} \sin\left(\frac{n\pi}{w(t)}w(t)\right) \exp\left(\frac{i(mvw^2 - 2E_n^i at)}{2\hbar w(t)}\right) \\ &= \sqrt{\frac{2}{w}} \sin(n\pi) \exp\left(\frac{i(mvw^2 - 2E_n^i at)}{2\hbar w(t)}\right) = 0. \end{aligned}$$



Therefore, $\Phi_n(x, t)$ satisfies the boundary conditions, and hence, $\Psi(x, t)$ satisfies the boundary conditions.

- (ii) At $t = 0$,

$$\begin{aligned} c_n &= \int_0^a \Phi_n(x, 0)^* \Psi(x, 0) dx = \int_0^a \left(\sqrt{\frac{2}{a}} \sin\left(\frac{n\pi}{a}x\right) e^{\frac{imvx^2}{2\hbar a}} \right)^* \sqrt{\frac{2}{a}} \sin\left(\frac{\pi x}{a}\right) dx \\ &= \frac{2}{a} \int_0^a \sin\left(n\frac{\pi x}{a}\right) e^{-\frac{imva}{2\pi^2\hbar}(\frac{\pi x}{a})^2} \sin\left(\frac{\pi x}{a}\right) dx. \end{aligned}$$

here, you need to prove that it is an orthonormal basis



Let $z = \pi x/a$, then

$$c_n = \frac{2}{a} \int_0^\pi \sin(nz) e^{-\frac{imva}{2\pi^2\hbar}z^2} \sin(z) \frac{dx}{dz} dz = \frac{2}{\pi} \int_0^\pi e^{-i\alpha z^2} \sin(nz) \sin(z) dz,$$



where $\alpha = \frac{mva}{2\pi^2\hbar}$.

- (iii) Without loss of generality, suppose the initial time is at $t = 0$ and the final time is at $t = \Delta t$. For the instantaneous eigenstate $\psi_n(x, t)$ at time t , it satisfies the TISE at some specific time

$$-\frac{\hbar^2}{2m} \frac{d^2}{dx^2} \psi_n(x, t) = E_n(t) \psi_n(x, t),$$



*Email: xchen106@ur.rochester.edu

and it satisfies the boundary conditions at time t :

$$\psi_n(x = 0, t) = 0, \quad \psi_n(x = w(t), t) = 0. \quad \checkmark$$

Solve the ODE and we have that $\psi_n(x, t) = N \sin\left(\frac{n\pi x}{w(t)}\right)$ and $E_n(t) = n^2\pi^2\hbar^2/2mw(t)^2$; if we require the instantaneous eigenstates are orthonormal, *i.e.*, $\int_0^{w(t)} \psi_n(x, t)^* \psi_m(x, t) dx = \delta_{nm}$, then we have that $N = \sqrt{2/w(t)}$, and hence,

$$\psi_n(x, t) = \sqrt{\frac{2}{w(t)}} \sin\left(\frac{n\pi x}{w(t)}\right). \quad \checkmark$$

Now we can use the instantaneous eigenstates to represent the solution

$$\Psi(x, t) = \sum_{n=1}^{\infty} a_n(t) \psi_n(x, t) e^{i\theta_n(t)}, \quad \checkmark$$

where $\theta_n = -\int_0^t E_n(t')/\hbar dt'$. From the adiabatic theorem, we have that

$$\begin{aligned} \frac{d}{dt} a_n(t) &\approx -a_n(t) \int_0^{w(t)} \psi_n(x, t) \frac{d}{dt} \psi_n(x, t) dx \\ &= -a_n(t) \frac{2}{w(t)} \frac{n\pi}{w(t)} \int_0^{w(t)} \sin\left(\frac{n\pi x}{w(t)}\right) \cos\left(\frac{n\pi x}{w(t)}\right) dx = 0. \end{aligned} \quad \checkmark$$

On the other hand, at $t = 0$, we have that $a_n(0) = \delta_{1n}$. Therefore,

$$a_n(t) = \delta_{1n}. \quad \checkmark$$

Note that the well expands to twice its original time in a time Δt , which means $v = a/\Delta t$. Therefore,

$$\Psi(x, t) = a_1(t) \psi_1(x, t) e^{i\theta_1(t)} = e^{i\theta(t)} \sqrt{\frac{2}{w(t)}} \sin\left(\frac{\pi x}{w(t)}\right), \quad \checkmark$$

where $w(t) = a(1 + t/\Delta t)$ and

$$\theta(t) = -\frac{1}{\hbar} \int_0^t \frac{\pi^2 \hbar^2}{2ma^2(1+t'/\Delta t)^2} dt' = -\frac{\pi^2 \hbar}{2ma^2} \frac{t}{1+t/\Delta t}. \quad \checkmark$$

On the other hand, if the velocity of expansion v is small enough, *i.e.*, $v \rightarrow 0$, from the exact solution, we will have that $\alpha = mva/2\pi^2\hbar \approx 0$ and

$$c_n(t) \approx \frac{2}{\pi} \int_0^\pi \sin(nz) \sin(z) dz = \delta_{1n}. \quad \checkmark$$

Hence,

$$\begin{aligned} \Psi(x, t) &= \sum_{i=1}^{\infty} c_n(t) \Phi_n(x, t) \approx \Phi_1(x, t) = \sqrt{\frac{2}{w}} \sin\left(\frac{\pi}{w(t)} x\right) \exp\left(\frac{i(mvx^2 - 2E_1^i at)}{2\hbar w(t)}\right) \\ &= e^{i\theta'(t)} \sqrt{\frac{2}{w}} \sin\left(\frac{\pi}{w(t)} x\right), \end{aligned} \quad \checkmark$$

where $\theta'(t) = \frac{mvx^2 - 2E_1^i at}{2\hbar w(t)} \approx -\frac{\pi^2 \hbar^2 at}{2ma^2 \hbar (1+t/\Delta t)} = -\frac{\pi^2 \hbar at}{2ma^2 (1+t/\Delta t)}$, which is consistent with the result from the adiabatic theorem at the limit of $v \rightarrow 0$.

- (iv) If the expansion of the well can be considered as a sudden perturbation, that is, the stationary state does not change as t goes from 0 to Δt , that is, $\Psi(x, \Delta t) \approx \Psi(x, 0) = \sqrt{2/a} \sin(\pi x/a)$;

On the other hand, since $[H(t_1), H(t_2)] = 0$ for all $t_1, t_2 \in [0, \Delta t]$, ✓

$$\Psi(x, \Delta t) = e^{-i\bar{H}\Delta t/\hbar} \sum_{n=1}^{\infty} c_n \Phi_n(x, 0) = e^{-iE^i t/\hbar} \sum_{n=1}^{\infty} c_n e^{-i(E_n - E^i)t/\hbar} \Phi_n(x, 0).$$

can you give me more details how you get the last equation?

where E^i is the energy at initial state. Notice that only terms satisfying $t > \hbar/(E_n - E^i)$ will make appreciable change, and for the sudden perturbation limit, we hence require $t \ll \hbar/(E_n - E^i)$ for all n , that is, $t \ll \hbar/\Delta\bar{H}$.

