



Numerical Computing in C & C++

REPORT

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Preface:

In order to avoid repetition of explanations within this report and comments on the .cpp files I have written below a series of explanations. These explanations are to explain parts of code which are 'basic' but is needed to understand in order to maintain a high level of accuracy when the each code is being run.

`static_cast long double` ('expression') – This particular bit of code allows to avoid accuracy loss when the expression takes a value in the form of int and needs to be redefined as a long double. This is because if we were carry out integer division we would lose the decimal point values as it wont be stored.

`using namespace std`

`cout cin`

`Const`

`#include <iostream>`

`#include <cmath>`

`exp`

Question 1:

Code:

```
#include <iostream>
#include <cmath>
#include <iomanip>
#define EP 1E-15
//this defines the epsilon
using namespace std

int main

    long double      1.0
    //initial guess is defined

    for int      0      2.0
    //for loop will continue until the value of x is more than 2(this won't occur due to the nature of the
function
    long double      exp
    //This outputs the result of the function
    long double      1.0
    //This long double stores number of iterations until the end of the loop or if the loop breaks

    if abs      EP
    //This if statement sets the conditions for the final value of x
    cout      setprecision 18      "The final value of x: "      endl
    //This outputs to 18 digits the final value of x
    cout      "The number of iterations: "      endl
    //This outputs the number iterations occurred stored from the double
    long double      exp
    //This calculates the error
    cout      "The error in the transcendental equation is: "      endl
    break
    //This breaks the entire for loop

    //This re-assignment of doubles allows for the recursive iteration to loop with the new value

    return 0
```

Output:

- A) From the code output, we know the answer is
- B) Since we implemented the long double variable 'iterations' which would increase by 1.0 every time the 'if' loop is executed. When the loop is broken from the *break* function the final value of the number of iterations is 61(as we output the long double 'iterations') which is shown in the code output.
- C) We calculate the transcendental error by using the equation mentioned in the question by inputting our final value of x. The error is: 9.932470248475982208e-16, this is what I expected as I had programmed the 'if' loop to break when $|x_{\text{new}} - x_{\text{old}}|$ is less than or equal to the stated epsilon. In this case, it will always be less than epsilon as it would be very improbable for x to be equal to 1e-15.

Question 2:

Code:

```
#include <iostream>
#include <cmath>
#include <valarray>
#include <iomanip>
using namespace std

long double dot(const valarray<long double>& a, const valarray<long double>& b)
{
    return sum(a*b);
}
//This valarray calculation returns the sum of A*B

//This answers q2a

//The function cdot is defined in week10 notes - compensated summation.cpp - KahanSum function
long double cdot(const valarray<long double>& a)
{
    long double sum = 0.0;
    long double c = 0.0;
    long double s;
    long double t;
    for(int i = 0; i < a.size(); i++)
    {
        s = a[i];
        t = s + c;
        c = (t - s) - c;
        sum = t + c;
    }
    return sum;
}

class normal_cal
{
//using class allows to extract different components out the function
public:
    int m;
    double operator()(const valarray<long double>& a) const
    {
        //This initialises the valarray into the class and uses function object to call on the valarray
        long double sum = 0.0;
        //We let normal be zero as we are going to add to it
        for(int i = 0; i < a.size(); i++)
        {
            //the for loop will go through the entirety of the valarray
            sum += a[i] * a[i];
        }
        //This double computes summation part of the normal
        return pow(static_cast<long double>(sum), 0.5);
    }
    //The function will now return the final value of the normal
};

int main()
{
    cout << setprecision(20);
    //We use the setprecision function to see the accuracy of the programme outputs
    long double a = pow(10.0, 6.0);
    //Defined from question to 10^6 from library cmath
    valarray<long double> a(a);
    //We initialise valarray a with size n
    for(double i = 1; i < a.size(); i++)
    {
        //This creates the valarray of the Euclidean n-valarray
        a[i] = static_cast<long double>(i);
    }
}
```

```

cout    "Product of dot: "    dot    endl
//Using the dot function we sum A and A together
long double    3.1415926535897932385
cout    "Difference in dot and the actual answer: "    dot    pow    2.0 6.0    endl
//This answers q2b

valarray long double    0.1
//This is the constant Euclidean constant n-valarray for the dot function
valarray long double    0.01
//This is the constant Euclidean constant n-valarray for the dot function
//We have to use 0.01 as that is c^2
long double    10000.0
// 10000 = (0.1^2)*(10^6)
// We use the numerical exact value instead to avoid accuracy loss when the programme calculates
the large number

cout    endl
cout    "Product of dot: "    dot    endl
//Using the dot function we sum x and x together to find the sum of nc^2 using the dot function
cout    "Product of cdot: "    cdot    endl
//Using the dot function we sum x_new to find the sum of nc^2 using the KahanSum method
cout    "Difference in cdot and the actual answer: "    cdot    endl
//This answers q2c

normal_cal    2
//This uses the function object and defines the int m = 2
long double
//This uses the function object and defines the valarray object valarray a
cout    endl    "The result of norm l2('valarray'a) is: "    endl
//This answers q2d

return 0

