

9.9.

b) Now do the same for all states but using the operators a and a^\dagger .

$$a = \frac{1}{\sqrt{2\hbar m\omega}}(-ip + m\omega x)$$

$$a^\dagger = \frac{1}{\sqrt{2\hbar m\omega}}(ip + m\omega x)$$

$$x = \sqrt{\frac{\hbar}{2m\omega}}(a^\dagger + a)$$

$$p = i\sqrt{\frac{\hbar m\omega}{2}}(a^\dagger - a)$$

$$\begin{aligned}\langle x \rangle &= \langle n|x|n \rangle = \sqrt{\frac{\hbar}{2m\omega}} \langle n|a^\dagger + a|n \rangle \\ &= \sqrt{\frac{\hbar}{2m\omega}} [\langle n|a^\dagger|n \rangle + \langle n|a|n \rangle] = \sqrt{\frac{\hbar}{2m\omega}} [\langle n|\sqrt{n+1}|n+1 \rangle + \langle n|\sqrt{n}|n-1 \rangle] \\ &= \sqrt{\frac{\hbar}{2m\omega}} [\sqrt{n+1}\langle n|n+1 \rangle + \sqrt{n}\langle n|n-1 \rangle] = 0 \quad \text{since } \langle n|m \rangle = \delta_{nm}\end{aligned}$$

$$\begin{aligned}\langle p \rangle &= \langle n|p|n \rangle = \sqrt{\frac{\hbar}{2m\omega}} \langle n|a^\dagger - a|n \rangle \\ &= \sqrt{\frac{\hbar}{2m\omega}} [\langle n|a^\dagger|n \rangle - \langle n|a|n \rangle] = \sqrt{\frac{\hbar}{2m\omega}} [\langle n|\sqrt{n+1}|n+1 \rangle - \langle n|\sqrt{n}|n-1 \rangle] \\ &= \sqrt{\frac{\hbar}{2m\omega}} [\sqrt{n+1}\langle n|n+1 \rangle - \sqrt{n}\langle n|n-1 \rangle] = 0 \quad \text{since } \langle n|m \rangle = \delta_{nm}\end{aligned}$$

Note also that $\langle n|a^2|n \rangle = 0$ and $\langle n|(a^\dagger)^2|n \rangle = 0$ in a similar manner, so that

$$\begin{aligned}\langle x^2 \rangle &= \langle n|x^2|n \rangle = \frac{\hbar}{2m\omega} \langle n|(a^\dagger + a)^2|n \rangle = \frac{\hbar}{2m\omega} \langle n|(a^\dagger)^2 + a^\dagger a + aa^\dagger + a^2|n \rangle \\ &= \frac{\hbar}{2m\omega} \langle n|a^\dagger a + aa^\dagger|n \rangle = \frac{\hbar}{2m\omega} \langle n|\sqrt{n}\sqrt{n} + \sqrt{n+1}\sqrt{n+1}|n \rangle \\ &= \frac{\hbar}{2m\omega} (2n+1) = \frac{\hbar}{m\omega} (n + \frac{1}{2}) \\ \langle p^2 \rangle &= \langle n|p^2|n \rangle = -\frac{\hbar m\omega}{2} \langle n|(a^\dagger - a)^2|n \rangle = -\frac{\hbar m\omega}{2} \langle n|(a^\dagger)^2 - a^\dagger a - aa^\dagger + a^2|n \rangle \\ &= \frac{\hbar m\omega}{2} \langle n|a^\dagger a + aa^\dagger|n \rangle = \frac{\hbar m\omega}{2} \langle n|\sqrt{n}\sqrt{n} + \sqrt{n+1}\sqrt{n+1}|n \rangle \\ &= \frac{\hbar m\omega}{2} (2n+1) = \hbar m\omega (n + \frac{1}{2})\end{aligned}$$