2. Landau-Zener theory gives a formula that predicts the nonadiabatic transit possibility. If the hamiltonian is

$$\begin{aligned} H &= H_0 + V, \\ H_0 &= E_1(t) |1\rangle\langle 1| + E_2(t) |2\rangle\langle 2|, \\ V &= \frac{\Delta}{2} |1\rangle\langle 2| + \frac{\Delta}{2} |2\rangle\langle 1|, \end{aligned}$$

where $E_1(t) = vt/2$, $E_2(t) = -vt/2$, and we start with $|\psi(t = -\infty)\rangle = |1\rangle$, then for $t \rightarrow +\infty$, Laudau-Zener theory gives the nonadiabatic transit possibility



$$P_{1 \rightarrow 2} = 1 - e^{\frac{\Delta^2 \pi}{2\hbar v}}.$$

can you give me more details with regard to this?

For instance, for the molecular dynamics Laudau-Zener indicates that the nonadiabatic transitions take place only when the surface gap attains a local minimum for each classical trajectory, which gives the theoretical basis of surface hopping algorithms. ✓

3. (i) The Floquet Hamiltonian is

$$\mathcal{H} = H(t) - i\hbar \frac{\partial}{\partial t}$$

The Floquet Hamiltonian is defined on the expanded Hilbert space that spanned by basis $\{|\alpha n\rangle\} := \{|\alpha\rangle|n\rangle\}$, where $|\alpha\rangle \in \{|E_1\rangle, |E_2\rangle\}$ and $\langle t|n\rangle = e^{int\omega}$. The Floquet mode $|U_\lambda\rangle$ satisfies ✓

$$\mathcal{H}|U_\lambda\rangle = \mathcal{E}_\lambda|U_\lambda\rangle.$$

Let

$$\langle \beta t|U_\lambda\rangle = \sum_{m=-\infty}^{\infty} U_{\beta\lambda}^{(m)} e^{im\omega t},$$

$$\langle \alpha|H(t)|\beta\rangle = \sum_{n=-\infty}^{\infty} H_{\alpha\beta}^{(n)} e^{in\omega t}.$$

Then,

$$\langle \alpha m|\mathcal{H}|\beta n\rangle = \sum_k H_{\alpha\beta}^{(k)} \frac{1}{T} \int_0^T e^{i(k-m+n)\omega t} + n\hbar\omega\delta_{\alpha\beta}\delta_{mn} = H_{\alpha\beta}^{(m-n)} + n\hbar\omega\delta_{\alpha\beta}\delta_{mn}$$

where $T = 2\pi/\omega$. On the other hand, from the matrix representation of $H(t)$ we have that ✓

$$\begin{aligned} H^{(0)} &= \begin{pmatrix} -\hbar\omega_0/2 & \\ & \hbar\omega_0/2 \end{pmatrix}, \\ H^{(-1)} &= \begin{pmatrix} & b \\ b & \end{pmatrix}, \\ H^{(1)} &= \begin{pmatrix} & b \\ b & \end{pmatrix}, \\ H^{(n)} &= 0, \text{ otherwise.} \end{aligned}$$

where we have applied the fact that $\cos \omega t = (e^{i\omega t} + e^{-i\omega t})/2$. Therefore, the matrix representation of \mathcal{H} under basis $\{|\alpha n\rangle\}$ is

$$\begin{pmatrix} \ddots & \ddots & & & & \\ \ddots & A - \hbar\omega & B & & & \\ & B & A & B & & \\ & & B & A + \hbar\omega & \ddots & \\ & & & & \ddots & \ddots \end{pmatrix}, \quad \checkmark$$

where $A = \frac{\hbar\omega_0}{2} \begin{pmatrix} -1 & \\ & 1 \end{pmatrix}$ and $B = b \begin{pmatrix} & 1 \\ 1 & \end{pmatrix}$. ✓

Truncate the basis by $\{|\alpha n\rangle\}_{n=-100}^{100}$ and diagonalize the representation matrix of \mathcal{H} numerically, and note the not very that from the original structure of \mathcal{H} , if \mathcal{E} is its eigenenergy then $\mathcal{E} + \hbar\omega$ is also its eigenenergy, then we only clear need to pick the one as quasienergy in a equivalence class that is defined by \sim , where $A \sim B$ if and only if there about this is an integer n such that $A = B + n\hbar\omega$. Here I chose the eigenvalues of the truncated matrix that most close to statement. 0 to represent the that quasienergies of the infinite matrix. The result is in Figure 1

What is A
and B
here.

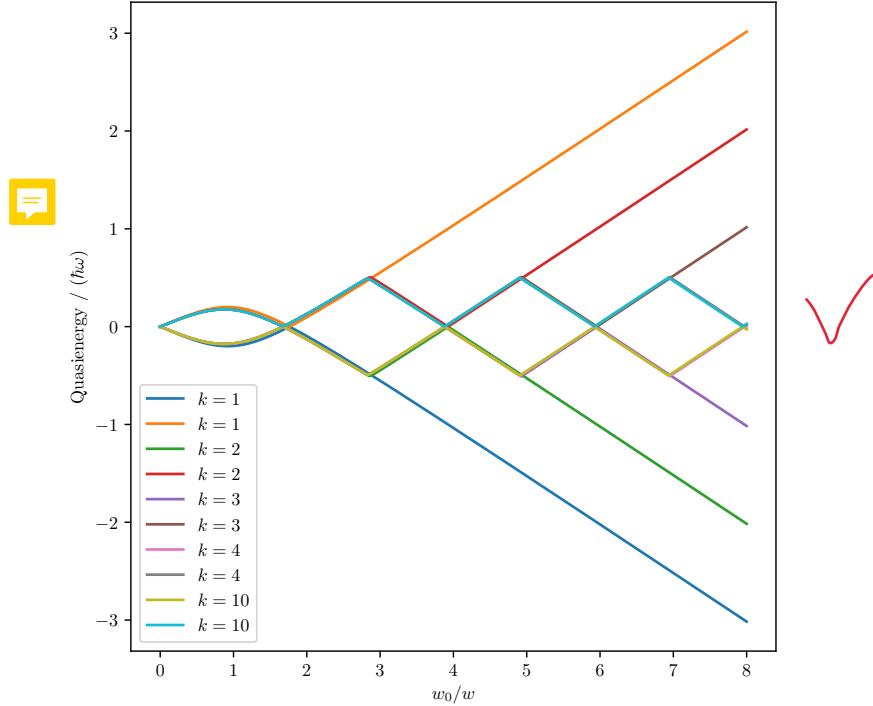


Figure 1: Quasienergies vs. ω_0

Note that the off-diagonal coupling mixes these states, and the non-zero off-diagonal terms will directly mix the the dressed states which differ in energy by $\hbar\omega$ and When $0 < \omega_0/\omega \lesssim 2$, $0 < \Delta$ to make $|E_1\rangle$ mix with $|E_2\rangle$ we need only one Floquet Brillouin zone since the only the dressed $|E_1\rangle$; when $2 \lesssim \omega_0/\omega \lesssim 4$, we need 2 Floquet Brillouin zone; and so on. This can be seen by selecting specific submatrix to diagonalize for the whole matrix, i.e., change the truncated basis by $\{|\alpha n\rangle\}_{n=-k}^k$, and the results are also plotted in Figure 1.