```

Output:

- A) I have defined the dot function before main() and the code explaining it is simply understood as it uses valarray multiplication with .sum() to return the final value.
- B) I first create the valarray A within main() with size n as specified in the question. Then using the dot function I had outputted the final result and the exact value difference which is shown in the output. We get for dot: 1.6449330668477264373, with the difference being:

Since the cdot function already exists in the KahanSum function in week 10 notes (Compensated Summation.cpp) I have just adjusted the valarray x to have the value c squared already instead of having to create two input values for cdot which reduces accuracy loss in the output. We also can see the difference in values from dot and cdot and the exact value. The result of cdot is :

- D)** Using the class `normal_cal` we can implement the double operator()... into the function in order to create a function object. The result of finding the $l_m(A)$ is:

<https://www.symbolab.com/solver/step-by-step/%5Csqrt%7B%5Cfrac%7B%5Cpi%5E%7B2%7D%7D%7B6%7D%7D?or=input>
 $l_m(A)^2 = \text{dot}(A,A)$ (Source: <https://www.symbolab.com/solver/step-by-step/1.2825494403132093879%5E%7B2%7D?or=input>) if we compute it within the .cpp file which is implied from the previous statement of square rooting the exact value.

Question 3

Code:

```
#include <iostream>
#include <cmath>
#include <valarray>
#include <iomanip>
using namespace std;

long double normal_calculation(int i, valarray<long double> &f, valarray<long double> &fddx)
//We have turned the normal_cal function object into a function so it's easier to use to avoid longer
//calculation times and less re-defining of valarrays when the function is called several times
{
    long double result = 0;
    for (int j = 0; j < f.size(); j++)
        result += f[j] * pow(abs(f[j] - f[j+1]), 2);

    long double result2 = pow(result, 1/2);
    //This takes the normal formula from 2d in the form of a function instead of function object