9.11 a) Normalize:

$$\begin{aligned} |\psi(t=0)\rangle &= A[|0\rangle + 2e^{i\pi/2}|1\rangle] \\ 1 = \langle\psi|\psi\rangle &= A^* \left(\langle 0| + 2e^{-i\pi/2} \langle 1| \right) A \left(|0\rangle + 2e^{i\pi/2} |1\rangle \right) \\ &= |A|^2 (1 + 4) = |A|^2 5 \\ \Rightarrow A &= \frac{1}{\sqrt{5}} \end{aligned}$$

b) Time evolution

$$\begin{aligned} |\psi(t)\rangle &= \frac{1}{\sqrt{5}} \left(e^{-iE_0t/\hbar} |0\rangle + 2e^{i\pi/2} e^{-iE_1t/\hbar} |1\rangle \right) \\ &= e^{-i\omega t/2} \frac{1}{\sqrt{5}} \left(|0\rangle + 2e^{i\pi/2} e^{-i\omega t} |1\rangle \right) \end{aligned}$$

c) Expectations values with ladder ops:

$$\begin{aligned} \langle x \rangle &= \langle \psi(t) | x | \psi(t) \rangle = \sqrt{\frac{\hbar}{2m\omega}} \langle \psi(t) | a^\dagger + a | \psi(t) \rangle \\ &= \sqrt{\frac{\hbar}{2m\omega}} e^{+i\omega t/2} \frac{1}{\sqrt{5}} \left(\langle 0| + 2e^{-i\pi/2} e^{+i\omega t} \langle 1| \right) (a^\dagger + a) e^{-i\omega t/2} \frac{1}{\sqrt{5}} \left(|0\rangle + 2e^{i\pi/2} e^{-i\omega t} |1\rangle \right) \\ &= \sqrt{\frac{\hbar}{2m\omega}} \frac{1}{5} \left[2e^{i\pi/2} e^{-i\omega t} \langle 0|a|1\rangle + 2e^{-i\pi/2} e^{+i\omega t} \langle 1|a^\dagger|0\rangle \right] \\ &= \sqrt{\frac{\hbar}{2m\omega}} \frac{2}{5} \left[ie^{-i\omega t} \sqrt{1} - ie^{+i\omega t} \sqrt{1} \right] = \sqrt{\frac{\hbar}{2m\omega}} \frac{4}{5} \sin \omega t \end{aligned}$$

Momentum expectation value:

$$\begin{aligned} \langle p \rangle &= \langle \psi(t) | p | \psi(t) \rangle = i\sqrt{\frac{m\omega\hbar}{2}} \langle \psi(t) | a^\dagger - a | \psi(t) \rangle \\ &= i\sqrt{\frac{m\omega\hbar}{2}} e^{+i\omega t/2} \frac{1}{\sqrt{5}} \left(\langle 0| + 2e^{-i\pi/2} e^{+i\omega t} \langle 1| \right) (a^\dagger - a) e^{-i\omega t/2} \frac{1}{\sqrt{5}} \left(|0\rangle + 2e^{i\pi/2} e^{-i\omega t} |1\rangle \right) \\ &= i\sqrt{\frac{m\omega\hbar}{2}} \frac{1}{5} \left[-2e^{i\pi/2} e^{-i\omega t} \langle 0|a|1\rangle + 2e^{-i\pi/2} e^{+i\omega t} \langle 1|a^\dagger|0\rangle \right] \\ &= i\sqrt{\frac{m\omega\hbar}{2}} \frac{2}{5} \left[-ie^{-i\omega t} \sqrt{1} - ie^{+i\omega t} \sqrt{1} \right] = \sqrt{\frac{m\omega\hbar}{2}} \frac{4}{5} \cos \omega t \end{aligned}$$

Ehrenfest's theorem in this case is

$$\langle p \rangle = m \frac{d}{dt} \langle x \rangle = m \frac{d}{dt} \left[\sqrt{\frac{\hbar}{2m\omega}} \frac{4}{5} \sin \omega t \right] = m \sqrt{\frac{\hbar}{2m\omega}} \frac{4}{5} [\omega \cos \omega t] = \sqrt{\frac{m\omega\hbar}{2}} \frac{4}{5} \cos \omega t$$

So it is satisfied.

