# 可视计算与交互概论-Lab4

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2024/11/23

### 1 Task 1: Inverse Kinematics

#### 1.1 Sub-Task 1:forward kinematic

思路:从 startindex 开始更新 ik 上的关节位置信息,考虑 i 关节,其父关节为 i-1,则有以下公式

```
global\_rotate[i] = global\_rotate[i-1] * local\_rotate[i]
```

 $global\_pos[i] = global\_pos[i-1] + global\_offset[i-1,i] = global\_rotate[i-1] * local\_offset[i-1,i] = global\_rotate[i-1] * local\_offset[i-1] * local\_offset[i-$ 

即 i 关节的全局旋转等于 i-1 关节的全局旋转叠加上 i 关节相对 i-1 关节的局部旋转; i 关节的全局位置等于 i-1 关节的全局位置加上 i 关节相对 i-1 关节的偏移 (考虑旋转)。

```
void ForwardKinematics(IKSystem & ik , int StartIndex) {
      if (StartIndex == 0)  {
          ik.JointGlobalRotation[0] = ik.JointLocalRotation[0];
          ik.JointGlobalPosition[0] = ik.JointLocalOffset[0];
           StartIndex
                                      = 1;
      }
      for (int i = StartIndex; i < ik.JointLocalOffset.size(); i++) {</pre>
          // your code here: forward kinematics, update JointGlobalPosition and
      Joint Global Rotation
          ik. Joint Global Rotation [i] = glm::normalize(glm::normalize(ik. Joint Global Rotation [i -
10
      1]) *glm::normalize(ik.JointLocalRotation[i]));
          glm:: vec4 \quad global\_offset = ik.JointGlobalRotation[i - 1] * glm:: vec4(ik.
11
      JointLocalOffset[i], 1.0);
          ik.JointGlobalPosition[i] = ik.JointGlobalPosition[i - 1] + glm::vec3(global offset);
      }
```

#### 1.2 Sub-Task 2:CCD IK

思路:对于 CCD IK,想法是从尾部节点开始,每次旋转都使得自己、末端关节节点、目标位置三者位于同一条直线上,这样使得每一次旋转后末端关节节点都处在与目标位置最近的位置 (对于目前正在处理的关节节点的所有可能局部旋转而言)。

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具体实现:对每一个关节节点,分别求其到 EndPosition 与 EndEffectorPosition 的向量,然后标准化得到 cur\_to\_tar 与 cur\_to\_effect,然后使用 glm::rotation 函数计算从 cur\_to\_tar 旋转至 cur\_to\_effect 的旋转四元数,并更新相应节点的局部旋转值,最后从这个节点开始对其子节点进行位姿更新,通过不断的迭代实现末端关节节点逼近目标位置。

```
void InverseKinematicsCCD (IKSystem & ik, const glm::vec3 & EndPosition, int maxCCDIKIteration,
      float eps) {
      ForwardKinematics (ik, 0);
      // These functions will be useful: glm::normalize, glm::rotation, glm::quat * glm::quat
      for (int CCDIKIteration = 0; CCDIKIteration < maxCCDIKIteration && glm::12Norm(ik.
     EndEffectorPosition() - EndPosition) > eps; CCDIKIteration++) {
          // your code here: ccd ik
          for (int i = ik.NumJoints() - 1; i >= 0; i--)
              glm::vec3 cur_joint_pos = ik.JointGlobalPosition[i];
              glm::vec3 end_effector_pos = ik.EndEffectorPosition();
10
              glm::vec3 cur_to_tar = glm::normalize(EndPosition - cur_joint_pos);
11
              glm::vec3 cur_to_effect = glm::normalize(end_effector_pos - cur_joint_pos);
12
13
14
              glm::quat rotation_quat = glm::rotation(cur_to_effect,cur_to_tar);
15
              ik.JointLocalRotation[i] = rotation_quat * ik.JointLocalRotation[i];
16
              ForwardKinematics(ik, i);
18
19
20
          }
      }
22
23 }
```

得到效果图如下:

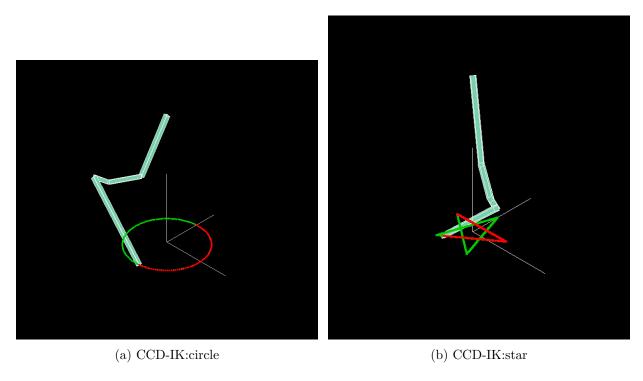


图 1: Sub-Task 2:CCD IK-results

### 1.3 Sub-Task 3:FABR IK

思路:与 CCD IK 不同,FABR IK 不通过计算旋转角度然后调整位姿,而是直接调整骨骼方向,改变关节位置。整个过程经历了两次更新,在 backward update 阶段,先假设能够末端关节位置位于目标位置,然后在此基础上根据 cur\_joint 的局部 offset 和 next\_position 到 cur\_joint 的方向更新 cur\_joint 的位置,直至更新得到一个新的根节点位置 p0';然后进行 forward update,保持 p0 位置不变,利用 p0',p1'等 backward update 得到的新关节位置计算得到新的关节位置 p0",p1"..... 经过这样两次 update,可以使得末端关节位置不断逼近目标位置。

```
void InverseKinematicsFABR(IKSystem & ik, const glm::vec3 & EndPosition, int
     maxFABRIKIteration, float eps) {
      ForwardKinematics (ik, 0);
      int nJoints = ik.NumJoints();
      std::vector<glm::vec3> backward_positions(nJoints, glm::vec3(0, 0, 0)), forward_positions(
     nJoints, glm::vec3(0, 0, 0));
      for (int IKIteration = 0; IKIteration < maxFABRIKIteration && glm::l2Norm(ik.
     EndEffectorPosition() - EndPosition) > eps; IKIteration++) {
          // task: fabr ik
          // backward update
          glm::vec3 next position
                                           = EndPosition;
          backward\_positions[nJoints - 1] = EndPosition;
10
          for (int i = nJoints - 2; i >= 0; i--) {
11
              // your code here
              glm::vec3 normalized_dir = glm::normalize(ik.JointGlobalPosition[i] -
13
     next_position);
              backward positions[i] = next position + normalized dir * glm::length(ik.
14
     JointLocalOffset[i + 1]);
              {\tt next\_position}
                                        = backward_positions[i];
          }
```

```
// forward update
    glm::vec3 now_position = ik.JointGlobalPosition[0];
     forward positions [0] = ik. Joint Global Position [0];
     for (int i = 0; i < nJoints - 1; i++) {
        // your code here
        glm::vec3 normalized_dir = glm::normalize(backward_positions[i + 1] - now_position
);
         forward_positions[i + 1] = now_position + normalized_dir * glm::length(ik.
JointLocalOffset[i + 1]);
                                  = forward positions[i + 1];
         now\_position
     ik.JointGlobalPosition = forward_positions; // copy forward positions to
joint_positions
}
// Compute joint rotation by position here.
for (int i = 0; i < nJoints - 1; i++) {
     ik. Joint Global Rotation [i] = glm::rotation (glm::normalize(ik. Joint Local Offset [i + 1]),
glm::normalize(ik.JointGlobalPosition[i + 1] - ik.JointGlobalPosition[i]));
ik.JointLocalRotation[0] = ik.JointGlobalRotation[0];
for (int i = 1; i < nJoints - 1; i++) {
     ik.JointLocalRotation[i] = glm::inverse(ik.JointGlobalRotation[i - 1]) * ik.
JointGlobalRotation[i];
ForwardKinematics (ik, 0);
```

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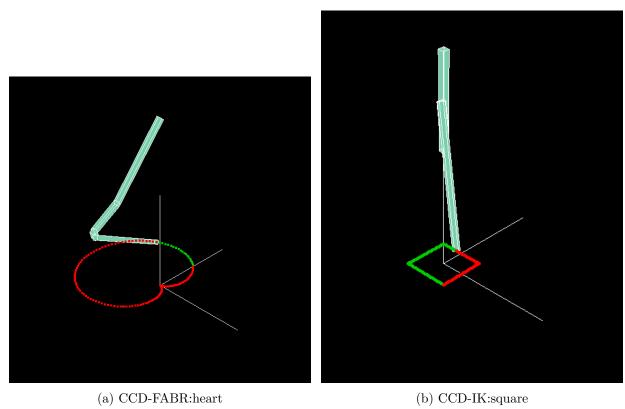