    return result2;
}

void fanalytical(int N, valarray<long double> &f, valarray<long double> &fddx)
//This initialises the valarray to the correct size for given N
{
    long double result = 0;
    //We calculate the long double here which is required to calculate other doubles
    for (int i = 0; i < N; i++)
    {
        long double xi = 2.0 * M_PI * i / N;
        //xi formula
        long double fxi = exp(16.0 * pow(xi, 2.0)) * pow(32.0, 32.0);
        //f(xi) formula
        long double fddxi = 2.0 * fxi;
        //fddx formula
        f[i] = fxi;
        fddx[i] = fddxi;
    }
}

// Write header, in week 9 lab code
#define SP << setw(30) << setprecision(10) << // save some repetition when writing
cout << "i:" SP "xi:" SP "f(xi):" SP "fddx(xi):" SP "ei:" endl
for (int i = 0; i < N; i++)
    cout << SP << f[i] << SP << fddx[i] << SP << endl
//This generates the headers when tabulating the outputs

long double e; int
```

```

valarray long double      1
valarray long double      1
valarray long double      1

long double      pow 2.0 static_cast long double      2.0
//We calculate the long double here which is required to calculate other doubles

for int      0      1
    double      //xi formula
    exp 16.0 pow      2.0 //f(xi) formula
    32.0 32.0 pow      2.0
//This series of for loops inputs the correct values into each respective valarray

cout setprecision 10
valarray long double      1
valarray long double      1
for int      0      1
    if      0
        0      2 2.0      1
        0      0      0
    else if
        2.0      1      2

    else
        1 2.0      1

//The above is already explained but we need the previous values to calculate the error
cout setprecision 15
long double      static_cast long double      static_cast long
double      1 normal_calculation 1
//This long double calculates the final value error from the value of n inputted
cout "The result of (N^2)<e>: " " N = " endl
return

int main

fanalytical 63

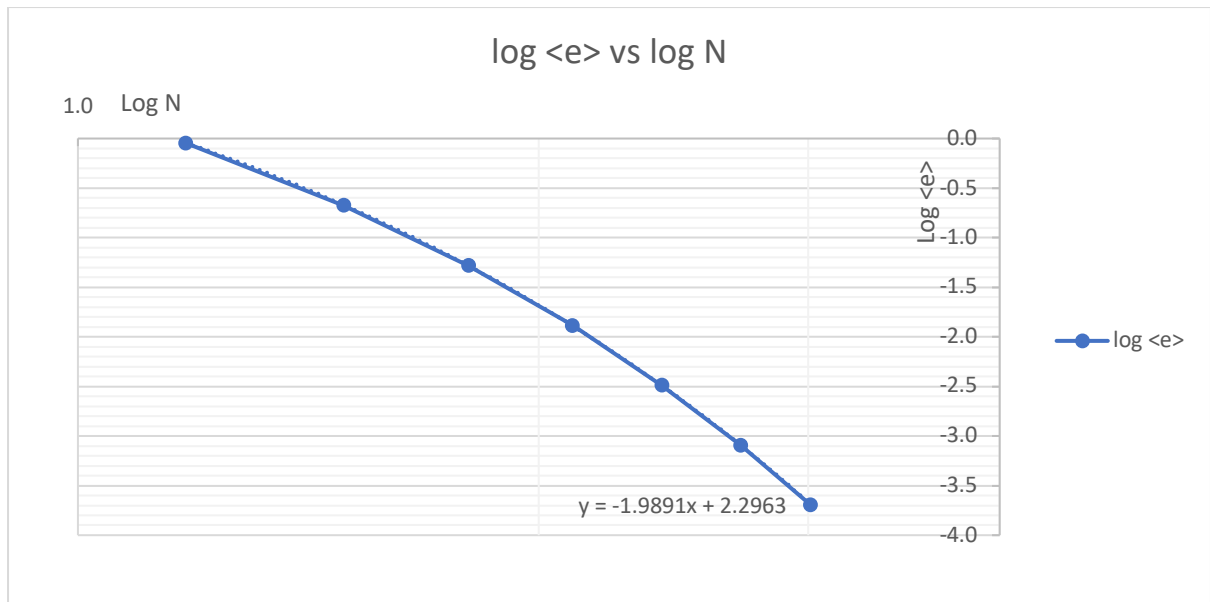
int 15
while 2048
    e
    2 1

return 0

