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//Question 2-----  
-----  
  
/*  
newton-raphson method for root finding.  
*/  
  
#include <iostream>  
#include <cstdlib>  
#include <cmath>  
#include <fstream>  
  
using std::endl;  
using std::cout;  
using std::ofstream;  
  
// define function:  
  
double func_y(double x){  
    return sqrt(pow(x,2)+2*x+1)-2*x*sin(x); /* <---- input the function that you want to  
find root on here . */  
}  
  
//define the first derivative of your function:  
  
double func_dydx(double s){  
    return (s+1)/sqrt(pow(s,2)+2*s+1)-2*sin(s)-2*s*cos(s); /* <---- replace it with the d  
ifferentiation of your function*/  
}  
  
int main()  
{  
    int N= 0;  
    int N_limit = 1000;           // max number of steps  
    double s = 2.0;              // initial guess for root,  
    double tol = 1.0e-7;         // tolerance  
  
    ofstream out_file;  
    const char out_file_name[] = "newton.dat"; // save the data in newton.dat  
  
    out_file.open(out_file_name);  
  
        while ((fabs(func_y(s)) >= tol) && (N < N_limit)){  
  
            s = s - func_y(s)/func_dydx(s);  
            ++N;  
  
out_file << N << " " << s << endl; // write the date to output file  
  
        }  
  
if(N >= N_limit){  
    cout << "iteration limit exceeded" << endl;  
    return 0;  
}  
  
if (N <= N_limit){  
    cout << "estimated root = " << s << " with " << N << " iterations." << endl;
```

```
    }

    return 0;
}

/*
bisection method for root finding
of function f(x) in closed interval [a,b].
*/

#include <iostream>
#include <cstdlib>
#include <cmath>
#include <fstream>

using std::endl;
using std::cout;
using std::ofstream;

// define function:

double func_f(double x){
    return tanh(x);
}

int main()
{
    int N= 0;
    int N_limit = 1000;           // max number of steps

    double a = -10.0;   // lower bound for [a,b]
    double b = 3.0;     // upper bound for [a,b]

    double tol = 1.0e-10;   // tolerance f

    ofstream out_file;
    const char out_file_name[] = "Bisection.dat"; // save data in Bisection.dat

    out_file.open(out_file_name);

    if (func_f(a) == 0.0) {
        cout << a << " is a root." << endl;
        return 0;
    }

    if (func_f(b) == 0.0) {
        cout << b << " is a root." << endl;
        return 0;
    }

    if (func_f(a)*func_f(b) > 0.0){
        cout << "[" << a << ", " << b << "]" do not bracket a root." << endl;
        return 0;
    }

    double x_low = a;
    double x_high = b;
    double x_mid;

    while ((fabs(x_high - x_low) >= tol) && (N < N_limit)){
```

```
x_mid = 0.5*(x_low + x_high);
if ( func_f(x_mid)*func_f(x_high) < 0){
    x_low = x_mid;
}
else {
    x_high = x_mid;
}
out_file << x_mid << " " << N << endl; // write date to output file
++N;
}

if(N > N_limit){
    cout << "iteration limit exceeded" << endl;
    return 0;
}

if (N <= N_limit){
    cout << "estimated root = " << x_mid << " with " << N << " iterations." << endl;
}

return 0;
}

// plots the .dat files on gnuplot.

//Question 2-----
-----

//1, Newton-Raphson

/*
newton-raphson.
*/

#include <iostream>
#include <cstdlib>
#include <cmath>
#include <fstream>

using std::endl;
using std::cout;
using std::ofstream;

// define function:

double func_y(double x){
    return sin(x); /* <---- replace it of your function */
}

//define first derivative of function:

double func_dydx(double s){
    return cos(s); /* <---- the differentiation of your function */
}

int main()
{
    int N= 0;
```

```
int N_limit = 1000;           // max number of trys
double s = 0.1;              // initial guess for root, /* <---- for the 2nd functio
n, replace this by 3, or -5 for the 3rd function. */
double tol = 1.0e-4;        // tolerance for finding zero

ofstream out_file;
const char out_file_name[] = "newton.dat"; // save data in fern.dat

out_file.open(out_file_name);

while ((fabs(func_y(s)) >= tol) && (N < N_limit)){

    s = s - func_y(s)/func_dydx(s);
    ++N;

out_file << N << " " << s << endl; // write date to output file

    }

if(N >= N_limit){
    cout << "iteration limit exceeded" << endl;
    return 0;
}

if (N <= N_limit){
    cout << "estimated root = " << s << " with " << N << " iterations." << endl;
}

return 0;
}

//and for Bisection Algorithm, the source code is:

/*
bisection method for root finding
of function f(x) in closed interval [a,b].
*/

#include <iostream>
#include <cstdlib>
#include <cmath>
#include <fstream>

using std::endl;
using std::cout;
using std::ofstream;

// define function:

double func_f(double x){
    return sin(x); // <----replace this with x*sqrt(fabs(x)) for the 2nd function.
}

int main()
{
    int N= 0;
    int N_limit = 1000;           // max number of trys

    double a = -2.9; // lower bound for [a,b]
    double b = 3.0; // upper bound for [a,b]
```