图 2: Sub-Task 3:FABR IK-results

#### 1.4 Sub-Task 4:Custom Draw

思路:类似参数方程,画直线使用一次函数,画曲线可以使用圆/椭圆的参数方程。最终得到 custom str:XTRY。

```
IKSystem::Vec3ArrPtr IKSystem::BuildCustomTargetPosition() {
       int nums
                       = 50;
       using Vec3Arr = std::vector<glm::vec3>;
       std::shared_ptr<Vec3Arr> custom(new Vec3Arr(nums*12));
       int
                                    index = 0;
6
       // X
       for (int i = 0; i < nums; i++) {
                            = 0.4 * i / nums;
            float x_val
9
            float y_val
                                 = 0.4 * i / nums;
10
            (*custom)[index++] = glm :: vec3(x_val+1.35, 0.0f, y_val);
11
12
       for (int i = 0; i < nums; i++) {
13
                            = 0.4 * i / nums;
            float x_val
14
            float y_val
                                = 0.4 - 0.4 * i / nums;
            (*custom)[index++] = glm :: vec3(x_val+1.35, 0.0f, y_val);
16
       }
17
18
       // T
19
       for (int i = 0; i < nums; i++) {
20
                               = 0.4 * i / nums;
            float x_val
21
22
            float y_val
            (*{\rm custom}\,)\,[\,{\rm index} + +] \,=\, {\rm glm}:: {\rm vec3}\,(\,{\rm x\_val} + 0.85\,,\ 0.0\,{\rm f}\,\,,\ {\rm y\_val}\,)\,;
23
       }
24
```

```
for (int i = 0; i < nums; i++) {
    float x_val
                        = 0.4 * i / nums;
    float y_val
    (*custom)[index++] = glm :: vec3(x val+0.85, 0.0f, y val);
}
    //R
    for (int i = 0; i < nums; i++) {
    float x_val
                        = 0.2;
    float y_val
                        = 0.4 * i / nums;
    (*custom)[index++] = glm :: vec3(x_val+0.5, 0.0f, y_val);
}
    for (int i = 0; i < nums; i++) {
    float x_val
                        = 0.2 - 0.2 * \sin(EIGEN_PI * i / nums);
                            = 0.3+0.1 * \cos(EIGEN_PI* i / nums);
        float y val
        (*custom)[index++] = glm :: vec3(x_val+0.5, 0.0f, y_val);
    for (int i = 0; i < nums; i++) {
        float x_val
                            = 0.2 - 0.2 * i / nums;
                            = 0.2 - 0.2 * i / nums;
        float y val
        (*custom)[index++] = glm :: vec3(x_val+0.5, 0.0f, y_val);
    }
// Y
    for (int i = 0; i < nums; i++) {
        float x_val
                            = 0.15;
                            = 0.25 * i / nums;
        float y_val
        (*custom)[index++] = glm :: vec3(x_val+0.05, 0.0f, y_val);
    }
    for (int i = 0; i < nums; i++) {
        float x_val
                            = 0.15 + 0.15 * i / nums;
        float y val
                            = 0.25 + 0.15 * i / nums;
        (*custom)[index++] = glm :: vec3(x_val+0.05, 0.0f, y_val);
    }
    for (int i = 0; i < nums; i++) {
        float x_val
                            = 0.15 - 0.15 * i / nums;
        float y val
                            = 0.25 + 0.15 * i / nums;
        (*custom)[index++] = glm :: vec3(x_val+0.05, 0.0f, y_val);
    }
custom->resize (index);
return custom;
```

得到效果图如下:

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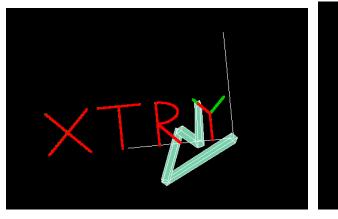
58

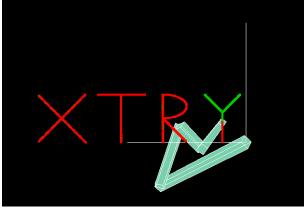
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63 64 }





(a) Cutom Draw-XTRY

(b) Cutom Draw-XTRY

图 3: Sub-Task 4:Cutom Draw-XTRY

彩蛋-真·doge, 在 wolframalpha 网站上发现了 doge-curve, 修改了其中的参数格式, 得到 doge:D

```
float doge_custom_x(float t) {
             float x_t = ***/too long-hidden
             return x_t + 550.0;
        float doge_custom_y(float t) {
             float y_t = ***/too long->hidden
             return y_t + 550.0;
       IKSystem::Vec3ArrPtr IKSystem::BuildCustomTargetPosition() {
10
       int nums
                         = 5000;
11
        using Vec3Arr = std::vector<glm::vec3>;
12
       std::shared_ptr<Vec3Arr> custom(new Vec3Arr(nums));
13
                                        index = 0;
14
       for (int i = 0; i < nums; i++) {
15
             \mathbf{float} \  \, \mathbf{x\_val} = 1.5\,\mathrm{e}{-3}\mathrm{f} \  \, * \  \, \mathrm{doge\_custom\_x}(92 \  \, * \  \, \mathrm{glm}::\mathrm{pi}{<}\mathbf{float}{>}() \  \, * \  \, \mathrm{i} \  \, / \  \, \mathrm{nums});
             float y_val = 1.5e-3f * doge_custom_y(92 * glm::pi < float > () * i / nums);
17
             if (std::abs(x_val) < 1e-3 || std::abs(y_val) < 1e-3) continue;
18
             (*custom)[index++] = glm :: vec3(1.6 f - x_val, 0.0 f, y_val - 0.2 f);
19
20
       custom->resize(index);
21
       return custom;
22
23
24
```

得到效果图如下:

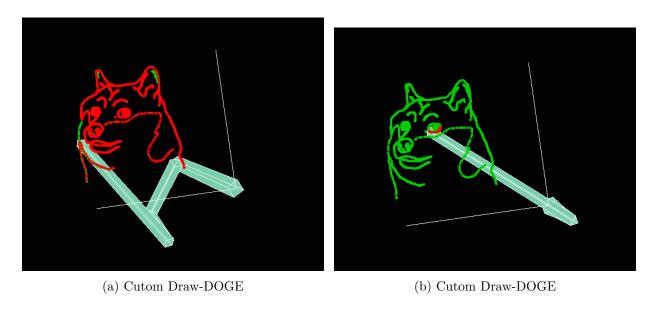


图 4: Sub-Task 4:Cutom Draw-DOGE

## 1.5 Sub-Task 4-1:Cutom Draw sample

思路:采样点不均匀的原因可能是在头发处坐标变化比较剧烈,如果采样密度不够高,就会导致头发区域的采样点比较稀疏,一个可行的方法是增加采样密度,下图分别是总采样 num=5000 与总采样 num=50000 的效果图区别:

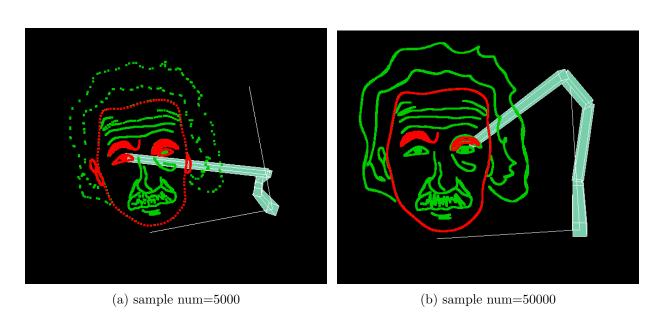


图 5: Sub-Task 4-1:Cutom Draw-sample

但这只是增加了总体的采样数目,会导致计算开销较大,一个可行的方案是根据图像的剧烈变化程度动态调整采样数目,在图像剧烈变化的地方 (可以用梯度图分析变化程度) 提高采样密度, 这样就可以使得采样点更加均匀。

## 1.6 Report

# 1.6.1 1. 如果目标位置太远,无法到达, IK 结果会怎样?

会把"臂"伸直指向目标位置,然后在 origin\_pos->tar\_pos 方向上的最远处作画. 其中 CCD-IK 会抽搐,留下无规律的散点,而 FABR-IK 会在最远处画出对应图像

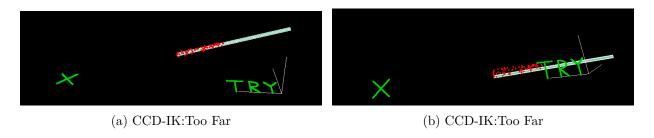


图 6: Task 1-Question 1-CCD-IK

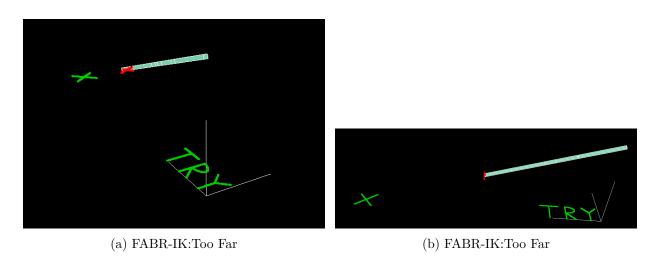


图 7: Task 1-Question 1-FABR-IK

### 1.6.2 2. 比较 CCD IK 和 FABR IK 所需要的迭代次数。

FABR IK 的迭代次数比 CCD IK 的迭代次数更少。

如下图是画 circle 时 CCD-IK 与 FABR-IK 的迭代输出结果,可以看出 CCD-IK 基本上到 max\_iteration=100 都还没有收敛,而 FABR-IK 只需要迭代个位数次就能早早完成收敛。

```
IK-CCD cur-iteration:63, Max_iteration_limit:100
IK-CCD cur-iteration:64, Max_iteration_limit:100
IK-CCD cur-iteration:65, Max_iteration_limit:100
IK-CCD cur-iteration:66, Max_iteration_limit:100
IK-CCD cur-iteration:67, Max_iteration_limit:100
IK-CCD cur-iteration:69, Max_iteration_limit:100
IK-CCD cur-iteration:70, Max_iteration_limit:100
IK-CCD cur-iteration:71, Max_iteration_limit:100
IK-CCD cur-iteration:72, Max_iteration_limit:100
IK-CCD cur-iteration:73, Max_iteration_limit:100
IK-CCD cur-iteration:74, Max_iteration_limit:100
IK-CCD cur-iteration:75, Max_iteration_limit:100
IK-CCD cur-iteration:75, Max_iteration_limit:100
IK-CCD cur-iteration:76, Max_iteration_limit:100
IK-CCD cur-iteration:78, Max_iteration_limit:100
IK-CCD cur-iteration:80, Max_iteration_limit:100
IK-CCD cur-iteration:81, Max_iteration_limit:100
IK-CCD cur-iteration:82, Max_iteration_limit:100
IK-CCD cur-iteration:83, Max_iteration_limit:100
IK-CCD cur-iteration:83, Max_iteration_limit:100
IK-CCD cur-iteration:84, Max_iteration_limit:100
IK-CCD cur-iteration:85, Max_iteration_limit:100
IK-CCD cur-iteration:89, Max_iteration_limit:100
IK-CCD cur-iteration:99, Max_iteratio
```

图 8: Task 1-Question 2-iterations time results

### 6.3 3. IK 多解导致前后两帧关节旋转抖动的情况

可能的解决方案: 限制旋转角度, 规定 IK 旋转只能在一个规定的小范围内进行。

# 2 Task 2: Mass-Spring System

整体思路:

对于 Implicit Euler:

$$x_{n+1} = x_n + dt \cdot v_{n+1}$$
$$v_{n+1} = v_n + dt \cdot M^{-1} f(x_{n+1})$$

经过分析计算,我们得到  $x_{n+1}=(x_n+dt(v_n+dt\cdot M^{-1}f_{ext}))+dt^2M^{-1}f_{int}(x_{n+1})$  求解这个方程相当于求解  $x_{n+1}=argmin_xg(x), g(x)=\frac{1}{2dt^2}|x-y|_M^2+E(x)$  其中  $|x-y|_M^2=(x-y)^TM(x-y), y=x_n+dt(v_n+dt\cdot M^{-1}f_{ext})$  进一步,我们对 g(x) 在当前尝试解  $x_i$  处左二阶泰勒展开,得到

$$g(x) = g(x_i) + \nabla g(x_i) \cdot (x - x_i) + \frac{1}{2}(x - x_i)^T H_g(x_i)(x - x_i) + O(||x - x_i||^3)$$