```


i:	xi:	f(xi):	fddx(xi):	ei:
0	-1.0000000000E+00	1.1253517470E-07	3.0325028050E-04	-1.9161538720E-04
1	-9.6825396830E-01	3.0582525730E-07	3.0325028050E-04	-1.9439510260E-05
2	-9.3650793650E-01	8.0473416630E-07	7.4090850990E-04	-4.3931269510E-05
3	-9.0476190480E-01	2.0503384740E-06	1.7485617630E-03	-9.5496127630E-05
4	-8.7301587300E-01	5.0581617430E-06	3.9853066720E-03	-1.9953322570E-04
5	-8.4126984130E-01	1.2082419050E-05	8.7701453610E-03	-4.0040497600E-04
6	-8.0952380950E-01	2.7945321210E-05	1.8629553820E-02	-7.7090936670E-04
7	-7.7777777780E-01	6.2583283910E-05	3.8187293640E-02	-1.4223185560E-03
8	-7.4603174600E-01	1.3570680330E-04	7.5510389870E-02	-2.5108658700E-03
9	-7.1428571430E-01	2.8493048890E-04	1.4397690970E-01	-4.2330422030E-03
10	-6.8253968250E-01	5.7925562040E-04	2.6459093940E-01	-6.7983992840E-03
11	-6.5079365080E-01	1.1402382870E-03	4.6839726220E-01	-1.0366788600E-02
12	-6.1904761900E-01	2.1732766480E-03	7.9822669370E-01	-1.4940336510E-02
13	-5.8730158730E-01	4.0107762770E-03	1.3084792130E+00	-2.0212618550E-02
14	-5.5555555560E-01	7.1669750380E-03	2.0611753540E+00	-2.5400518920E-02
15	-5.2380952380E-01	1.2400448030E-02	3.1163492080E+00	-2.9115670420E-02
16	-4.9206349210E-01	2.0774610570E-02	4.5153801800E+00	-2.9362383040E-02
17	-4.6031746030E-01	3.3699420810E-02	6.2573972800E+00	-2.3760098410E-02
18	-4.2857142860E-01	5.2930501930E-02	8.2715389850E+00	-1.0059826070E-02
19	-3.9682539680E-01	8.0497727160E-02	1.0391247510E+01	1.3058370220E-02
20	-3.6507936510E-01	1.1853736110E-01	1.2339938960E+01	4.5049998080E-02
21	-3.3333333330E-01	1.6901331540E-01	1.3738815290E+01	8.2718057280E-02
22	-3.0158730160E-01	2.3333539250E-01	1.4145669560E+01	1.1994722350E-01
23	-2.6984126980E-01	3.1191362430E-01	1.3127321060E+01	1.4830119450E-01
24	-2.3809523810E-01	4.0372170860E-01	1.0358281690E+01	1.5862304310E-01
25	-2.0634920630E-01	5.0596897830E-01	5.7267759290E+00	1.4343306840E-01
26	-1.7460317460E-01	6.1398775300E-01	-5.7970953320E-01	9.9533673940E-02
27	-1.4285714290E-01	7.2142229040E-01	-8.0392287710E+00	2.9969057300E-02
28	-1.1111111110E-01	8.2075480830E-01	-1.5832696630E+01	-5.5495213470E-02
29	-7.9365079370E-02	9.0413096780E-01	-2.2958712730E+01	-1.4184488810E-01
30	-4.7619047620E-02	9.6436909490E-01	-2.8408464310E+01	-2.1208606360E-01
31	-1.5873015870E-02	9.9597687240E-01	-3.1362817260E+01	-2.5148112580E-01
32	1.5873015870E-02	9.9597687240E-01	-3.1362817260E+01	-2.5148112580E-01
33	4.7619047620E-02	9.6436909490E-01	-2.8408464310E+01	-2.1208606360E-01
34	7.9365079370E-02	9.0413096780E-01	-2.2958712730E+01	-1.4184488810E-01
35	1.1111111110E-01	8.2075480830E-01	-1.5832696630E+01	-5.5495213470E-02
36	1.4285714290E-01	7.2142229040E-01	-8.0392287710E+00	2.9969057300E-02
37	1.7460317460E-01	6.1398775300E-01	-5.7970953320E-01	9.9533673940E-02
38	2.0634920630E-01	5.0596897830E-01	5.7267759290E+00	1.4343306840E-01

