Newtonian gravitation for C++ programmers

S. Halayka*

Friday 29th November, 2024 09:19

Abstract

...

1 Typedefs

```
typedef long double real_type;
```

or

```
#include <boost/multiprecision/cpp_bin_float.hpp>
using namespace boost::multiprecision;

typedef number<
        backends::cpp_bin_float<
                237,
                backends::digit_base_2,
                void,
                std::int32_t,
                -262142,
                262143>,
        et_off> cpp_bin_float_oct;

typedef cpp_bin_float_oct real_type;
```

2 Constants

```
const real_type pi = 4.0 * atan(1.0);
const real_type G = 6.67430e-11;
const real_type c = 299792458;
const real_type c2 = c * c;
const real_type c3 = c * c * c;
const real_type c4 = c * c * c * c;
const real_type h = 6.62607015e-34;
const real_type hbar = h / (2.0 * pi);
```

---

*sjhalayka@gmail.com

3 Brute force: integer field line count

Where r is the receiver radius, R is the distance from the centre of the emitter, β is the get intersecting line count function, and n is the field line count, the gradient is:

$$\alpha = \frac{\beta(R + \epsilon) - \beta(R)}{\epsilon}. \tag{1}$$

The gradient strength is:

$$g = \frac{-\alpha}{r^2}. \tag{2}$$

```cpp
long long unsigned int get_intersecting_line_count(
        const vector<vector_3>& unit_vectors,
        const vector_3& sphere_location,
        const real_type sphere_radius)
{
        long long unsigned int count = 0;

        vector_3 cross_section_edge_dir(sphere_location.x, sphere_radius, 0);
        cross_section_edge_dir.normalize();

        vector_3 receiver_dir(sphere_location.x, 0, 0);
        receiver_dir.normalize();

        const real_type min_dot = cross_section_edge_dir.dot(receiver_dir);

        for (size_t i = 0; i < unit_vectors.size(); i++)
                if (unit_vectors[i].dot(receiver_dir) >= min_dot)
                        count++;

        return count;
}
```

```cpp
int main(int argc, char** argv)
{
        // Field line count
        const size_t n = 1000000000;

        cout << "Allocating memory for field lines" << endl;
        vector<vector_3> unit_vectors(n);

        for (size_t i = 0; i < n; i++)
        {
                unit_vectors[i] = RandomUnitVector();

                static const size_t output_mod = 10000;

                if (i % output_mod == 0)
                        cout << "Getting pseudorandom locations: "
                        << static_cast<float>(i) / n << endl;
        }

        string filename = "newton.txt";
```

```cpp
        ofstream out_file(filename.c_str());
        out_file << setprecision(30);

        const real_type start_distance = 10.0;
        const real_type end_distance = 100.0;
        const size_t distance_res = 1000;

        const real_type distance_step_size =
                (end_distance - start_distance)
                / (distance_res - 1);

        for (size_t step_index = 0; step_index < distance_res; step_index++)
        {
                const real_type r =
                        start_distance +
                        step_index * distance_step_size;

                const vector_3 receiver_pos(r, 0, 0);
                const real_type receiver_radius = 1.0;

                const real_type epsilon = 1.0;

                vector_3 receiver_pos_plus = receiver_pos;
                receiver_pos_plus.x += epsilon;

                const long long signed int collision_count_plus =
                        get_intersecting_line_count(
                                unit_vectors,
                                receiver_pos_plus,
                                receiver_radius);

                const long long signed int collision_count =
                        get_intersecting_line_count(
                                unit_vectors,
                                receiver_pos,
                                receiver_radius);

                const real_type gradient =
                        static_cast<real_type>
                        (collision_count_plus - collision_count)
                        / epsilon;

                const real_type gradient_strength =
                        -gradient
                        / (receiver_radius * receiver_radius);

                cout << "r: " << r << " gradient strength: "
                << gradient_strength << endl;

                out_file << r << " " << gradient_strength << endl;
        }

        out_file.close();
```

```
        return 0;
}
```

While this method works, it is both memory and processor intensive. This method is meant to be a stepping stone for the next section. See Fig 1.