%*****

% Program 9.20

% Feel free to play with this code

%*****

clear

% Calculate and plot orthogonal polynomials

% PARAMETERS:

Order=25; %number of energy eigenfunctions used (as order

```

%increases the approximation improves!)
Xoffset=5;      %amount of offset
dx=0.05;        %spacing of interval
L=8;            %size of region we are interested in
X=-L:dx:L;      %support on x axis - fills X vector from -L to L
Period = 5;     %for later on in code - how many period to look at
TimeInterval = 0.1; %Steps for animation of plots later...
number_points=size(X,2); %how many points used on x-axis? "size" fcn

Phi_x_0_offset=(pi^(-0.25))*(exp(-((X-Xoffset).^2)/2)); % Build offset gnd state

for n=0:Order;

Phi_x_n(n+1,:)=(pi^(-0.25))*(((2.^n)*factorial(n))^(-0.5))*hermite(n,X).*(exp(-
((X.^2)/2)));

C(n+1)=dx*Phi_x_n(n+1,:)*(Phi_x_0_offset)'; %Even though strange looking, this is
%the integral that finds the c_n
%coefficients of the overlap
%integral. (It is an inner product!) Result is a
%row vector with c_n values.
Check_C_n(n+1)=(Xoffset^n/(2^n*factorial(n))^(0.5))*exp(-(Xoffset^2/4));

end
Check=C
Check2=Check_C_n
Sumcheck=(C*C');
%Checks the squared sums of the Cn's to see how close they come to
%saturating the normalization condition. Should be close to 1 if we are
%good.
ApproximationFunction=Phi_x_n'*C';
%This is simple, elegant and correct. The Matrix Phi_x_n' operates on the
%column vector C' via matrix multiplication to produce our approximation
%function.

%ApproximationFunction=sum((C'*ones(1,number_points)).*Phi_x_n);

%The above is a very short bit of code that multiplies the c_n coefficients by
%the eigenfunctions to produce our approximation of the wavefunction. It does the
same thing
%as the for loop below. Working inside out, it does it by producing column matrix of
%coefficients, and taking outer product of it with matrix of ones of the same size as
%support X. This produces a matrix of n rows by size of X columns matrix. Element
by
%element muliplication (.* ) with the the Phi_x_n produces a matrix of legendre_lth

```

```
%order*c_n. Sum just sums up columns to produce full approximation function,
NewF2.
```

```
%Code below does the same thing as the single line above, but in more pedestrian
way
```

```
%ApproximationFunction=0*X;
%for i=1:7;
%ApproximationFunction = ApproximationFunction + Funct2(i,:)*C(i);
%end
```

```
%Calculation of the Expectation Value of the Energy
```

```
EnergySpectrum=0.5:1:(Order+0.5);
%Creates an energy spectrum vector with energies in units of  $\hbar\omega$ 
```

```
Expected_Energy=(C.^2)*EnergySpectrum'
%Uses the energy spectrum vector to calculate the expectation value of the
%energy -->  $\text{Sum}((c_n^2)*E_N)$ 
```

```
%Calculation of the time dependent state
for T=0:TimeInterval:2*pi*Period
    %Time is T. Period is set up above and is how many period
```

```
    Factor=C.*(exp(-1i*EnergySpectrum*T));
    %Factor is  $c_n \exp(-i\omega t) = c_n \exp(-iE_n t / \hbar)$ 
```

```
    Approx_Phi_x_t=Phi_x_n'*Factor';
    %This is the time dependent approximation function at time = T
    ProbDensity=Approx_Phi_x_t.*conj(Approx_Phi_x_t);
    %Now plot the functions and the envelope
    figure(1)
    %plot(X,ProbDensity,'b',X,0.5*X.*X,'-r');
```

```
plot(X,ProbDensity,X,real(Approx_Phi_x_t),'r',X,imag(Approx_Phi_x_t),'g',X,0.5*X.*X,'-r');
```

```
axis([-Xoffset-2,Xoffset+2,-1,1]);
%slows things down so we can see the animation
pause(0.1);
%
```

```
end
```

```
%Propogator_with_Time=Propogator*exp(t)
```