其中, $\nabla g(x_i)$  表示  $g(x_i)$  的梯度, $H_g(x_i)$  表示 g(x) 在  $x_i$  处的海瑟矩阵。然后我们忽略三阶小量,对式子两边求梯度,再代入  $x_{i+1}$ , 得到

$$\nabla q(x_{i+1}) = \nabla q(x_i) + H_q(x_i)(x_{i+1} - x_i)$$

令这条等于 0(求极值点), 得到  $x_{i+1} = x_i - H_q^{-1}(x_i) \nabla g(x_i)$ 

$$\nabla g(x_i) = \frac{1}{dt^2} M(x_i - y_i) + \nabla E(x_i)$$
$$H_g(x_i) = \frac{1}{dt^2} M + H(x_k)$$

其中  $H(x_k)$  是弹性势能 E(x) 的海瑟矩阵,有  $H(x_k) = \Sigma_{(i,j)} H_{ij}(x_k)$  在完成  $\nabla g(x_i)$  与  $H_g(x_i)$  的求解后,我们得到  $x_{i+1} - x_i = -H_g^{-1}(x_i) \nabla g(x_i)$  然后更新  $x_{i+1}$  并以此为基础更新  $v_{i+1}$  具体实现:

首先初始化 x0,v0,f0(外力矩阵, 此处为简单的重力加速度 mg/m), 然后根据公式计算 y0

利用函数 eval\_grad\_g 计算出 grad\_g, 主要围绕公式  $\nabla g(x_i) = \frac{1}{dt^2} M(x_i - y_i) + \nabla E(x_i)$ , 其中  $\nabla E(x_i)$  总能量关于所有质点的位置的梯度就是把关于每个质点的梯度拼接起来,而能量关于质点的梯度相当于质点 受到的力 (取负号表示方向相反)。

```
Eigen:: VectorXf eval_intern_f(MassSpringSystem & system, Eigen:: VectorXf & x_vec)
      std::vector<glm::vec3> forces(system.Positions.size(), glm::vec3(0));
      std :: vector < glm :: vec3 > x_vec3 = eigen2glm(x_vec);
      for (auto const spring : system.Springs) {
          auto const
                           p0 = spring.AdjIdx.first;
                           p1 = spring.AdjIdx.second;
          auto const
          if (system.Fixed[p0] && system.Fixed[p1]) continue;
          glm :: vec3 const x01 = x_vec3[p1] - x_vec3[p0]; // use updated x_vec
          glm :: vec3 const e01 = glm :: normalize(x01);
                               = (system. Stiffness * (glm::length(x01) - spring.RestLength)) *
          glm :: vec3
     e01;
          forces[p0] += f;
          forces [p1] -= f;
      return glm2eigen (forces);
16
18 Eigen:: VectorXf eval_grad_g (MassSpringSystem & system, Eigen:: VectorXf & x_vec, Eigen::
     VectorXf & y_vec, float ddt)
19 {
      Eigen:: VectorXf grad_g_ret = system.Mass * (x_vec - y_vec) / (ddt*ddt);
20
```

```
grad_g_ret -= eval_intern_f(system, x_vec);
return grad_g_ret;
}
```

进一步,我们计算  $H_g(x_i)$ , 主要围绕公式  $H_g(x_i) = \frac{1}{dt^2}M + H(x_k)$  前面带 M 的部分好处理,直接给 SparseMatrix 对角线赋值即可

后面的 Hessian 矩阵部分利用公式

$$H_{e-ij} = k_{ij} \frac{(x_i - x_j)(x_i - x_j)^T}{||x_i - x_j||^2} + k_{ij} \left(1 - \frac{l_{ij}}{||x_i - x_j||}\right) \left(I - \frac{(x_i - x_j)(x_i - x_j^T)}{||x_i - x_j||^2}\right)$$