4   Heuristic: real field line count

Where r is the receiver radius, R is the distance from the centre of the emitter, β is the get intersecting line count function, and n is the field line count, the gradient is:

$$\alpha = \frac{\beta(R + \epsilon) - \beta(R)}{\epsilon}. \tag{3}$$

Here we assume that the maximum number of field lines is given by the holographic principle:

$$n = \frac{Ac^3}{4G\hbar \log 2}. \tag{4}$$

The gradient strengths are:

$$g = \frac{-\alpha}{r^2} \approx \frac{n}{2R^3}, \tag{5}$$

$$g_N = \frac{gRc\hbar \log 2}{2\pi M} = \frac{nc\hbar \log 2}{4\pi MR^2} = \frac{Ac^4}{16\pi GMR^2} = \frac{GM}{R^2}. \tag{6}$$

We will use g_N in the next section.

```
real_type get_intersecting_line_count (
        const real_type n,
        const vector_3& sphere_location ,
        const real_type sphere_radius )
{
        const real_type big_area =
                4 * pi * sphere_location.x * sphere_location.x;

        const real_type small_area =
                pi * sphere_radius * sphere_radius ;

        const real_type ratio =
                small_area / big_area ;

        return n * ratio ;
}
```

```
int main(int argc , char** argv)
{
        const real_type emitter_radius = 1.0;

        const real_type emitter_area =
                4.0 * pi * emitter_radius * emitter_radius ;

        // Field line count
```

```cpp
// re: holographic principle:
const real_type n =
        (c3 * emitter_area)
        / (log(2.0) * 4.0 * G * hbar);

const real_type emitter_mass = c2 * emitter_radius / (2.0 * G);

// 1.73502e+70 is the 't Hooft-Susskind constant:
// the number of field lines for a black hole of
// unit Schwarzschild radius
//
//const real_type G_ =
//        (c3 * pi)
//        / (log(2.0) * hbar * 1.73502e+70);

const string filename = "newton.txt";
ofstream out_file(filename.c_str());
out_file << setprecision(30);

const real_type start_distance = 10.0;
const real_type end_distance = 100.0;
const size_t distance_res = 1000;

const real_type distance_step_size =
        (end_distance - start_distance)
        / (distance_res - 1);

for (size_t step_index = 0; step_index < distance_res; step_index++)
{
        const real_type r =
                start_distance + step_index * distance_step_size;

        const vector_3 receiver_pos(r, 0, 0);
        const real_type receiver_radius = 1.0;

        const real_type epsilon = 1.0;

        vector_3 receiver_pos_plus = receiver_pos;
        receiver_pos_plus.x += epsilon;

        // https://en.wikipedia.org/wiki/Directional_derivative
        const real_type collision_count_plus =
                get_intersecting_line_count(
                        n,
                        receiver_pos_plus,
                        receiver_radius);

        const real_type collision_count =
                get_intersecting_line_count(
                        n,
                        receiver_pos,
                        receiver_radius);

        const real_type gradient =
```

```
                            ( collision_count_plus − collision_count )
                            / epsilon ;

                const real_type gradient_strength =
                            −gradient
                            / ( receiver_radius ∗ receiver_radius );

                const real_type gradient_strength_ =
                            n / (2.0 ∗ pow( receiver_pos.x, 3.0 ));

                const real_type newton_strength =
                            n ∗ c ∗ hbar ∗ log (2.0)
                            /
                            (pow( receiver_pos.x, 2.0)
                                    ∗ emitter_mass ∗ 4.0 ∗ pi );

                const real_type newton_strength_ =
                            c4 ∗ emitter_area
                            / (16.0 ∗ pi ∗ G
                                    ∗ pow( receiver_pos.x, 2.0) ∗ emitter_mass );

                const real_type newton_strength__ =
                            G ∗ emitter_mass / pow( receiver_pos.x, 2.0);

                const real_type newton_strength___ =
                            gradient_strength_ ∗ receiver_pos.x
                            ∗ c ∗ hbar ∗ log (2)
                            / (2 ∗ pi ∗ emitter_mass );

                //cout << newton_strength___ / newton_strength << endl;

                cout << "r: " << r << " gradient strength: "
                            << gradient_strength << endl;

                out_file << r << " " << gradient_strength << endl;
        }

        out_file.close ();

        return 0;
}
```

This method is faster and less memory intensive when compared to the integer field count method. This method is meant to be a stepping stone for the next section.

5   Application: modeling Mercury's orbit using numerical integration

In essence, the numerical calculation of the Newtonian orbit of Mercury is as follows:

- Place Mercury at the aphelion to start

- Calculate the orbit path by repeatedly taking steps in time

The constant time slice is:

```
const real type dt = 10000; // 2.77777 hours
```

The initial conditions are:

```
vector_3 Mercury_pos(0, 69817079000.0, 0); // Aphelion location
vector_3 Mercury_vel(-38860, 0, 0); // Aphelion velocity
```

The orbit code is as follows. Here we use Eq. 6 (e.g. $g_N = GM/R^2$) to calculate the acceleration from Newtonian gravitation:

```
vector_3 Newtonian_acceleration(
        const real_type emitter_mass,
        const vector_3& pos, // Receiver pos
        const real_type G)
{
        // Sun's position is fixed at the origin
        vector_3 grav_dir = vector_3(0, 0, 0) - pos;
        const real_type distance = grav_dir.length();
        grav_dir.normalize();

        vector_3 accel = grav_dir * G * emitter_mass / pow(distance, 2.0);

        return accel;
}
```