I think I understand your point, but this sentence is not clear to me

- (ii) The time averaged transition probability in Eq. 19 in Shirley 1965 can be written in our notations as

$$\bar{P}_{\alpha \rightarrow \beta} = \sum_n \sum_{\lambda} |\langle \beta n | U_{\lambda} \rangle \langle U_{\lambda} | \alpha 0 \rangle|^2.$$
✓

This can be justified from the definition of $\bar{P}_{\alpha \rightarrow \beta}$

$$P_{\alpha \rightarrow \beta} = |U_{\beta \alpha}(0, t)|^2 = \left| \langle \beta | e^{-iHt/\hbar} | \alpha \rangle \right|^2.$$
✓

On the other hand, note that $\langle t | \alpha 0 \rangle = |\alpha\rangle$ for all t . Insert the spectrum decomposition of identity operator in the space spanned by $\{|n\rangle\}$ and we have

$$\langle \beta | U(0, t) | \alpha \rangle = \langle \beta | \sum_n |n\rangle \langle n | U \langle t | \alpha 0 \rangle = \sum_n \langle t | n \rangle \langle \beta n | U | \alpha 0 \rangle = \sum_n e^{in\omega t} \langle \beta n | \sum_{\lambda} |U_{\lambda}\rangle e^{-i\varepsilon_{\lambda} t/\hbar} \langle U_{\lambda} | \alpha 0 \rangle.$$
✓

Therefore,

$$P_{\alpha \rightarrow \beta} = \left| \sum_n \sum_{\lambda} e^{i(n\omega - \varepsilon_{\lambda}/\hbar)t} \langle \beta n | U_{\lambda} \rangle \langle U_{\lambda} | \alpha 0 \rangle \right|^2$$
✓

If we take the time average, then contributions of the crossing phases are zeros, and the diagonal phases are ones. Hence

$$\bar{P}_{\alpha \rightarrow \beta} = \sum_n \sum_{\lambda} |\langle \beta n | U_{\lambda} \rangle \langle U_{\lambda} | \alpha 0 \rangle|^2.$$
✓

- (iii) See Figure 2. Here the size of the truncated basis k is 10 and has been verified consistency with the case that $k = 100$.

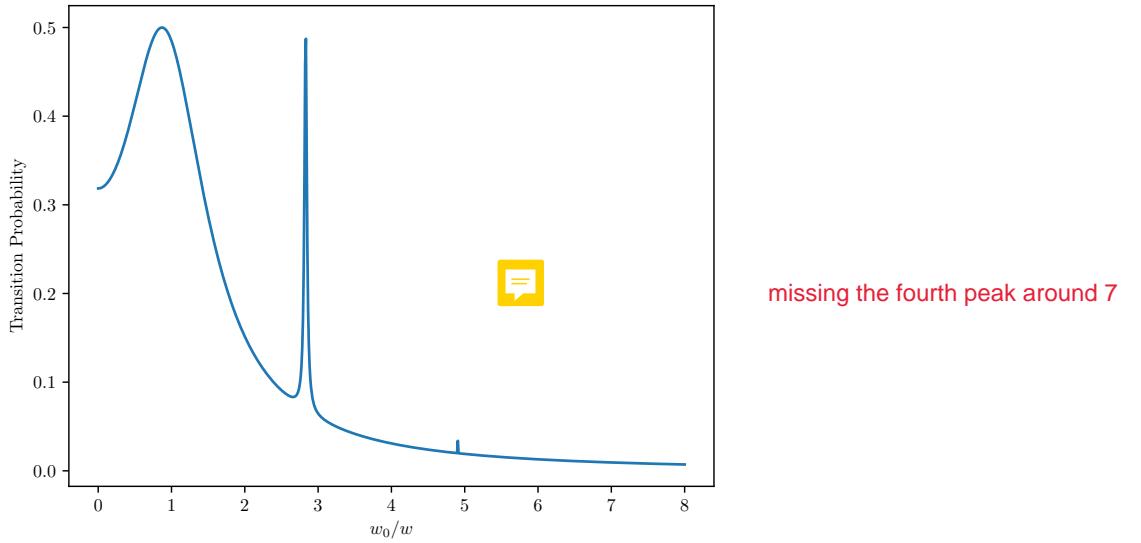


Figure 2: Transition probability vs. ω_0/ω

- (iv) It is clear that there are peaks in Figure 2 at $\omega_0/\omega \approx 1, 3, 5$, and the peaks are almost invisible as ω_0/ω gets larger. This is because when ω_0/ω is nearly odd, the system is more likely to absorb correct amount of energy in laser and then the transition happens, while when ω_0/ω is nearly even, since the Floquet modes share almost the same energy, it is not that likely to absorb correct amount of energy in laser.

the peaks appear because of the avoid crossing points created at that frequencies.

4. (i) Let $\{|n\rangle\}_{n=1}^{\infty}$ be an orthonormal and complete basis. Then

$$\begin{aligned}\text{tr}(\hat{A}\hat{B}) &= \sum_n \langle n|\hat{A}\hat{B}|n\rangle = \sum_n \langle n| \hat{A} \sum_m |m\rangle \langle m| \hat{B} |n\rangle = \sum_n \sum_m \langle n|\hat{A}|m\rangle \langle m|\hat{B}|n\rangle \\ &= \sum_m \sum_n \langle m|\hat{B}|n\rangle \langle n|\hat{A}|m\rangle = \sum_m \langle m|\hat{B}\hat{A}|m\rangle = \text{tr}(\hat{B}\hat{A});\end{aligned}$$

and

$$\langle \hat{A} \rangle = \langle \Psi | \hat{A} | \Psi \rangle = \langle \Psi | \hat{A} \sum_n |n\rangle \langle n| \Psi \rangle = \sum_n \langle \Psi | \hat{A} |n\rangle \langle n| \Psi \rangle = \sum_n \langle n| \Psi \rangle \langle \Psi | \hat{A} |n\rangle = \sum_n \langle n| \hat{\rho} \hat{A} |n\rangle = \text{tr}(\hat{\rho} \hat{A}).$$

how about the general case, here you assume you have a pure state

- (ii) For $|E_1\rangle$:

$$\rho_1 = (1 \ 0)^\dagger (1 \ 0) = \begin{pmatrix} 1 & 0 \\ 0 & 0 \end{pmatrix};$$

for $|E_2\rangle$:

$$\rho_2 = (0 \ 1)^\dagger (0 \ 1) = \begin{pmatrix} 0 & 0 \\ 0 & 1 \end{pmatrix};$$

for $(|E_1\rangle + i|E_2\rangle)/\sqrt{2}$:

$$\rho_3 = \frac{1}{2} \begin{pmatrix} 1 \\ i \end{pmatrix} (1 \ -i) = \frac{1}{2} \begin{pmatrix} 1 & -i \\ i & 1 \end{pmatrix};$$

for a mixture with P_1 to be in $|E_1\rangle$ and P_2 to be in $(|E_1\rangle + i|E_2\rangle)/\sqrt{2}$:

$$\rho_4 = P_1 \rho_1 + P_2 \rho_3 = \begin{pmatrix} P_1 + P_2/2 & -iP_2/2 \\ iP_2/2 & P_2/2 \end{pmatrix};$$

- (iii) Consider a two-level system $\rho = \sum_n P_n |\phi_n\rangle\langle\phi_n|$, where $|\phi_n\rangle = c_{n1} |E_1\rangle + c_{n2} |E_2\rangle$ is normalized, and $|E_1\rangle$ and $|E_2\rangle$ construct an orthonormal basis of the Hilbert space, P_n is the (positive) possibility of finding the system in $|\phi_n\rangle$. It is clear that ρ is hermite since $\rho^\dagger = \sum_n P_n^* |\phi_n\rangle\langle\phi_n| = \sum_n P_n |\phi_n\rangle\langle\phi_n| = \rho$. And hence, we can diagonalize ρ such that

$$\rho = w_1 |\psi_1\rangle\langle\psi_1| + w_2 |\psi_2\rangle\langle\psi_2|,$$

where w_1 and w_2 are nonnegative, and $\{|\psi_1\rangle, |\psi_2\rangle\}$ spans the same Hilbert space as $\{|E_1\rangle, |E_2\rangle\}$, and is orthonormal. Note that

$$\text{tr } \rho = w_1 + w_2 = 1;$$

and therefore, $\rho^2 = w_1^2 |\psi_1\rangle\langle\psi_1| + w_2^2 |\psi_2\rangle\langle\psi_2|$

$$\text{tr } \rho^2 = w_1^2 + w_2^2 = w_1^2 + (1 - w_1)^2 = 2w_1^2 - 2w_1 + 1.$$

Since $0 \leq w_1 \leq 1$, therefore, $1/2 \leq \text{tr } \rho^2 \leq 1$. Note that when $w_1 = w_2 = 1/2$, $\text{tr } \rho^2$ gets its minimal $1/2$ and when $w_1 = 0$ or $w_1 = 1$, $\text{tr } \rho^2$ gets its maximal 1 .