得到一个 3\*3 的  $H_{e-ij}$  小矩阵,然后进一步计算对应的  $H_{ij}(x_k)$  大矩阵,其 shape 为 (n\*n)\*(3\*3), 即 n\*n 个 3\*3 的块,其中  $H_{ij}(i,j) = H_{ij}(j,i) = -H_{e-ij}, H_{ij}(i,i) = H_{ij}(j,j) = H_{e-ij}$ 

```
// Hessian Matrix
2 std::vector<Eigen::Triplet<float>>> Hessian_triplets;
3 std::vector<glm::vec3>
                                        x \text{ vec3} = eigen2glm(x \text{ vec});
  for (auto const spring : system.Springs) {
      auto const p0 = spring.AdjIdx.first;
      auto const p1 = spring.AdjIdx.second;
      glm :: vec3 \ const \ x01 = x_vec3[p1] - x_vec3[p0];
      float len_x01 = glm :: length(x01);
      glm::mat3 identity matrix {1.0 f};
11
12
      glm :: mat3 \times matrix(0.0 f);
13
      for (int i = 0; i < 3; i++)
           for (int j = 0; j < 3; j++)
17
18
               x_{matrix}[i][j] = x01[i] * x01[j];
19
      }
21
22
      glm::mat3 hessian_tmp_matrix = x_matrix / (len_x01*len_x01);
      hessian_tmp_matrix += (1.0 f - spring.RestLength / len_x01) * (identity_matrix - x_matrix/(
24
      len_x01*len_x01));
      hessian_tmp_matrix *= system.Stiffness;
25
```

```
for (int i = 0; i < 3; i++)
27
28
          for (int j = 0; j < 3; j++)
29
30
               Hessian_triplets.emplace_back(p0 * 3 + i, p0 * 3 + j, hessian_tmp_matrix[i][j]);
31
     //Warning: from size_t to
               Hessian_triplets.emplace_back(p1 * 3 + i, p1 * 3 + j, hessian_tmp_matrix[i][j]);
32
               Hessian_triplets.emplace_back(p0 * 3 + i, p1 * 3 + j, -hessian_tmp_matrix[i][j]);
33
               Hessian_triplets.emplace_back(p1 * 3 + i, p0 * 3 + j, -hessian_tmp_matrix[i][j]);
34
          }
36
37
38
40 Eigen::SparseMatrix<float> Hessian_mat = CreateEigenSparseMatrix(3 * pos_num, Hessian_triplets
     );
```

最后把得到的 m\_mat 与 Hessian\_mat 相加即得到待求的  $H_g(x_i)$  完整的 eval\_grad2\_g 函数如下:

```
void eval_grad2_g(MassSpringSystem & system, Eigen:: VectorXf & x_vec, float ddt, Eigen::
      SparseMatrix < float > & dst_matrix)
2 {
      int pos_num = system.Positions.size();
      Eigen::SparseMatrix<float> grad2_g_ret(3 * pos_num, 3 * pos_num);
      // M Matrix
      std::vector<Eigen::Triplet<float>> m_triplets;
                                           tmp m = system.Mass / (ddt * ddt);
      for (int i = 0; i < 3*pos_num; i++)
          m_triplets.emplace_back(i, i,tmp_m);
12
13
      Eigen::SparseMatrix<float> m_mat = CreateEigenSparseMatrix(size_t(3) * pos_num, m_triplets)
14
15
      // Hessian Matrix
16
      std::vector<Eigen::Triplet<float>>> Hessian_triplets;
17
      std::vector<glm::vec3>
                                           x_{vec3} = eigen2glm(x_{vec});
      for (auto const spring : system.Springs) {
19
          auto const p0 = spring.AdjIdx.first;
20
          auto const p1 = spring.AdjIdx.second;
21
           if (system.Fixed[p0] && system.Fixed[p1]) continue;
22
          glm :: vec3 \ const \ x01 = x_vec3[p1] - x_vec3[p0];
23
24
          float len_x01 = glm :: length(x01);
25
26
          glm::mat3 identity_matrix {1.0f};
```

```
//printf(\%f \%f \%f n\%, identity\_matrix[0][0], identity\_matrix[1][1], identity\_matrix
                            [2][2]);
                                                 glm::mat3 x matrix(0.0f);
                                                  for (int i = 0; i < 3; i++)
                                                                      for (int j = 0; j < 3; j++)
                                                                                        x_{matrix}[i][j] = x01[i] * x01[j];
                                                  }
                                                 glm::mat3 hessian_tmp_matrix = x_matrix / (len_x01*len_x01);
                                                 hessian_tmp_matrix += (1.0f - spring.RestLength / len_x01) * (identity_matrix -
                          x = \frac{x}{1} \cdot \frac{1}{1} \cdot 
                                                  hessian_tmp_matrix *= system.Stiffness;
                                                  for (int i = 0; i < 3; i++)
                                                                     for (int j = 0; j < 3; j++)
                                                                                         Hessian_triplets.emplace_back(p0 * 3 + i, p0 * 3 + j, hessian_tmp_matrix[i][j
                           ); //Warning: from size_t to
                                                                                         Hessian_triplets.emplace_back(p1 * 3 + i, p1 * 3 + j, hessian_tmp_matrix[i][j
                           ]);
                                                                                         Hessian\_triplets.emplace\_back(p0 * 3 + i, p1 * 3 + j, -hessian\_tmp\_matrix[i][j]
                           ]);
                                                                                         Hessian_triplets.emplace_back(p1 * 3 + i, p0 * 3 + j, -hessian_tmp_matrix[i][j
                           ]);
                                                                      }
                              }
                              Eigen::SparseMatrix<float> Hessian_mat = CreateEigenSparseMatrix(3 * pos_num,
                           Hessian_triplets);
                              dst_matrix = m_mat+Hessian_mat;
60 }
```