Here we show the Euler integration, which is extremely simple. The acceleration is calculated, then it is added (e.g. integrated) to the velocity. Once that's done, the velocity is added to the position.

```
void proceed_Euler(
        vector_3& pos,
        vector_3& vel,
        const real_type G,
        const real_type dt)
{
        vector_3 accel =
                Newtonian_acceleration(
                        emitter_mass,
                        pos,
                        G);

        vel += accel * dt;
        pos += vel * dt;
}
```

And so the passage of time is computed as:

```
void idle_func(void)
{
        proceed_Euler(Mercury_pos, Mercury_vel, G, dt);
}
```

On the other hand, rather than using Euler integration, the order-4 symplectic integration does a better job at conserving energy, but at a speed cost:

```
void proceed_symplectic_order_4 (
        vector_3& pos,
        vector_3& vel,
        real_type G,
        real_type dt)
{
        static const real_type cr2 =
                pow(2.0, 1.0 / 3.0);

        static const real_type c[4] =
        {
                1.0 / (2.0 * (2.0 - cr2)),
                (1.0 - cr2) / (2.0 * (2.0 - cr2)),
                (1.0 - cr2) / (2.0 * (2.0 - cr2)),
                1.0 / (2.0 * (2.0 - cr2))
        };

        static const real_type d[4] =
        {
                1.0 / (2.0 - cr2),
                -cr2 / (2.0 - cr2),
                1.0 / (2.0 - cr2),
                0.0
        };

        pos += vel * c[0] * dt;
        vel += Newtonian_acceleration(
                        emitter_mass,
                        pos,
                        G) * d[0] * dt;

        pos += vel * c[1] * dt;
        vel += Newtonian_acceleration(
                        emitter_mass,
                        pos,
                        G) * d[1] * dt;

        pos += vel * c[2] * dt;
        vel += Newtonian_acceleration(
                        emitter_mass,
                        pos,
                        G) * d[2] * dt;

        pos += vel * c[3] * dt;
        // last element d[3] is always 0
}
```

See Fig. 2.

6   Final code

A final code, which models the orbit of Mercury, is at:

```
https://github.com/sjhalayka/mercury_orbit_glut
```

References

[1] Misner et al. Gravitation. (1970)

[2] 't Hooft. Dimensional reduction in quantum gravity. (1993)

[3] Susskind. The World as a Hologram. (1994)

[4] Fiedler. Fix Your Timestep! (2004)
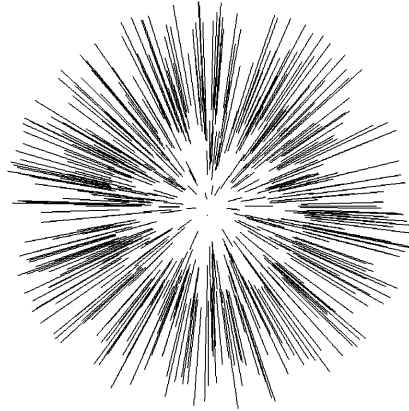
[5] Fiedler. Integration Basics. (2004)

Figure 1: Example of an isotropic emitter. The emitter is spherical. The field line starting locations are placed pseudorandomly on a 2-sphere, and the normals (e.g. field line directions) are calculated using the same sphere.
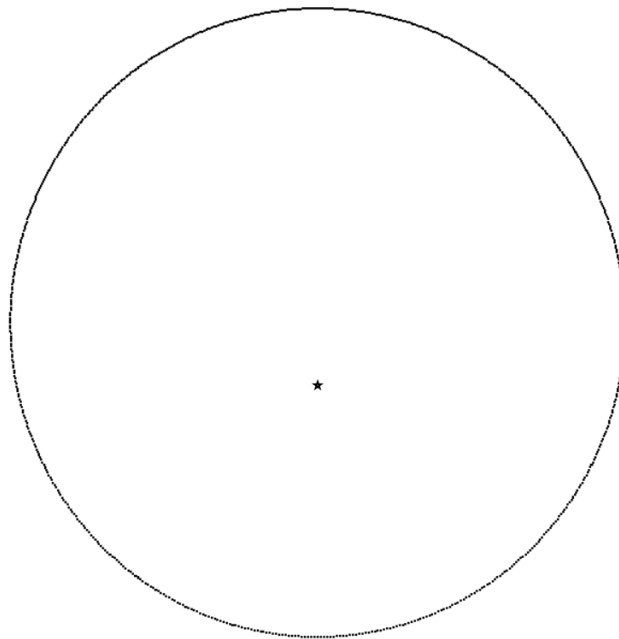

Figure 2: Mercury in orbit around the Sun.