The entropy of such a two-level system is

$$\begin{aligned}S &= -\text{tr}(\rho \ln \rho) = -\text{tr}((w_1 |\psi_1\rangle\langle\psi_1| + w_2 |\psi_2\rangle\langle\psi_2|)(\ln w_1 |\psi_1\rangle\langle\psi_1| + \ln w_2 |\psi_2\rangle\langle\psi_2|)) \\ &= -\text{tr}(w_1 \ln w_1 |\psi_1\rangle\langle\psi_1| + w_2 \ln w_2 |\psi_2\rangle\langle\psi_2|) = -w_1 \ln w_1 - w_2 \ln w_2 = -w_1 \ln w_1 - (1 - w_1) \ln(1 - w_1),\end{aligned}$$

where $0 \leq w_1 \leq 1$. Note that when $w_1 = w_2 = 1/2$, S gets its maximal $\ln 2$ and when $w_1 = 0$ or $w_1 = 1$, $\text{tr } \rho^2$ gets its minimal 0 .

We can relate the two quantities by defining the Rényi entropy

$$R_\alpha = \frac{1}{1-\alpha} \ln \text{tr } \rho^\alpha,$$

and notice that

$$R_2 = \frac{1}{1-2} \ln \text{tr} \rho^2 = -\ln \text{tr} \rho^2, \quad \checkmark$$

and

$$R_1 = \lim_{\alpha \rightarrow 1} \frac{1}{1-\alpha} \ln \text{tr} \rho^\alpha = -\lim_{\alpha \rightarrow 1} \frac{\text{tr} \rho^\alpha \ln \rho}{\text{tr} \rho^\alpha} = -\frac{\text{tr} \rho \ln \rho}{\text{tr} \rho} = -\text{tr} \rho \ln \rho = S. \quad \checkmark$$

- (iv) The entropy is not time-independent: consider a spin–boson model (SBM) starting with the initial state $(|E_1\rangle + |E_2\rangle)/\sqrt{2} \otimes |0\rangle$, where $|0\rangle$ is at the ground state of the harmonic oscillator. The numerical simulation gives Figure 3.

The entropy of a closed system
is time-independent

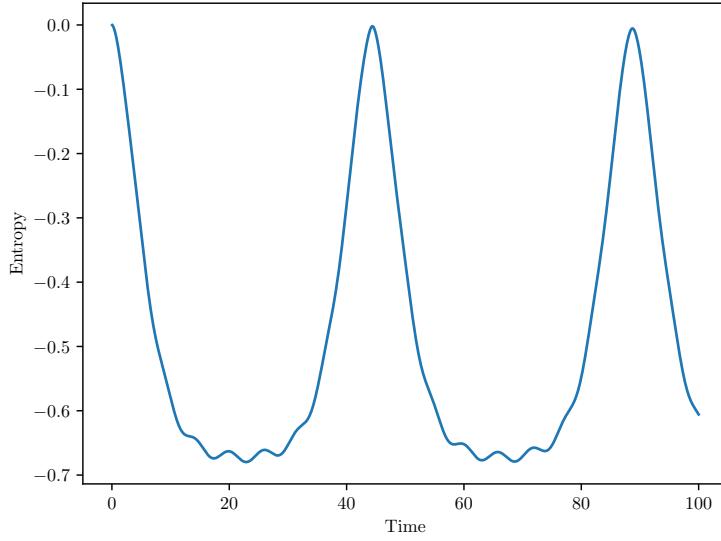


Figure 3: Time-dependency of entropy for spin–boson model.

5. (i) Recall the definition of $A_W(q, p) = \int_{-\infty}^{\infty} dy \langle q - y/2 | \hat{A}(q, p) | q + y/2 \rangle e^{ipy/\hbar}$. Therefore,

$$\begin{aligned} q_W &= \int_{-\infty}^{\infty} dy \langle q - y/2 | \hat{q} | q + y/2 \rangle e^{ipy/\hbar} \\ &= \int_{-\infty}^{\infty} dy \int_{-\infty}^{\infty} dx \delta(x - q + y/2)(q + y/2)\delta(x - q - y/2)e^{ipy/\hbar} \\ &= \int_{-\infty}^{\infty} dy (q + y/2)\delta(y)e^{ipy/\hbar} = q e^{ip \cdot 0/\hbar} = q, \end{aligned} \quad \checkmark$$

and similarly,

$$\begin{aligned} p_W &= \int_{-\infty}^{\infty} dy \langle q - y/2 | \hat{p} | q + y/2 \rangle e^{ipy/\hbar} \\ &= \int_{-\infty}^{\infty} dy \int_{-\infty}^{\infty} dx \delta(x - q + y/2)i\hbar \frac{\partial}{\partial x} \delta(x - q - y/2)e^{ipy/\hbar} \\ &= i\hbar \int_{-\infty}^{\infty} dy \int_{-\infty}^{\infty} dx \delta(x - q + y/2)\delta'(x - q - y/2)e^{ipy/\hbar} \\ &= i\hbar \int_{-\infty}^{\infty} dy \delta'(-y)e^{ipy/\hbar} = -i\hbar \int_{-\infty}^{\infty} dy \delta'(y)e^{ipy/\hbar} \\ &= -i\hbar \left(e^{ipy/\hbar} \right)' \Big|_{y=0} = p. \end{aligned} \quad \checkmark$$

Lemma 1. For any operators $\hat{A}(\hat{q}, \hat{p})$ and $\hat{B}(\hat{q}, \hat{p})$, $aA_W + bB_W = (aA + bB)_W$, where a, b are any complex numbers.

Proof. Note that

$$\begin{aligned}
A_W &= \int_{-\infty}^{\infty} dy \langle q - y/2 | \hat{A} | q + y/2 \rangle e^{ipy/\hbar}, \\
B_W &= \int_{-\infty}^{\infty} dy \langle q - y/2 | \hat{B} | q + y/2 \rangle e^{ipy/\hbar}, \\
(aA + bB)_W &= \int_{-\infty}^{\infty} dy \langle q - y/2 | aA + bB | q + y/2 \rangle e^{ipy/\hbar} \\
&= \int_{-\infty}^{\infty} dy \left(a \langle q - y/2 | \hat{A} | q + y/2 \rangle + b \langle q - y/2 | \hat{B} | q + y/2 \rangle \right) e^{ipy/\hbar} \\
&= a \int_{-\infty}^{\infty} dy \langle q - y/2 | \hat{A} | q + y/2 \rangle e^{ipy/\hbar} + b \int_{-\infty}^{\infty} dy \langle q - y/2 | \hat{B} | q + y/2 \rangle e^{ipy/\hbar} \\
&= aA_W + bB_W.
\end{aligned}$$
✓ ✓ ✓

□

Lemma 2. $(\hat{q}^n)_W = q^n$, $(\hat{p}^n)_W = p^n$.