39	2.3809523810E-01	4.0372170860E-01	1.0358281690E+01	1.5862304310E-01
40	2.6984126980E-01	3.1191362430E-01	1.3127321060E+01	1.4830119450E-01
41	3.0158730160E-01	2.3333539250E-01	1.4145669560E+01	1.1994722350E-01
42	3.3333333330E-01	1.6901331540E-01	1.3738815290E+01	8.2718057280E-02
43	3.6507936510E-01	1.1853736110E-01	1.2339938960E+01	4.5049998080E-02
44	3.9682539680E-01	8.0497727160E-02	1.0391247510E+01	1.3058370220E-02
45	4.2857142860E-01	5.2930501930E-02	8.2715389850E+00	-1.0059826070E-02
46	4.6031746030E-01	3.3699420810E-02	6.2573972800E+00	-2.3760098410E-02
47	4.9206349210E-01	2.0774610570E-02	4.5153801800E+00	-2.9362383040E-02
48	5.2380952380E-01	1.2400448030E-02	3.1163492080E+00	-2.9115670420E-02
49	5.5555555560E-01	7.1669750380E-03	2.0611753540E+00	-2.5400518920E-02
50	5.8730158730E-01	4.0107762770E-03	1.3084792130E+00	-2.0212618550E-02
51	6.1904761900E-01	2.1732766480E-03	7.9822669370E-01	-1.4940336510E-02
52	6.5079365080E-01	1.1402382870E-03	4.6839726220E-01	-1.0366788600E-02
53	6.8253968250E-01	5.7925562040E-04	2.6459093940E-01	-6.7983992840E-03
54	7.1428571430E-01	2.8493048890E-04	1.4397690970E-01	-4.2330422030E-03
55	7.4603174600E-01	1.3570680330E-04	7.5510389870E-02	-2.5108658700E-03
56	7.7777777780E-01	6.2583283910E-05	3.8187293640E-02	-1.4223185560E-03
57	8.0952380950E-01	2.7945321210E-05	1.8629553820E-02	-7.7090936670E-04
58	8.4126984130E-01	1.2082419050E-05	8.7701453610E-03	-4.0040497600E-04
59	8.7301587300E-01	5.0581617430E-06	3.9853066720E-03	-1.9953322570E-04
60	9.0476190480E-01	2.0503384740E-06	1.7485617630E-03	-9.5496127630E-05
61	9.3650793650E-01	8.0473416630E-07	7.4090850990E-04	-4.3931269510E-05
62	9.6825396830E-01	3.0582525730E-07	3.0325028050E-04	-1.9439510260E-05
63	1.0000000000E+00	1.1253517470E-07	3.0325028050E-04	-1.9161538720E-04



B)

Code:

```
#include <iostream>
#include <valarray>
#include <cmath>
#include <iomanip>
#include <random>
using namespace std

long double dot(const valarray<long double>& a, const valarray<long double>& b)
{
    return sum(a*b);
}
//This function is recalled from question 2a

long double monteCarloEstimate(long double a, long double b, long double c)
//Function to execute Monte Carlo integration on predefined function
{
    long double sum = 0.0;
    const int n = 31;
    //seed value
    mt19937_64 rng;
    //We are generating a seed with a real random value
    uniform_real_distribution<long double> dis(a, b);
    //This generates random values uniformly between the lower and upper bound

    for (int i = 0; i < n; ++i)
    {
        //Select a random number within the limits of integration
        long double x = dis(rng);
        //Add the f(x) value to the running sum using the correct function
        sum += pow(x, c);
    }

    long double result = static_cast<long double>(sum/n);
    //This calculates the final value of the Integration estimate using this method
    return result;
}

int main()
{
    long double n = 63.0;
    //This is the value of n defined in the question
    long double delta_x = static_cast<long double>(1/n);
    //We calculate delta x using the formula given, we use static_cast as this a fraction to avoid accuracy loss

    valarray<long double> x(n);
    for (int i = 0; i < x.size(); ++i)
        x[i] = double(i*delta_x);
    //This calculates the gridpoints for 4a/4b and inputs them into a valarray

    valarray<long double> f(n);
    for (int i = 0; i < f.size(); ++i)
        f[i] = pow(x[i], 4.0);
    //This calculates the f(i) points for 4a/4b and inputs them into a valarray

    valarray<long double> y(n);
    for (int i = 0; i < y.size(); ++i)
        y[i] = static_cast<long double>(f[i]*delta_x);
    if (i == 0)
        //This inputs the values into the valarray for when i is 0
    }
```

```

else if

    //This inputs the values into the valarray for when i is 63

else
    2.0
    //This inputs the values into the valarray

//This for loop inputs in the correct values for the weights creates the values inputted into the
valarray

valarray long double 1
for int 0 size
    long double static_cast long double 48.0
    if 0
        17.0
        //This inputs the values into the valarray for when i is 0

    else if 1
        59.0
        //This inputs the values into the valarray for when i is 1

    else if 2
        43.0
        //This inputs the values into the valarray for when i is 2

    else if 3
        49.0
        //This inputs the values into the valarray for when i is 3

    else if 3
        49.0
        //This inputs the values into the valarray for when i is 60

    else if 2
        43.0
        //This inputs the values into the valarray for when i is 61

    else if 1
        59.0
        //This inputs the values into the valarray for when i is 62

    else if
        17.0
        //This inputs the values into the valarray for when i is 63

    else
        48.0
        //This inputs the values into the valarray

//This for loop inputs in the correct values for the weights creates the values inputted into the
valarray

long double 3.1415926535897932385

valarray long double 1
for int 0 size
    static_cast long double
    //This calculates the theta values for 4c and inputs them into a valarray

