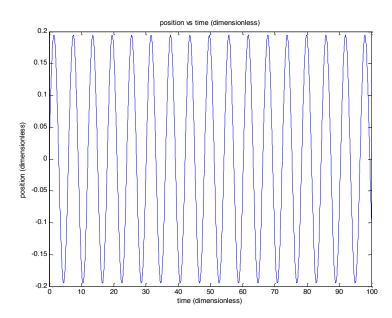
Below: all time steps are of 5.00000E-002.

Below:

Using gamma = 3.0;

initial\_v = 0.2;

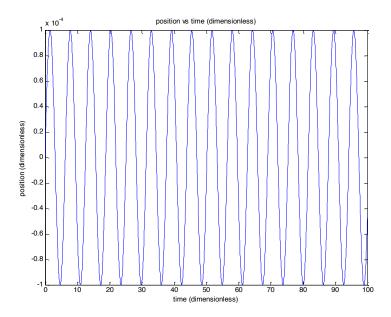


Below:

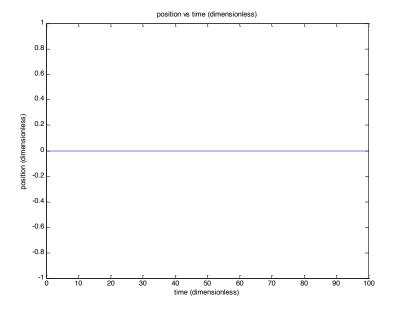
## Gamma = 3.0

initial\_x = 0.0;

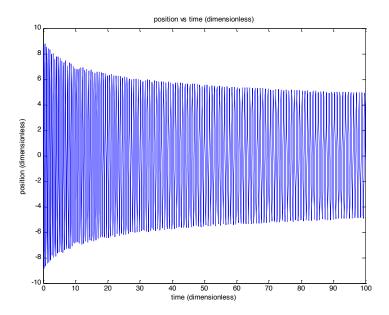
initial\_v = 0.0001;



Below v = 0;



V= 100



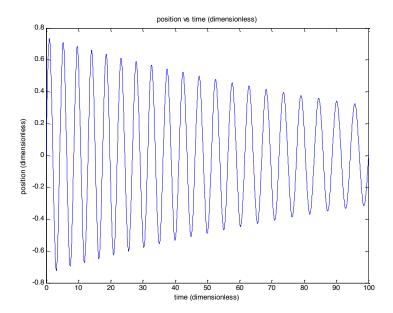
Frequency increases proportionally to the increase in initial velocity with initial position fixed at 0 and constant gamma.

## Q1b)

```
double period_initial = 10.0;
final_t = period_initial*10.0;
h = 0.05;
number_of_steps = final_t/h;
initial_x = 0.0;
initial_v = 1.0000;
double damp =0.02;
```

$$dydt[0] = y[1];$$
  
 $dydt[1] = -y[0]-3.0*pow(y[0],3)-damp*y[1];$ 

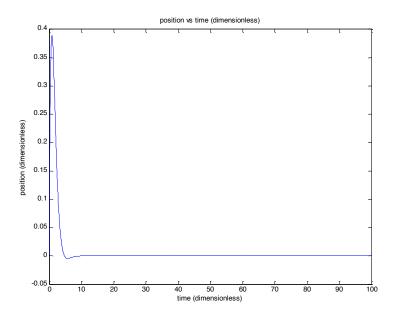
#### below is underdamped



Below is underdamped: double damp =1.7;

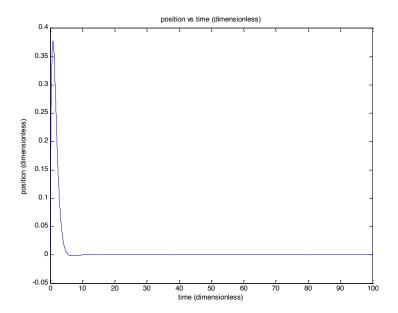
$$dydt[0] = y[1];$$

dydt[1] = -y[0]-3.0\*pow(y[0],3)-damp\*y[1];