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然后我们利用得到的 grad2\_g 与 grad\_g 列出方程  $H_q(x_i)(x_{i+1}-x_i) = - \nabla g(x_i)$  调用 ComputeSimplicialLLT 计算得  $\delta x$ , 然后更新  $x_{i+1}$  和  $v_{i+1}$  即可。

```
Eigen::VectorXf grad_g = eval_grad_g(system, x0, y0, ddt);
Eigen::SparseMatrix<float> grad2_g(3 * pos_num, 3 * pos_num);
eval_grad2_g(system, x0, ddt, grad2_g);
Eigen:: VectorXf x1_x0 = ComputeSimplicialLLT(grad2_g, -grad_g);
```

```
Eigen::VectorXf final_x = x1_x0 + x0;

std::vector<glm::vec3> final_x_vec3 = eigen2glm(final_x);

std::vector<glm::vec3> final_v = eigen2glm(v0 + ddt * (eval_intern_f(system, final_x)/system.

Mass + f0_extern_grav));

for (std::size_t i = 0; i < system.Positions.size(); i++) {

    if (system.Fixed[i]) continue;

    system.Velocities[i] = final_v[i];

    system.Positions[i] = final_x_vec3[i];

}
```

### 完整代码:

```
\textbf{float} \  \, \text{eval\_g} \, (\, \text{MassSpringSystem} \, \, \& \, \, \text{system} \, \, , \  \, \text{Eigen} :: \, \text{VectorXf} \, \, \& \, \, \text{x\_vec} \, , \, \, \, \text{Eigen} :: \, \text{VectorXf} \, \, \& \, \, \text{y\_vec} \, , \, \\ \textbf{float} \, \, \text{float} \, \, \, \text{float} \, \, \text{float}
                      ddt)
 2 {
                    Eigen::VectorXf tmp_vec = x_vec-y_vec;
                                                                           eval\_g\_ret = ((tmp\_vec * system.Mass).dot(tmp\_vec)) / (2.0 f*ddt * ddt);
                     for (auto const spring : system.Springs)
                                  auto const p0 = spring.AdjIdx.first;
                                  auto const p1 = spring.AdjIdx.second;
                                   if (system.Fixed[p0] && system.Fixed[p1]) continue;
                                  glm::vec3 const x01 = system.Positions[p1] - system.Positions[p0];
                                   float
                                                                                         deta length = glm::length(x01) - spring.RestLength;
                                   float
                                                                                        tmp_E = 1.0 f / 2.0 f * system. Stiffness * deta_length*deta_length;
14
                                  eval\_g\_ret += tmp\_E;
16
                    return eval_g_ret;
17
18 }
19
20 Eigen:: VectorXf eval_intern_f(MassSpringSystem & system, Eigen:: VectorXf & x_vec)
       {
21
                    std::vector<glm::vec3> forces(system.Positions.size(), glm::vec3(0));
22
                     std :: vector < glm :: vec3 > x_vec3 = eigen2glm(x_vec);
23
                     for (auto const spring : system.Springs) {
                                                                                        p0 = spring.AdjIdx.first;
                                  auto const
                                                                                        p1 = spring.AdjIdx.second;
                                  auto const
26
                                   if (system.Fixed[p0] && system.Fixed[p1]) continue;
27
                                  glm :: vec3  const  x01 = x_vec3[p1] - x_vec3[p0]; // use updated <math>x_vec
28
29
                                  glm:: vec3 const e01 = glm:: normalize(x01);
30
                                  glm :: vec3
                                                                                               = (system.Stiffness * (glm::length(x01) - spring.RestLength)) *
31
                   e01;
                                   forces[p0] += f;
32
                                   forces[p1] = f;
33
34
                    return glm2eigen(forces);
35
```

```
37 Eigen:: VectorXf eval_grad_g (MassSpringSystem & system, Eigen:: VectorXf & x_vec, Eigen::
      VectorXf & y_vec, float ddt)
38 {
      Eigen:: VectorXf grad_g_ret = system.Mass * (x_vec - y_vec) / (ddt*ddt);
39
      grad_g_ret = eval_intern_f(system,x_vec);
40
      return grad_g_ret;
41
42 }
43
44 void eval_grad2_g(MassSpringSystem & system, Eigen:: VectorXf & x_vec, float ddt, Eigen::
      SparseMatrix < float > & dst_matrix)
  {
45
      int pos_num = system.Positions.size();
46
47
      Eigen::SparseMatrix<float> grad2_g_ret(3 * pos_num, 3 * pos_num);
48
49
      // M Matrix
50
      std::vector<Eigen::Triplet<float>> m_triplets;
51
                                            tmp \ m = system.Mass / (ddt * ddt);
      for (int i = 0; i < 3*pos_num; i++)
53
54
           m_triplets.emplace_back(i, i,tmp_m);
55
56
      Eigen::SparseMatrix<float> m_mat = CreateEigenSparseMatrix(size_t(3) * pos_num, m_triplets)
57
58
      // Hessian Matrix
59
      std::vector<Eigen::Triplet<float>>> Hessian_triplets;
60
      std::vector<glm::vec3>
                                            x_{vec3} = eigen2glm(x_{vec});
61
      for (auto const spring : system. Springs) {
62
           auto const p0 = spring.AdjIdx.first;
63
           auto const p1 = spring.AdjIdx.second;
64
           if (system.Fixed[p0] && system.Fixed[p1]) continue;
65
           glm :: vec3 \ const \ x01 = x_vec3[p1] - x_vec3[p0];
66
           float len_x01 = glm :: length(x01);
68
69
           glm::mat3 identity_matrix {1.0f};
70
           //\text{printf}(\%f \%f \%f \%n), identity_matrix [0][0], identity_matrix [1][1], identity_matrix
71
      [2][2]);
72
           glm::mat3 x_matrix(0.0 f);
73
74
           for (int i = 0; i < 3; i++)
76
               for (int j = 0; j < 3; j++)
77
78
                   x_{matrix}[i][j] = x01[i] * x01[j];
79
80
           }
82
```

```
glm::mat3 hessian_tmp_matrix = x_matrix / (len_x01*len_x01);
                  hessian_tmp_matrix += (1.0f - spring.RestLength / len_x01) * (identity_matrix -
       x_{matrix}/(len_x01*len_x01);
                  hessian tmp matrix *= system. Stiffness;
                  for (int i = 0; i < 3; i++)
                  {
                           for (int j = 0; j < 3; j++)
                                    Hessian_triplets.emplace_back(p0 * 3 + i, p0 * 3 + j, hessian_tmp_matrix[i][j
        ); //Warning: from size_t to
                                    Hessian_triplets.emplace_back(p1 * 3 + i, p1 * 3 + j, hessian_tmp_matrix[i][j
       ]);
                                    Hessian\_triplets.emplace\_back(p0\ *\ 3\ +\ i\ ,\ p1\ *\ 3\ +\ j\ ,\ -hessian\_tmp\_matrix[\ i\ ][\ j\ ]
        ]);
                                    Hessian_triplets.emplace_back(p1 * 3 + i, p0 * 3 + j, -hessian_tmp_matrix[i][j
       ]);
                           }
                  }
         }
         Eigen::SparseMatrix<float> Hessian mat = CreateEigenSparseMatrix(3 * pos num,
        Hessian_triplets);
         dst_matrix = m_mat+Hessian_mat;
void AdvanceMassSpringSystem (MassSpringSystem & system, float const dt) {
         // Implicit Euler algorithm
         // x_{k+1}=x_{k}+ddt*v_{k+1}
         // v_{k+1}=v_{k}+ddt*F(x_{k+1})/M
                                    steps = 1;
         int const
         float const ddt = dt / steps;
         // without extern iterations
         for (std::size\_t \ s = 0; \ s < steps; \ s++) {
                  int pos_num = system.Positions.size();
                  // initialize
                                                                  = glm2eigen(system.Positions);
                  Eigen::VectorXf x0
                  Eigen:: VectorXf v0 = glm2eigen(system. Velocities);
                  Eigen::VectorXf f0_extern_grav = glm2eigen(std::vector<glm::vec3>(system.Positions.
        size(), glm::vec3(0, -system.Gravity, 0));
                  Eigen::VectorXf system_m_vectorxf = glm2eigen(std::vector<glm::vec3>(system.Positions.
        size(), glm::vec3(system.Mass,system.Mass,system.Mass)));
                  Eigen:: VectorXf y0 = x0 + ddt * (v0 + ddt * f0_extern_grav);
                  Eigen::VectorXf grad_g = eval_grad_g(system, x0, y0, ddt);
                  \label{eq:continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous_continuous
                  eval_grad2_g(system, x0, ddt, grad2_g);
```