```

```

valarray long double          1
for int 0          size
    4.0 static_cast double 4 cos 2.0
//This calculates the gridpoints for 4c and inputs them into a valarray

valarray long double          1
for int 0          size
    pow 4.0          0.5
//This calculates the f(i) for 4c and inputs them into a valarray

valarray long double          1
for int 0          size
    if 0
        static_cast double 2.0 pow 2.0
        //This inputs the values into the valarray for when i is 0

    else if
        static_cast double 2.0 pow 2.0
        //This inputs the values into the valarray for when i is 63

    else
        long double          0
        for double 1.0 32
            static_cast double 2.0 cos 2.0          4.0 pow 2.0 1.0
            //This generates the summation part of the CC equation

            2.0 static_cast double 2.0 1.0
            //This inputs the values for the valarray for the correct values of i

//This for loop inputs in the correct values for the weights creates the values inputted into the
valarray

long double          2.0
long double          dot
long double          dot
long double          dot
//We use the function from 2a to calculate the porduct of two valarrays
long double monteCarloEstimate 0.0 4.0 10000.0
//This calls the function defined to calculate the MC estimate for 4d

//The below outputs all the results for question 4
cout setprecision 10
cout "i_real: "          endl          endl
cout "i_trapezium: "          endl
cout "i_trapezium - i_real = "          endl          endl
cout "i_simpson: "          endl
cout "i_simpson - i_real = "          endl          endl
cout "i_CC: "          endl
cout "i_trapezium - i_real = "          endl          endl
cout "i_MC: "          endl
cout "i_MC - i_real = "          endl

return 0

```

- A)** Using the dot function from 2a, we get the answer for 4a in the first two rows of the output past `i_real` which gives our I_{exact} . The error is what I expected as this method is not the most accurate method. This is shown by $I_{\text{trapezium}} =$ $I_{\text{trapezium}} - I_{\text{exact}} =$

Using the dot function from 2a, we get the answer for 4b in the second two rows of the output past `i_real`. The error is what I expected as this method is not the most accurate method but it is more accurate compared to the trapezium method. This is shown by $I_{\text{Simpson}} = 6$.

$I_{\text{Simpson}} - I_{\text{exact}} =$

- C)** Using the dot function from 2a, we get the answer for 4c in the third two rows of the output past `i_real`. The error is what I expected as this method is the most accurate method. This is shown by $I_{\text{ClenshawCurtis}} =$ $I_{\text{ClenshawCurtis}} - I_{\text{exact}} =$ Extended Simpson method.