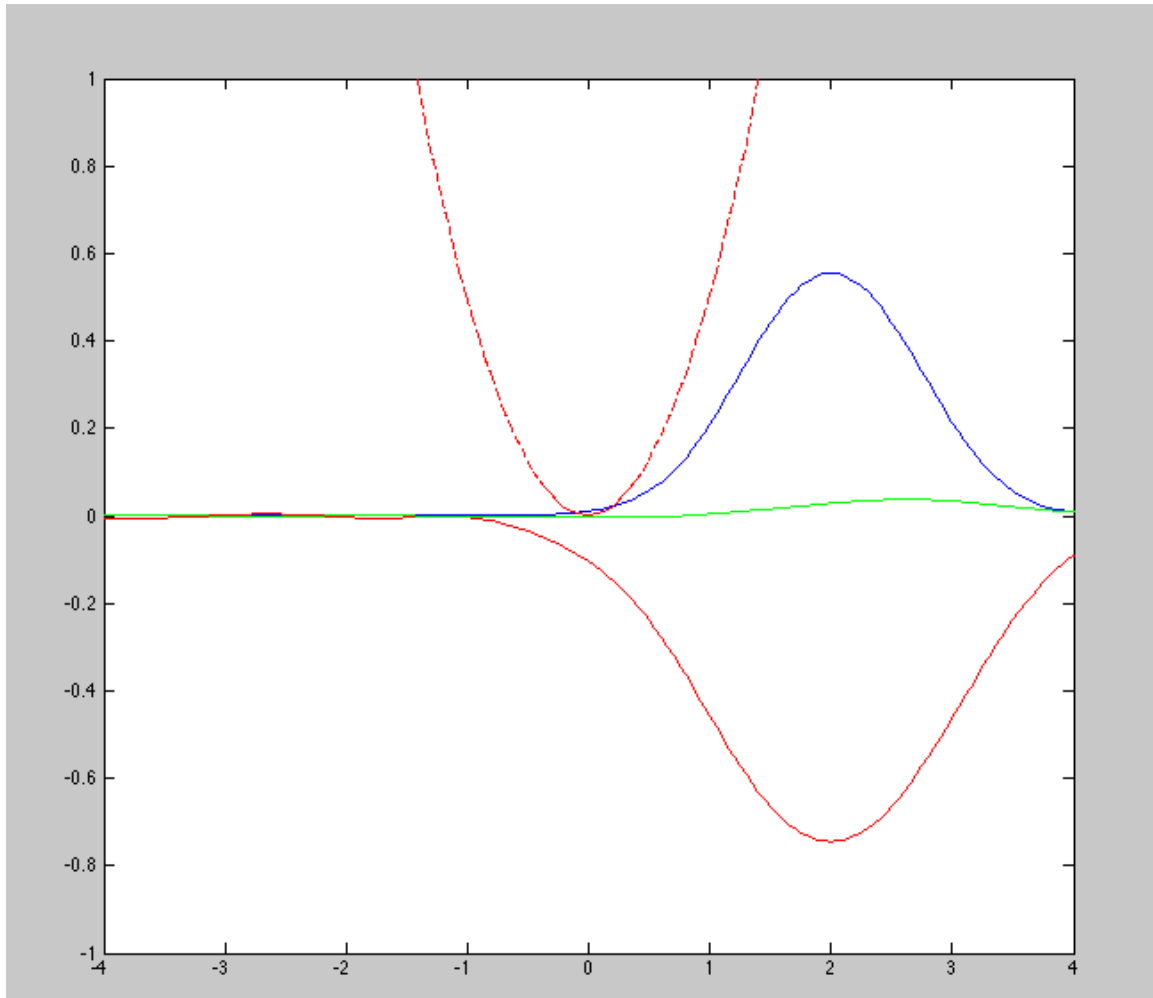
```
%TimeDepApproxFcn=sum((C*ones(1,number_points)).*Phi_x_n)
```