## Below is overdamped:

```
initial_x = 0.0;
initial_v = 1.0000;
double damp =1.8;
  dydt[0] = y[1];
  dydt[1] = -y[0]-3.0*pow(y[0],3)-damp*y[1];
```

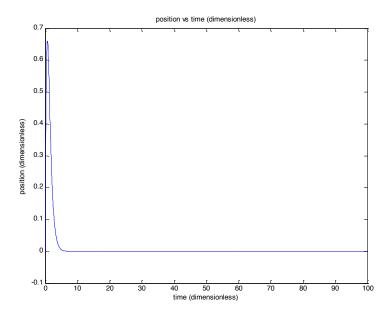


Below; initial v is increased from 1 to 2:

double damp =2.0;

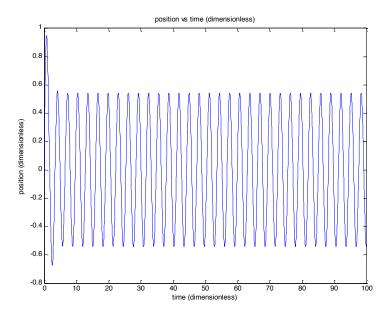
dydt[0] = y[1];

dydt[1] = -y[0]-3.0\*pow(y[0],3)-damp\*y[1];



The value that results in critical damping increases from 1.8 to 2.0 when changing initial v from 1.0 to 2.0.

```
Q1c)  
Below:  
double damp =2.0;  
double f = 2.5;  
double omega = 2.0;  
dydt[0] = y[1];  
dydt[1] = -y[0]-3.0*pow(y[0],3)-damp*y[1]+f*cos(omega*t);
```



# Program code for part c below:

//Code modified by Vincent Chiu. Based on code by professor Kristin Schleich.

//Pendulum with NO Driving force, damping of 10.0.