Using the dot function from 2a, we get the answer for 4d in the last two rows of the output past `i_real`. The error is what I expected as this method is a more accurate method compared to the Extended Simpson method. However, the accuracy is highly dependent on the seed value implemented into the random number generator in the function 'monteCarloEstimate'. $I_{\text{MonteCarlo}} =$ $I_{\text{MonteCarlo}} - I_{\text{exact}} =$

Question 5a, 5b

Code:

```
#include <iostream>
#include <iomanip>
#include <cmath>
#include <valarray>
using namespace std

valarray<long double> F(const long double, const valarray<long double>
/* du/dt = f(t,u) = { u0
    3(u1)^2 + (u1)
    since a =0,c=1(given) we get the above
*/
    valarray<long double> size
    // u[0] = dq/dt, u[1] = dp/dt
    //Insert the relevant equations into the f valarray
    0 1
    1 3.0 pow static_cast<long double> 0 2.0 0
    return
    //This function will return the correct values of dp/dt and dq/dt by using a valarray<valarray<... in
    main()

valarray<long double> RK2(const long double, const valarray<long double> const long double
    valarray<long double> const long double const valarray<long double>
    unsigned long long size
    //This allows for us to get the value of the size of the valarray we are using to stay consistent
    valarray<long double>
    //This initialises valarrays of the correct size

    //Since we call on function f and don't declare any specific valarray we are able to manipulate the
    function to our use different valarrays
    0.5 0.5
    return
    //The above formula is used the lecture notes - Page 5 of ODEs

int main

    long double 0.
    long double 10.0
    int 100000
    long double static_cast<long double>
    valarray<long double> 2 // [0],T[1],...,T[n],T[n+1]
    //Given from question

    valarray<valarray<long double>> 2
    valarray<long double> 2
    valarray<long double> 2

    //set initial data such that U(0) = {q(0),p(0)} = {-0.5,0}
    0 0.5 0.

    for(int 0
        static_cast<long double>
        //given in question
        1 RK2 F
        //We use the +1 as we already have been given the initial conditions, we do not want to overwrite
        it - we also find the 0th value of each valarray without the need for 2 'for' loops
        static_cast<long double> 0.5 j 1 j 1 static_cast<long
    double j 0 j 0 j 0 static_cast<long double> 0.5 j 0 j 0
```



```

//E(t) = 0.5(p^2) + (V(q)) = 0.5(p*p) - (q*q*q) -(0.5*q*q)
j 0 static_cast<long double> 0.5 cosh 0.5 cosh 0.5
//error = q(t) - (-0.5sech^2(0.5*t)) = q(t) - (-0.5/cosh^2(0.5*t))
//We have to use cosh identity instead of sech as only cosh is cmath library

// Write header
#define SP << setw(26) << setprecision(10) << // save some repetition when writing
cout << "t" SP "q(t)" SP "p(t)" SP "E(t)" SP "e(t)" << endl

int 0
while
    cout << SP << i << SP << i << SP << SP << endl
    10000
//We use a while function and this equation to output the only relevant time intervals for
i=0,i=1...i=10

return 0

```

Output:

t	q(t)	p(t)	E(t)	e(t)
0	-5.0000000000E-01	0.0000000000E+00	0.0000000000E+00	0.0000000000E+00
1	-3.9322386600E-01	1.8171549550E-01	-5.2467299170E-12	5.2732825060E-10
2	-2.0998717020E-01	1.5992500170E-01	-2.0739801250E-11	6.0979269530E-10
3	-9.0353319320E-02	8.1783148900E-02	-2.5493658060E-11	1.3709330730E-10
4	-3.5325412760E-02	3.4054671400E-02	-2.5990213230E-11	-3.3179652270E-10
5	-1.3296114370E-02	1.3118134550E-02	-2.6023080210E-11	-1.0259183440E-09
6	-4.9330212880E-03	4.9086209670E-03	-2.6024900570E-11	-2.7049412680E-09
7	-1.8204495770E-03	1.8171181970E-03	-2.6024994840E-11	-7.2171302700E-09
8	-6.7049481720E-04	6.7000626130E-04	-2.6024999610E-11	-1.9475678440E-08
9	-2.4681149980E-04	2.4664508290E-04	-2.6024999850E-11	-5.2800230780E-08
10	-9.0935003930E-05	9.0640063020E-05	-2.6024999860E-11	-1.4338845530E-07