Proof.

$$\begin{aligned}
(\hat{q}^n)_W &= \int_{-\infty}^{\infty} dy \langle q - y/2 | \hat{q}^n | q + y/2 \rangle e^{ipy/\hbar} \\
&= \int_{-\infty}^{\infty} dy \int_{-\infty}^{\infty} dx \delta(x - q + y/2) (q + y/2)^n \delta(x - q - y/2) e^{ipy/\hbar} \\
&= \int_{-\infty}^{\infty} dy (q + y/2)^n \delta(y) e^{ipy/\hbar} = q^n,
\end{aligned}$$
✓ ✓ ✓

and

$$\begin{aligned}
(\hat{p}^n)_W &= \int_{-\infty}^{\infty} dy \langle q - y/2 | \hat{p}^n | q + y/2 \rangle e^{ipy/\hbar} \\
&= \int_{-\infty}^{\infty} dy \int_{-\infty}^{\infty} dx \delta(x - q + y/2) (i\hbar)^n \frac{\partial^n}{\partial x^n} \delta(x - q - y/2) e^{ipy/\hbar} \\
&= (i\hbar)^n \int_{-\infty}^{\infty} dy \int_{-\infty}^{\infty} dx \delta(x - q + y/2) \delta^{(n)}(x - q - y/2) e^{ipy/\hbar} \\
&= (i\hbar)^n \int_{-\infty}^{\infty} dy \delta^{(n)}(-y) e^{ipy/\hbar} = (-i\hbar)^n \int_{-\infty}^{\infty} dy \delta^{(n)}(y) e^{ipy/\hbar} \\
&= (-i\hbar)^n \left. \left(e^{ipy/\hbar} \right)^{(n)} \right|_{y=0} = p^n.
\end{aligned}$$
✓ ✓ ✓ ✓ ✓ ✓ ✓

□

Therefore, if \hat{T} and \hat{W} converge to their Taylor series, then

$$\begin{aligned}
H_W &= (T(p) + V(q))_W = \left(\sum_{n=0}^{\infty} \frac{1}{n!} \frac{d^n T}{dp^n} p^n + \sum_{n=0}^{\infty} \frac{1}{n!} \frac{d^n V}{dq^n} q^n \right)_W \\
&= \sum_{n=0}^{\infty} \frac{1}{n!} \frac{d^n T}{dp^n} p_W^n + \sum_{n=0}^{\infty} \frac{1}{n!} \frac{d^n V}{dq^n} q_W^n = \sum_{n=0}^{\infty} \frac{1}{n!} \frac{d^n T}{dp^n} p^n + \sum_{n=0}^{\infty} \frac{1}{n!} \frac{d^n V}{dq^n} q^n = T(p) + V(q) = H(q, p).
\end{aligned}$$
✓

good!

(ii) For a pure state, $\hat{\rho} = |\psi\rangle\langle\psi|$, and

$$\begin{aligned}\rho_W &= \frac{1}{2\pi\hbar} \int_{-\infty}^{\infty} dy \langle q - y/2 | \hat{\rho} | q + y/2 \rangle e^{ipy/\hbar} \\ &= \frac{1}{2\pi\hbar} \int_{-\infty}^{\infty} dy \psi(q + y/2)^* \psi(q - y/2) e^{ipy/\hbar},\end{aligned}$$

Then


$$\begin{aligned}\lim_{\hbar \rightarrow 0} \rho_W &= \frac{1}{2\pi\hbar} \int_{-\infty}^{\infty} dy \langle q - y/2 | \hat{\rho} | q + y/2 \rangle e^{ipy/\hbar} \\ &= \frac{1}{2\pi} \int_{-\infty}^{\infty} dy \psi(q + y/2)^* \psi(q - y/2) \lim_{\hbar \rightarrow 0} e^{ipy/\hbar} / \hbar,\end{aligned}$$

You should also need to consider the dependence of the wavefunction on \hbar

and note that $e^{ipy/\hbar}$ oscillates when $\hbar \rightarrow +0$. Therefore, ρ_W is singular for a pure quantum state.

(iii) We need the following lemma.

Lemma 3. For any operator $\hat{A}(\hat{q}, \hat{p})$, $(A^\dagger)_W = (A_W)^*$.

Proof. Note that

$$\begin{aligned}(A^\dagger)_W &= \int_{-\infty}^{\infty} dy \langle q - y/2 | \hat{A}^\dagger | q + y/2 \rangle e^{ipy/\hbar} \\ &= \int_{-\infty}^{\infty} dy \langle q + y/2 | \hat{A} | q - y/2 \rangle^* e^{ipy/\hbar} \\ &= \int_{-\infty}^{\infty} dy \langle q - y/2 | \hat{A} | q + y/2 \rangle^* e^{-ipy/\hbar} \\ &= (A_W)^*\end{aligned}$$

□

$$\begin{aligned}\frac{d(\hat{O}_H)_W}{dt} &= \left(-\frac{i}{\hbar} [\hat{O}_H, \hat{H}_H] \right)_W \\ &= -\frac{i}{\hbar} ((\hat{O}_H \hat{H}_H)_W - (\hat{H}_H \hat{O}_H)_W) \\ &= -\frac{i}{\hbar} ((\hat{O}_H \hat{H}_H)_W - ((\hat{O}_H \hat{H}_H)_W)^*).\end{aligned}$$

Recall that $(\hat{A}\hat{B})_W = A_W e^{-\frac{i\hbar}{2}\vec{\Lambda}} B_W$, where $\vec{\Lambda} = \overleftarrow{\frac{\partial}{\partial p}} \overleftarrow{\frac{\partial}{\partial q}} - \overrightarrow{\frac{\partial}{\partial q}} \overrightarrow{\frac{\partial}{\partial p}}$, therefore,

$$\begin{aligned}\frac{d(\hat{O}_H)_W}{dt} &= -\frac{i}{\hbar} ((\hat{O}_H)_W e^{-\frac{i\hbar}{2}\vec{\Lambda}} (\hat{H}_H)_W - ((\hat{O}_H)_W e^{-\frac{i\hbar}{2}\vec{\Lambda}} (\hat{H}_H)_W)^*) \\ &= -\frac{i}{\hbar} ((\hat{O}_H)_W e^{-\frac{i\hbar}{2}\vec{\Lambda}} (\hat{H}_H)_W - (\hat{O}_H)_W^* e^{\frac{i\hbar}{2}\vec{\Lambda}} (\hat{H}_H)_W^*) \\ &= -\frac{2}{\hbar} (\hat{O}_H)_W \sin \frac{\hbar \vec{\Lambda}}{2} (\hat{H}_H)_W.\end{aligned}$$

In the last step we have applied the fact that $(\hat{O}_H)_W$ is real for any Hermitian \hat{O}_H since $(\hat{O}_H)_W = (\hat{O}_H^\dagger)_W = ((\hat{O}_H)_W)^*$.

(iv) When $\hbar \rightarrow 0$, $(\hat{O}_H)_W = O^c + \hbar O_1 + o(\hbar)$, $(\hat{H}_H)_W = H^c + \hbar H_1 + o(\hbar)$, and $\sin \frac{\hbar \vec{\Lambda}}{2} = \frac{\hbar \vec{\Lambda}}{2} + o(\hbar)$, then,

$$\begin{aligned}\frac{d(\hat{O}_H)_W}{dt} &= -\frac{2}{\hbar} (\hat{O}_H)_W \sin \frac{\hbar \vec{\Lambda}}{2} (\hat{H}_H)_W \\ \Rightarrow \frac{dO^c}{dt} + o(1) &= -2O^c \frac{\hbar \vec{\Lambda}}{2} H^c + o(1) \\ \Rightarrow \frac{dO^c}{dt} &= \{H^c, O^c\}_P + o(1),\end{aligned}$$

where $\{\cdot, \cdot\}_P$ is the Poisson bracket.