//duffing oscillator final project

#include <stdio.h>

#include <math.h>

#include <stdlib.h>

#include <assert.h>

```
void initialise (double *, double *, double *, int *);
void derivatives(double, double *, double *);
void euler(double *, double *, int, double, double, double *);
void output( FILE *, double, double, double *, double);
void runge_kutta_2(double *, double *, int, double, double, double *, void (*)(double, double*,
double*));
void runge_kutta_4(double *, double *, int, double , double *, void (*)(double, double *, double
*));
int main()
{
  printf("Welcome to Chiu Industries. Pendulum Calculator. NO Driving Force Mode. \n");
  FILE *output_file;
  //declarations of variables
  int number_of_steps;
  double initial_x, initial_v, final_t, E_initial;
  double h;
  double y[2], dydt[2], yout[2];
  double t;
  output_file = fopen("finalq1c.dat", "w");
  //read in input data from screen
  //initialise (&initial_x, &initial_v, &final_t, &number_of_steps);
```

```
// double length_rod = 1.0;
// double gravitational_constant = 1.0;
// double angular_frequency_initial = sqrt(gravitational_constant/length_rod);
//double period_initial = 2.0*M_PI/angular_frequency_initial;
double period_initial = 10.0;
final_t = period_initial*10.0;
h = 0.05;
number_of_steps = final_t/h;
initial_x = 0.0;
initial_v = 2.0000;
//number_of_steps = 100000;
assert(number_of_steps >0);
//initialise position and velocity
y[0] = initial_x;
y[1] = initial_v;
```

```
E_{initial} = 0.5*y[0]*y[0] + 0.5*y[1]*y[1];
  //h=final_t/number_of_steps;
  output(output_file, h, t, y, E_initial);
  t=0.0;
  while(t<=final_t)
  {
    derivatives(t,y,dydt);
    //euler(y, dydt, 2, t, h, yout);
    runge_kutta_4(y, dydt, 2, t, h, yout, derivatives);
    y[0]=yout[0];
    y[1]=yout[1];
    output(output_file, h, t, y, E_initial);
    t+=h;
  }
  fclose(output_file);
  return 0;
} //end main program
void initialise(double *initial_x, double *initial_v, double *final_t, int *number_of_steps)
```

```
{
  printf("Read in from screen the initial position x, initial velocity v, final time and number of steps \n");
  scanf("%lf %lf %lf %d", initial_x, initial_v, final_t, number_of_steps);
  return;
} // end of function initialise
// this function provides the first derivative of y[0] and y[1]; i.e.
// it provides the rhs of the couple first order equations of motion
//for the harmonic oscillator with omega=1
void derivatives(double t, double *y, double *dydt)
{
  //double damping_constant = 2.0;
  //double length_rod = 1.0;
  //double gravitational_constant = 1.0;
  //double mass = 1.0;
  // double angular_frequency_initial = sqrt(gravitational_constant/length_rod);
  // double b = damping_constant*angular_frequency_initial/(mass*gravitational_constant);
  //dydt[0] = y[1];
```

```
//dydt[1] = -y[0];
//dydt[1] = -b*y[1]-sin(y[0]);
  //dydt[1] = -y[0]-3.0*pow(y[0],3);
  double damp =2.0;
  double f = 2.5;
  double omega = 2.0;
  dydt[0] = y[1];
  dydt[1] = -y[0]-3.0*pow(y[0],3)-damp*y[1]+f*cos(omega*t);
}// end of function derivatives
// This function computes the first derivative with centered algorithm
void euler(double *y, double *dydt, int n, double t, double h, double *yout)
{
  int i;
  for(i=0; i<n; i++)
  {
    yout[i]=y[i] + h*dydt[i];
  }
} // end of euler integrator
```

```
// function to write out the final results
void output(FILE *output_file, double h, double t, double *y, double E_initial)
{
  fprintf(output_file, "%12.5E \t %12.10E \t %12.10E \t %12.10E \t %12.10E \n", h, t, y[0], y[1],
0.5*y[0]*y[0]+0.5*y[1]*y[1]-E_initial);
  return;
}//end of function output
void runge_kutta_2(double *y, double *dydt, int n, double x, double h, double *yout, void
(*derivatives)(double, double*, double *))
{
  int i;
  double xh;
  double *dyt, *yt;
  // allocate space for local vectors
  dyt = (double *)malloc(n*sizeof(double));
  yt = (double *)malloc(n*sizeof(double));
  xh = x+h/2.0;
  for (i=0; i<n; i++)
  {
    yt[i] = y[i]+h/2.0*dydt[i];
  }
```

```
//computation of y_t+1/2h = yt +h/2 dy/dt using k1 = h dy/dt
  (*derivatives)(xh,yt,dyt);
// find k2 which is h* dyt
  for(i=0; i<n; i++)
  {
    yout[i] = y[i] +h*dyt[i];
  }// increment yout using yout = y+h k2
  free(dyt);
  free(yt);
}//end of function runge katta 2
void runge_kutta_4(double *y, double *dydx, int n, double x, double h, double *yout, void
(*derivatives)(double, double *, double *))
{
  int i;
  double
            xh,hh,h6;
  double *dym, *dyt, *yt;
  // allocate space for local vectors
  dym = (double *) malloc(n*sizeof(double));
  dyt = (double *) malloc(n*sizeof(double));
```

```
yt = (double *) malloc(n*sizeof(double));
hh = h*0.5;
h6 = h/6.;
xh = x+hh;
for (i = 0; i < n; i++)
{
  yt[i] = y[i] + hh*dydx[i];
}
(*derivatives)(xh,yt,dyt); // computation of k2, eq. 3.60
for (i = 0; i < n; i++)
{
  yt[i] = y[i] + hh*dyt[i];
}
(*derivatives)(xh,yt,dym); // computation of k3, eq. 3.61
for (i=0; i < n; i++)
{
  yt[i] = y[i]+h*dym[i];
  dym[i] += dyt[i];
}
(*derivatives)(x+h,yt,dyt); // computation of k4, eq. 3.62
     now we upgrade y in the array yout
for (i = 0; i < n; i++)
{
  yout[i] = y[i]+h6*(dydx[i]+dyt[i]+2.0*dym[i]);
}
```

```
free(dym);
free(dyt);
free(yt);
} // end of function Runge-kutta 4
```