```

double tol = 1.0e-4;           // tolerance for finding zero

ofstream out_file;
const char out_file_name[] = "Bisection.dat"; // save data in fern.dat

out_file.open(out_file_name);

if (func_f(a) == 0.0) {
    cout << a << " is a root." << endl;
    return 0;
}

if (func_f(b) == 0.0) {
    cout << b << " is a root." << endl;
    return 0;
}

if (func_f(a)*func_f(b) > 0.0){
    cout << "[" << a << ", " << b << "]" do not bracket a root." << endl;
    return 0;
}

double x_low = a;
double x_high = b;
double x_mid;

while ((fabs(x_high - x_low) >= tol) && (N < N_limit)){
    x_mid = 0.5*(x_low + x_high);
    if ( func_f(x_mid)*func_f(x_high) < 0){
        x_low = x_mid;
    }
    else {
        x_high = x_mid;
    }
    out_file << x_mid << " " << N << endl; // write date to output file
    ++N;
}

if(N > N_limit){
    cout << "iteration limit exceeded" << endl;
    return 0;
}

if (N <= N_limit){
    cout << "estimated root = " << x_mid << " with " << N << " iterations." << endl;
}

return 0;
}

/*
iterations table:

                                newton           Bisection

sin(x)                          2                16

x*sqrt(fabs(x))                  1                16

*/

```

```
//Question 3-----  
-----  
  
/*  
  
    numerov method to find eigenvalues and eigenfunctions of  
        a particle in a potential well.  
  
*/  
#include <iostream>  
#include <cmath>  
#include <fstream>  
  
using namespace std;  
  
const double V0 = -83.0;           // depth of potential well in MeV  
const double E_min = -83.0;       // energy band to find  
const double E_max = -65.0;       // energy eigenvalue  
const double eps = 1E-6;          // matching tolerance  
double h = 0.004;                 // step size  
                                     // well is assumed 4 fm wide  
                                     // universe is 8 fm wide  
                                     // 2000 total steps  
  
double k2(int i, double E);        /* returns potential at x */  
double diff(double E);             /* difference of derivatives */  
void write_data(double E);         /* saved data for final plot */  
  
int main()  
{  
    double E_mid, E_lo, E_hi;  
    int i=0;                         // counter for iterations  
  
    E_lo = E_min;  
    E_hi = E_max;  
  
    do  
    {  
        ++i;  
        E_mid = (E_lo + E_hi)/2.0;    // first guess for energy eigenvalue  
        if (diff(E_mid)*diff(E_hi) < 0.0) E_lo = E_mid; // the bisection algorithm  
        else E_hi = E_mid;  
    }while(fabs(diff(E_mid))>eps);  
  
    cout << "Eigenvalue E = " << E_mid << endl;  
    cout << "after " << i << " iterations." << endl;  
    write_data(E_mid);  
    return 0;  
}  
  
/*----- end of main program -----*/  
  
/* function returns difference between left and right wavefunction */  
double diff(double E)  
{  
    double psi_now, psi_last, psi_next, psi_left, psi_right;  
    int i;  
  
    psi_last = 0.0;
```

```

psi_now = 0.00001;
for (i=1; i<=1500; ++i)          /* left side first */
{
    psi_next = (2*psi_now*(1 - ((5./12.)*h*h*k2(i,E)))-
                (1. + ((1./12.)*h*h)*k2(i-1,E))*psi_last)/(1.+ ((h*h)/12.)*k2(i+1,E))
;
    psi_last = psi_now;
    psi_now = psi_next;
}
psi_left = psi_now;              /* value at matching point */

psi_last = 0.0;                 /* reset starting conditions */
psi_now = 0.00001;
for (i=1; i<500; ++i)          /* now the right side */
{
    psi_next = (2*psi_now*(1 - ((5./12.)*h*h*k2(i,E)))-
                (1. + ((1./12.)*h*h)*k2(i-1,E))*psi_last)/(1.+ ((h*h)/12.)*k2(i+1,E))
;

    psi_last = psi_now;
    psi_now = psi_next;
}
psi_right = psi_now;            /* value at matching point */
return((psi_left - psi_right));
}

/*-----*/

// function returns k^2 depending on position i
double k2(int i, double E)
{
    const double units_convert = 0.4829;          // MeV^1 fm^-2
    if (i<500) return(units_convert*E);           // outside the well
    if (i>=500) return (units_convert*(E-(V0))); // inside the well
}
/*-----*/

/* write data for eigenfunction into files left.dat, right.dat */
void write_data(double E)
{
    double psi_last, psi_now, psi_next;
    int i;

    ofstream out_file_left;
    ofstream out_file_right;

    out_file_left.open("left.dat"); // save data in files
    out_file_right.open("right.dat");

    psi_last = 0.0;
    psi_now = 0.00001;
    for (i=1; i<=1500; i++) // left side first
    {
        psi_next = (2*psi_now*(1- ((5./12.)*h*h*k2(i,E)))-
                    (1 + ((1./12.)*h*h)*k2(i-1,E))*psi_last)/(1 + ((h*h)/12.)*k2(i+1,E));

        out_file_left << (i-1000)*h << "\t" << psi_next/3.0 << endl;

        psi_last = psi_now;
        psi_now = psi_next;
    }

    psi_last = 0.0;              /* reset starting conditions */

```

```
psi_now = 0.00001;
for (i=1; i<500; i++)      /* now the right side */
{
    psi_next=(2*psi_now*(1- ((5./12.)*h*h*k2(i,E)))-
              (1 + ((1./12.)*h*h)*k2(i-1,E))*psi_last)/(1 + ((h*h)/12.)*k2(i+1,E));

    out_file_right << (1000 -i)*h << "\t" << psi_next/3.0 << endl;

    psi_last = psi_now;
    psi_now = psi_next;
}
out_file_left.close();
out_file_right.close();
cout << "data saved in left.dat and right.dat" << endl;
}
```