(v) When we use the wigner representation and take the limit of $\hbar \rightarrow 0$ then we recover the classical mechanics, which means that the classical mechanics is a specific representation of quantum mechanics under the limit that Planck constant is small enough.

6. Code is available at <https://github.com/vINyLogY/QD-hw2>.

(A) (i) Here we only need to focus on monitoring the eigenstate with $n = 100$. Therefore the dimensionality of the grid basis should be at least 101 (assume we start counting n from 0). We choose an overkill (checked convergence by simply increasing these parameters) dimensionality 1001 and $q_0 = 25$, and the results are showed in Figure 4.

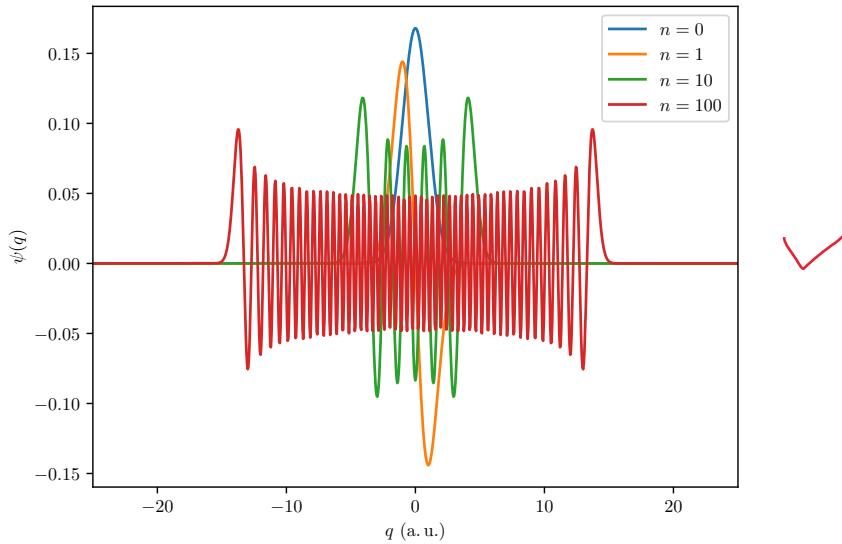


Figure 4: $\psi(x)$ for a harmonic oscillator.

The following results are focused with $q_0 = 25$.

(ii) See Figure 5.

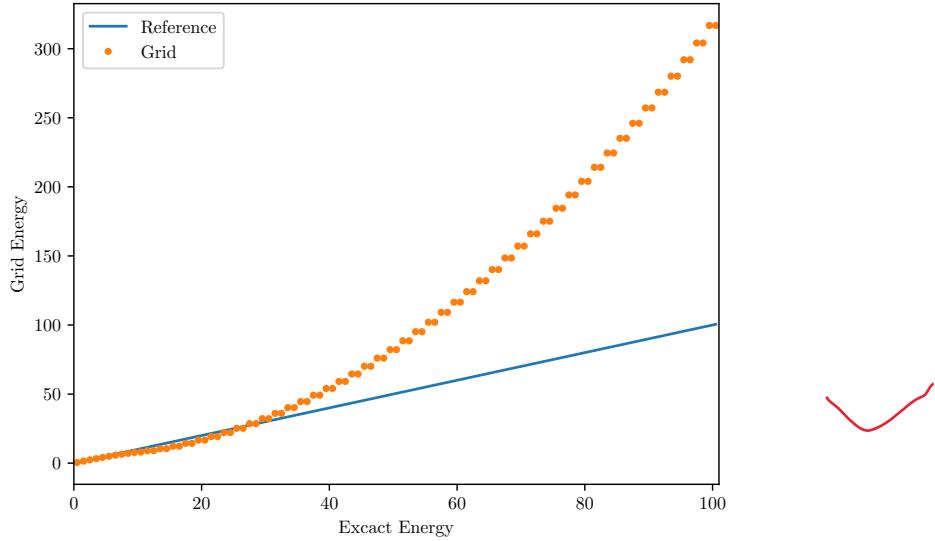


Figure 5: Energy of each eigenstates for a harmonic oscillator.

Note that the grid method only gets the exact answer for those states with small eigenvalues. This is because for the excited states with large energy, the wavefunction oscillates crazy as showed in Figure 4, and hence we need more dense grid to characterize the wavefunction.

- (iii) We need to change the dimensionality of grid basis in this case. For the eigenstate $n = 32$, we need at least 33 grids. We calculated the energy difference between the one from the grid representation and the exact one, change the dimensionality of grid basis, *i.e.*, different Δq . See Figure 6.

You are asked to plot the error in the determination of the energy of the

state versus q and determine the q necessary to accurately capture the eigenenergy. this q would be more obvious if you
plot $\log(\text{error})$ vs Δq

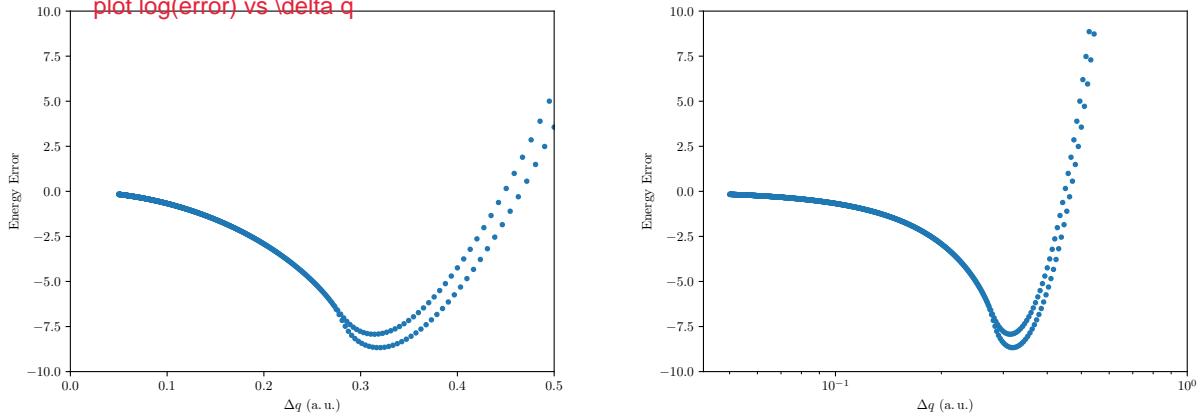


Figure 6: Energy error vs. Δq .

Therefore, $\Delta q < 0.1$ a. u. is an adequate choice.

- (B) (i) Use Sine-DVR¹ here. As the concerns in grid basis, we choose the dimensionality to be 999 and $q_0 = 25$. See Figure 7.

¹See <https://www.pci.uni-heidelberg.de/tc/usr/mctdh/lit/NumericalMethods.pdf>.

have you checked your result with smaller q_0 ?

