Newtonian gravitation for C++ programmers

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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 dt = 10000; // 2.77777 hours

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;
```

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```
const real_type c4 = c * c * c * c;
const real_type h = 6.62607015e-34;
const real_type hbar = h / (2.0 * pi);
```

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:

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

```
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;
}
```

```
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";
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. Thus, 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:

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

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

```
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
// 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

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:

```
vector_3 Newtonian_acceleration(
        const real_type emitter_mass,
        const vector_3& pos, // Receiver pos
        const real_type G)
{
        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:

```
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] Fiedler. Fix Your Timestep! (2004)

[2] Fiedler. Integration Basics. (2004)

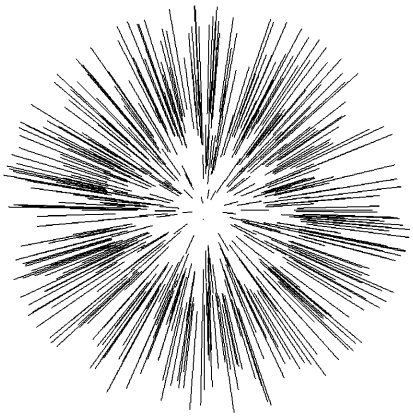

Figure 1: Example of an isotropic emitter. The emitter is spherical. The field line starting locations are placed pseudorandomly on a 2-sphere, and the normals (e.g. field line directions) are calculated using the same sphere.

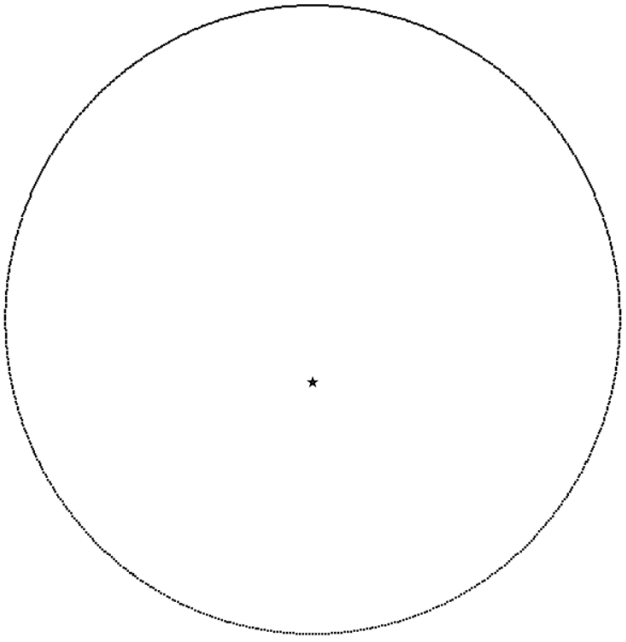

Figure 2: Mercury in orbit around the Sun.