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```
Eigen::VectorXf x1_x0 = ComputeSimplicialLLT(grad2_g, -grad_g);

Eigen::VectorXf final_x = x1_x0 + x0;
std::vector<glm::vec3> final_x_vec3 = eigen2glm(final_x);
std::vector<glm::vec3> final_v = eigen2glm(v0 + ddt * (eval_intern_f(system, final_x)/system.Mass + f0_extern_grav));

for (std::size_t i = 0; i < system.Positions.size(); i++) {
    if (system.Fixed[i]) continue;
    system.Velocities[i] = final_v[i];
    system.Positions[i] = final_x_vec3[i];
}
</pre>
```

### 得到效果图:

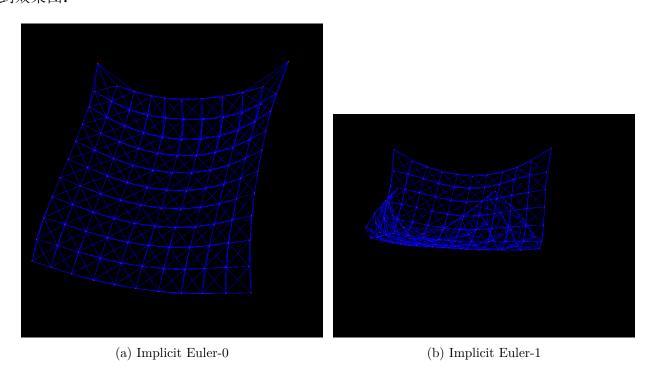


图 9: Task 2: Mass-Spring System-Implicit Euler

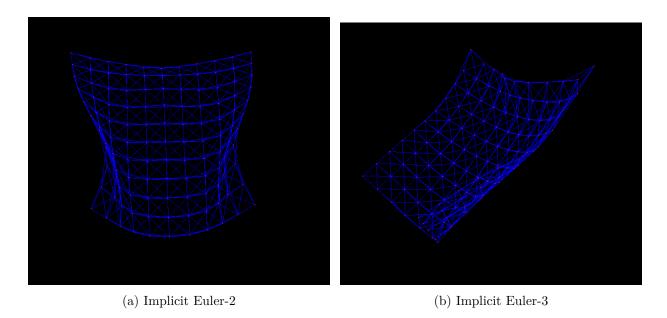


图 10: Task 2: Mass-Spring System-Implicit Euler