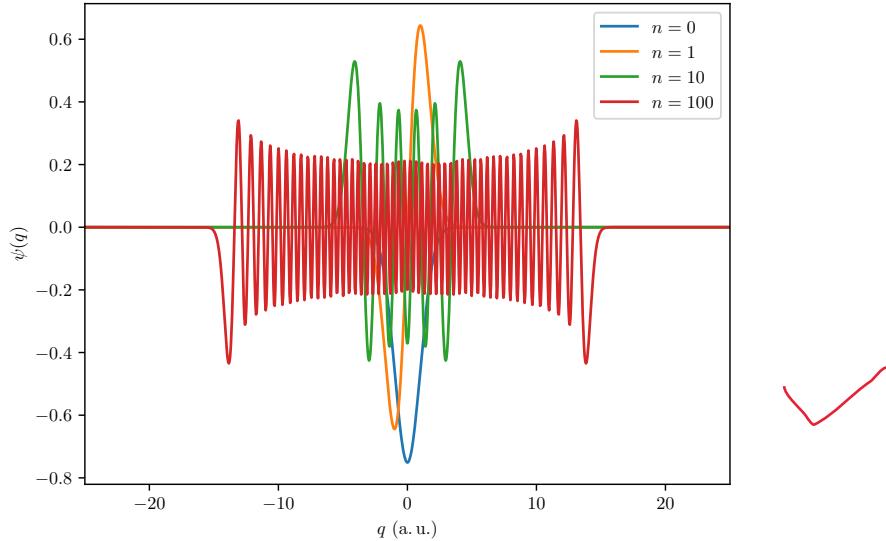


Figure 7: Energy of each eigenstates for a harmonic oscillator.

Note that the $\psi(q)$ has up to a different phase from the results by grid basis.

(ii) See Figure 8.

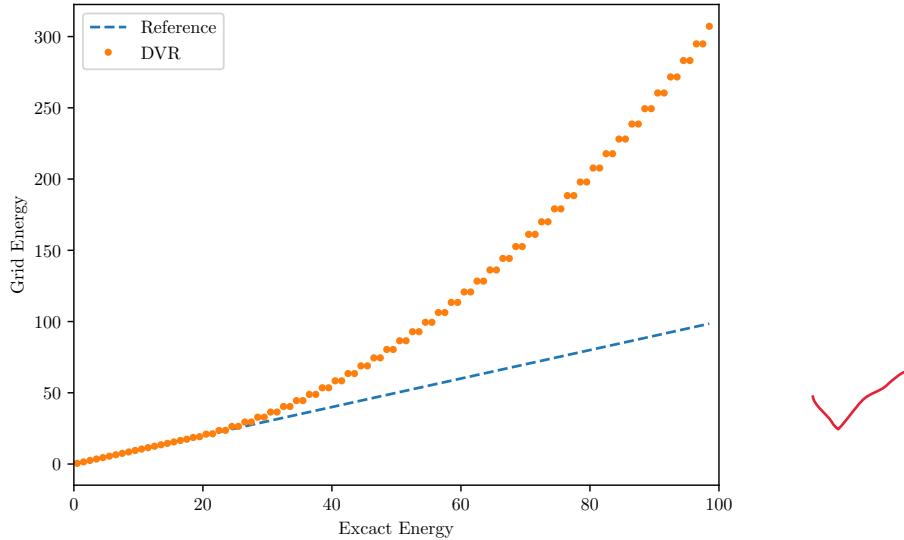


Figure 8: Energy of each eigenstates for a harmonic oscillator.

Similarly, the DVR method only gets the exact answer for those states with small eigenvalues. This is because for the excited states with large energy, they need the basis includes the functions that corresponding to larger momentum, therefore, we need more dimensionality for the finite basis representation, or smaller grid interval, to represent such high-energy states.

- (iii) We need to change the dimensionality of grid basis in this case. For the eigenstate $n = 32$, we need at least 33 grids. We calculated the energy difference between the one from the grid representation and the exact one, change the dimensionality of grid basis, *i.e.*, different Δq . See Figure 9.

You are asked to plot the error in the determination of the energy of the state versus q and determine the q necessary to accurately capture the eigenenergy. This q would be more obvious if you plot $\log(\text{error})$ vs Δq .

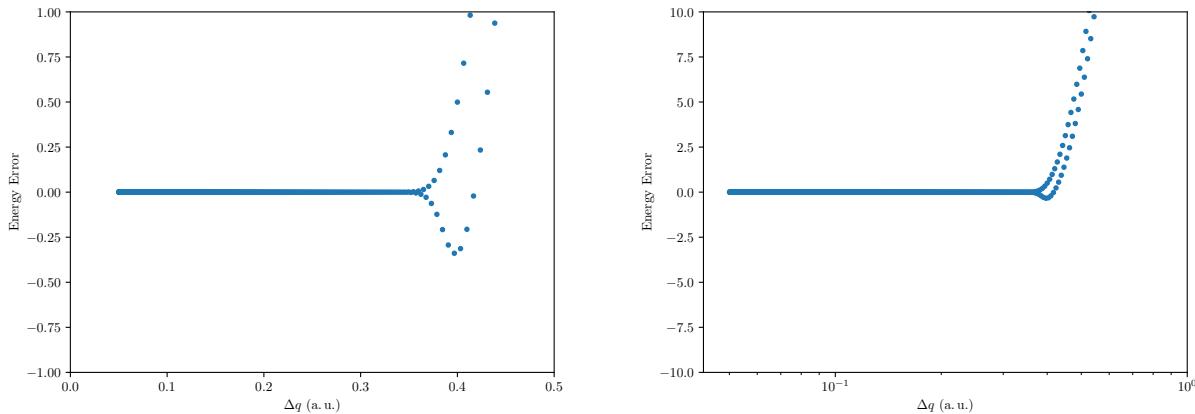


Figure 9: Energy error vs. Δq .

Therefore, $\Delta q < 0.1$ a.u. is an adequate choice.

- (iv) From Figure 5 and Figure 8 we can find that DVR can get exact energy for states with relatively larger energy when compared to grid method. ~~Also, from Figure 6 and Figure 9 we can find that DVR can get exact energy with a relatively small dimensionality, or larger grid interval.~~
- (C) (i) Suppose under 1-D DVR basis $\{|\chi_i\rangle\}$ the kinetic matrix in one dimension i is T_i , the potential matrix in one dimension i is V_i , then the kinetic matrix of 2-D is $T_1 \otimes T_2$, and the potential matrix is a diagonal matrix $V_1 \otimes V_2$. Hence we can diagonalize the hamiltonian in the 2-D basis.
- (ii) See Figure 10. It is clear that for such a system, the exact eigenenergies are $E = \hbar\omega(n + 1)$, where $n = n_a + n_b$, and $n_a, n_b \in \mathbb{N}$. This means that for $E = \hbar\omega$, the degeneracy is 1 ($n = 0 + 0$); for $E = 2\hbar\omega$, the degeneracy is 2 ($n = 1 + 0 = 0 + 1$); for $E = 3\hbar\omega$, the degeneracy is 3 ($n = 2 + 0 = 0 + 2 = 1 + 1$); for $E = 4\hbar\omega$, the degeneracy is 4 ($n = 3 + 0 = 0 + 3 = 1 + 2 = 2 + 1$); and so on. From Figure 10 we can also find that the parameters (upper and lower bound are both 25 and -25 for each dimension; the dimensionality of basis for each degrees of freedom is 99) is sufficient for the lowest 10 eigenstates.
- (iii) See Figure 11.