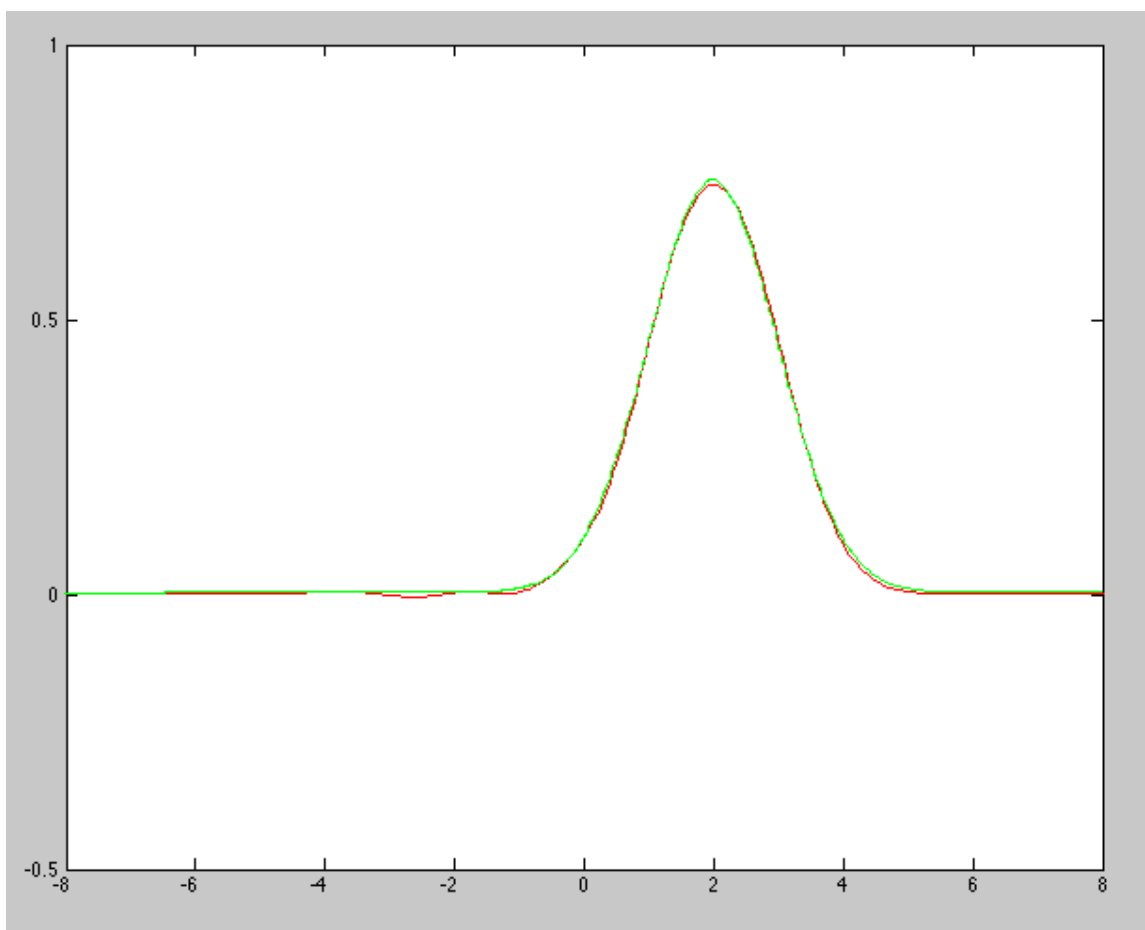
```
figure(2)
```

```
plot(X,ApproximationFunction,'r',X,Phi_x_0_offset,'--g');
```

```
axis([min(X),max(X),-0.5,1]);
```

```
shg
```





```
>> Plot9_20b
```

```
Check =
```

```
0.3679  0.5203  0.5203  0.4248  0.3004  0.1900  0.1097  0.0586  0.0293
```

```
Check2 =
```

```
0.3679  0.5203  0.5203  0.4248  0.3004  0.1900  0.1097  0.0586  0.0293
```

```
Expected_Energy =
```

```
2.4977
```