Question 5c

Code:

```
#include <iostream>
#include <iomanip>
#include <cmath>
#include <valarray>
using namespace std

valarray<long double> F(const long double, const valarray<long double>
/* du/dt = f(t,u) = { u0
    3(u1)^2 + (u1)
    since a =0,c=1(given) we get the above
*/
    valarray<long double> size
    // u[0] = dq/dt, u[1] = dp/dt
    //Insert the relevant equations into the f valarray
    0 1
    1 3.0 pow static_cast<long double> 0 2.0 0
    return
    //This function will return the correct values of dp/dt and dq/dt by using a valarray<valarray<... in
    main()

//We use the pre-defined function Hermit-2 as it uses the trapezium rule without having to alter the
function F values
valarray<long double> H2(const long double, const valarray<long double>, const long double
    valarray<long double> const long double const valarray<long double>
    long double //This increases the time by delta_t starting from t_initial
    unsigned int 10 //This is the maximum number of self-consistent iterations (10) that will be
    used
    valarray<long double> //This sets initial guess for the implicit ODE
    for int 0
        //This 'for' loop carries out the self-consistent iterations
        0.5 //trapezium rule is now applied to the new values

    return

int main

    long double 0.
    long double 10.0
    int 100000
    long double static_cast<long double>
    valarray<long double> 2 // [0], T[1], ..., T[n], T[n+1]
    valarray<valarray<long double>> 2
    valarray<long double> 2
    valarray<long double> 2
    //Given from question

    // Write header
#define SP << setw(26) << setprecision(10) << // save some repetition when writing
    cout << "t" SP "q(t)" SP "p(t)" SP "E(t)" SP "e(t)" << endl

    //set initial data such that U(0) = {q(0),p(0)} = {-0.5,0}
    0 0.5 0.

    for int 0
        static_cast<long double>
        //given in question
```

```

1  H2          F
//We use the +1 as we already have been given the initial conditions, we do not want to overwrite
it - we also find the 0th value of each valarray without the need for 2 'for' loops
static_cast<long double>(0.5) j 1 j 1 static_cast<long
double> j 0 j 0 j 0 static_cast<long double>(0.5) j 0 j 0
//E(t) = 0.5(p^2) + (V(q)) = 0.5(p*p) - (q*q*q) -(0.5*q*q)
j 0 static_cast<long double>(0.5) cosh 0.5 cosh 0.5
//error = q(t) - (-0.5sech^2(0.5*t)) = q(t) - (-0.5/cosh^2(0.5*t))
//We have to use cosh identity instead of sech as only cosh is cmath library

int 0
while
cout SP i 0 SP i 1 SP SP endl
10000
//We use a while function and this equation to output the only relevant time intervals for
i=0,i=1...i=10

return 0

```

Output:

t	q(t)	p(t)	E(t)	e(t)
0	-5.0000000000E-01	0.0000000000E+00	0.0000000000E+00	0.0000000000E+00
1	-3.9322386670E-01	1.8171549530E-01	1.0518493570E-11	-2.4935283190E-10
2	-2.0998717100E-01	1.5992500250E-01	4.1511998940E-11	-1.9208806810E-10
3	-9.0353319240E-02	8.1783149770E-02	5.1020583580E-11	2.2312316660E-10
4	-3.5325411610E-02	3.4054672630E-02	5.2013756370E-11	8.1606356960E-10
5	-1.3296111260E-02	1.3118137480E-02	5.2079493920E-11	2.0833425880E-09
6	-4.9330131700E-03	4.9086288420E-03	5.2083134830E-11	5.4129834200E-09
7	-1.8204279210E-03	1.8171396040E-03	5.2083323390E-11	1.4438908260E-08
8	-6.7043636860E-04	6.7006446620E-04	5.2083332920E-11	3.8972944540E-08
9	-2.4665303330E-04	2.4680331010E-04	5.2083333390E-11	1.0566625470E-07
10	-9.0504655690E-05	9.1070174610E-05	5.2083333420E-11	2.8695978550E-07