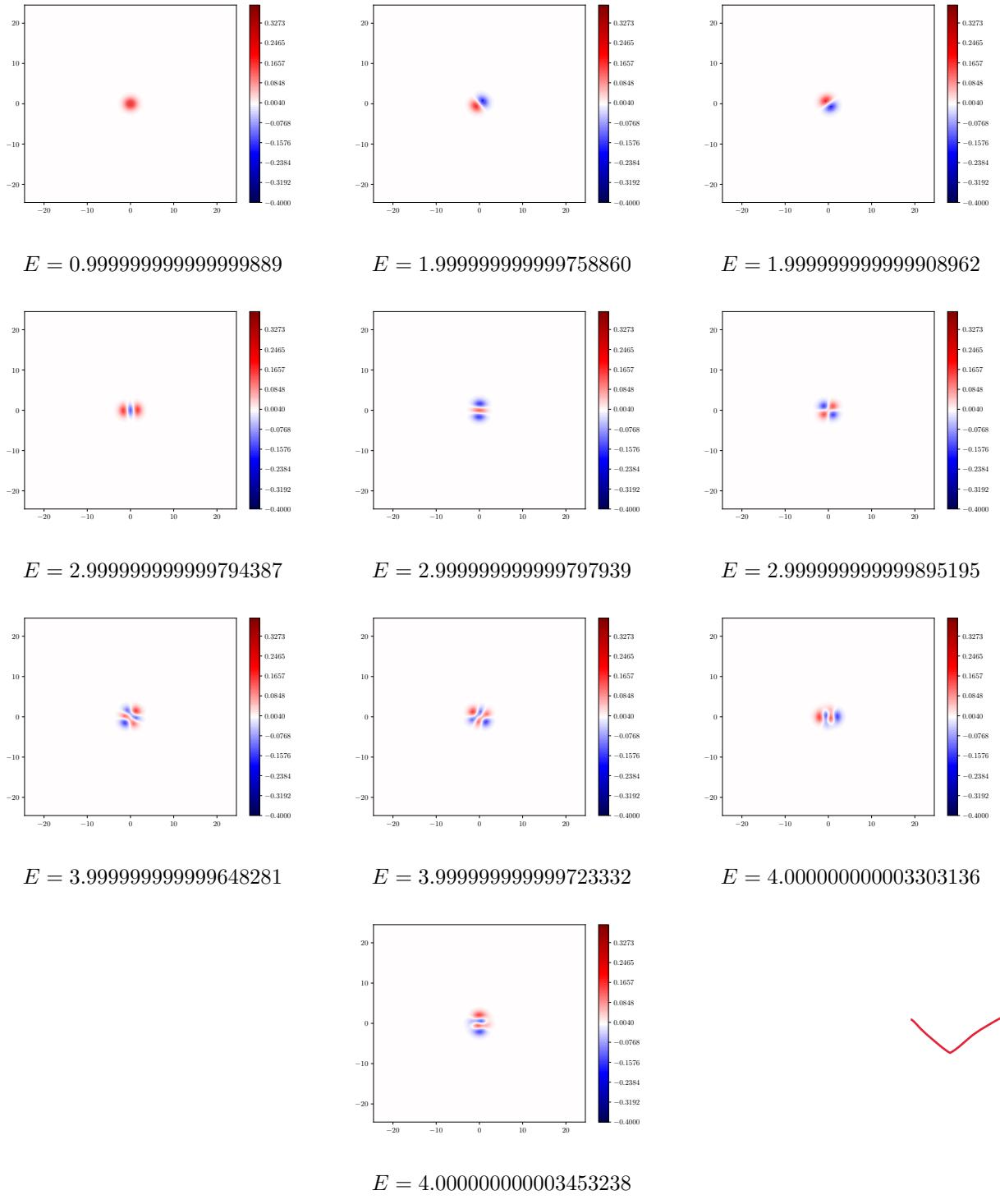


Figure 10: Eigenstates and their eigenenergies for a 2-D harmonic oscillator.

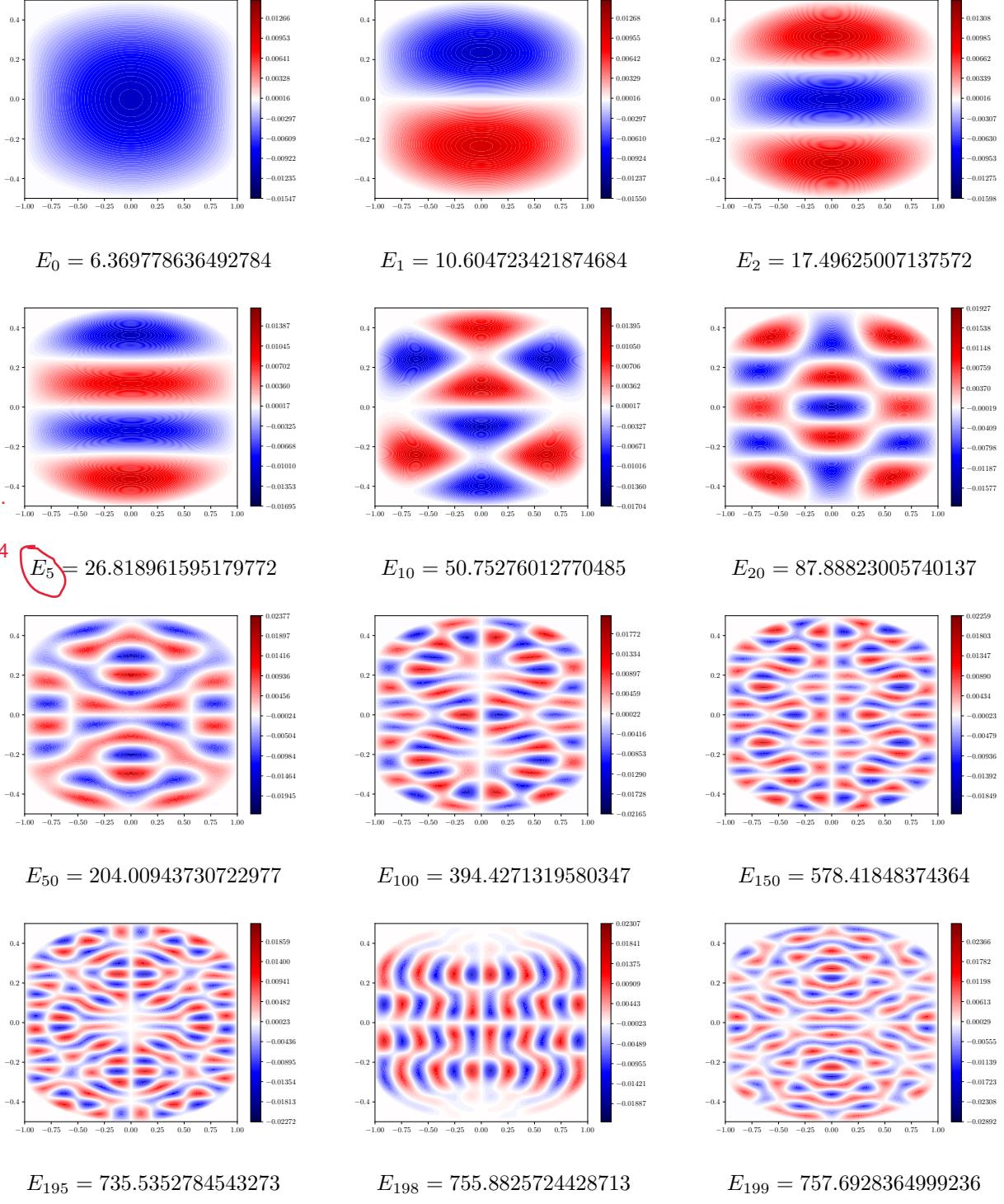


Figure 11: Eigenstates and their eigenenergies for a 2-D potential well.