

VIO第三章作业分享





纲要



▶1. 使用LM算法估计曲线参数

- ▶1.1 绘制阻尼因子µ随着迭代变化的曲线图
- ▶1.2 将曲线函数改成 $y = ax^2 + bx + c$
- ▶1.3 实现更优秀的阻尼因子策略
- ▶2. 公式推导雅可比F, G中的两项
- ▶3. 证明式(9)



1.1 绘制阻尼因子µ随着迭代变化的曲线图

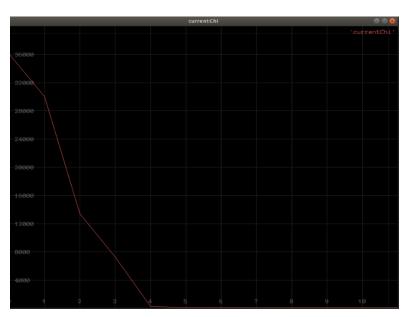
这里使用Pangolin来绘制 损失函数currentChi 和 阻尼因子currentLambda 随着迭代步骤iter变化的曲线图。 我们在backend/problem.cc中 Problem类内写入draw_curve函数

```
void Problem::draw_curve(const vector<double, std::allocator<double>> &currentValue_vec, string name){
   pangolin::DataLog log;
    std::vector<std::string> labels;
   log.SetLabels(labels):
   pangolin::Plotter plotter( default_log: &log, left: 0, right: currentValue_vec.size() - 1, bottom: 0, top: *maxValue, tickx: 1, ticky: *maxValue/20)
   unsigned iter = 0;
            log.Log( v: currentValue_vec[iter]);
```

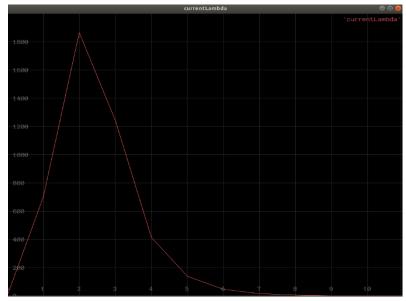


1.1 绘制阻尼因子µ随着迭代变化的曲线图

损失函数变化:



阻尼因子变化:





1.2 将曲线函数改成 $y = ax^2 + bx + c$

我们改变残差函数构建和残差对变量的雅可比:

```
// 误差模型 模板参数:观测值维度,类型,连接顶点类型
  virtual void ComputeResidual() override
      Vec3 abc = verticies_[0]->Parameters(); // 估计的参数
  virtual void ComputeJacobians() override
```

```
Test CurveFitting start...

iter: 0 , chi= 719.475 , Lambda= 0.001

iter: 1 , chi= 91.395 , Lambda= 0.000333333

problem solve cost: 2.68313 ms

makeHessian cost: 2.0453 ms

-----After optimization, we got these parameters :

1.61039  1.61853 0.995178

-----ground truth:

1.0, 2.0, 1.0
```

我们发现只有两步迭代, 迭代结束后的残差仍然较大, 得到的曲线参数离真实值误差较大。



1.3 实现更优秀的阻尼因子策略

新建变量update_strategy_用来存放用户想要输入的LM算法参数更新方法,可输入1,2或3:

首先是往Hessian矩阵中加入阻尼因子 λ 和去除 λ ,根据参数更新方法不同,有不同的做法:

$$\left[\mathbf{J}^{\mathsf{T}}\mathbf{W}\mathbf{J} + \lambda\mathbf{I}\right]\mathbf{h}_{\mathsf{lm}} = \mathbf{J}^{\mathsf{T}}\mathbf{W}(\mathbf{y} - \hat{\mathbf{y}}) , \qquad (12)$$

$$\left[\mathbf{J}^\mathsf{T} \mathbf{W} \mathbf{J} + \lambda \; \mathsf{diag}(\mathbf{J}^\mathsf{T} \mathbf{W} \mathbf{J}) \right] \mathbf{h}_{\mathsf{lm}} = \mathbf{J}^\mathsf{T} \mathbf{W} (\mathbf{y} - \hat{\mathbf{y}}) \; . \tag{}$$

```
cout << "Please enter a LM parameter update strategy (1, 2 or 3):" << endl;
cin >> update_strategy_;
```

```
void Problem::AddLambdatoHessianLM() {
   ulong size = Hessian .cols():
   for (ulong i = 0; i < size; ++i) {
void Problem::RemoveLambdaHessianLM() {
   ulong size = Hessian_.cols();
   assert(Hessian_.rows() == Hessian_.cols() && "Hessian is not square");
   // TODO:: 这里不应该减去一个,数值的反复加减容易造成数值精度出问题?而应该保存叠加 ambda前的值,在这里直接赋值
   for (ulong i = 0; i < size; ++i) {
```



1.3 实现更优秀的阻尼因子策略

方法一:

1. $\lambda_0 = \lambda_o$; λ_o is user-specified [5]. use eq'n (13) for \mathbf{h}_{lm} and eq'n (16) for ρ if $\rho_i(\mathbf{h}) > \epsilon_4$: $\mathbf{p} \leftarrow \mathbf{p} + \mathbf{h}$; $\lambda_{i+1} = \max[\lambda_i/L_{\downarrow}, 10^{-7}]$; otherwise: $\lambda_{i+1} = \min[\lambda_i L_{\uparrow}, 10^7]$;

$$\begin{array}{ll} \rho_{i}(\mathbf{h}_{\mathsf{lm}}) & = & \frac{\chi^{2}(\mathbf{p}) - \chi^{2}(\mathbf{p} + \mathbf{h}_{\mathsf{lm}})}{(\mathbf{y} - \hat{\mathbf{y}})^{\mathsf{T}}(\mathbf{y} - \hat{\mathbf{y}}) - (\mathbf{y} - \hat{\mathbf{y}} - \mathbf{J}\mathbf{h}_{\mathsf{lm}})^{\mathsf{T}}(\mathbf{y} - \hat{\mathbf{y}} - \mathbf{J}\mathbf{h}_{\mathsf{lm}})} \\ & = & \frac{\chi^{2}(\mathbf{p}) - \chi^{2}(\mathbf{p} + \mathbf{h}_{\mathsf{lm}})}{\mathbf{h}_{\mathsf{lm}}^{\mathsf{T}}(\lambda_{i}\mathbf{h}_{\mathsf{lm}} + \mathbf{J}^{\mathsf{T}}\mathbf{W}(\mathbf{y} - \hat{\mathbf{y}}(\mathbf{p})))} & \text{if using eq'n (12) for } \mathbf{h}_{\mathsf{lm}} \ (15) \\ & = & \frac{\chi^{2}(\mathbf{p}) - \chi^{2}(\mathbf{p} + \mathbf{h}_{\mathsf{lm}})}{\mathbf{h}_{\mathsf{lm}}^{\mathsf{T}}(\lambda_{i}\mathsf{diag}(\mathbf{J}^{\mathsf{T}}\mathbf{W}\mathbf{J})\mathbf{h}_{\mathsf{lm}} + \mathbf{J}^{\mathsf{T}}\mathbf{W}(\mathbf{y} - \hat{\mathbf{y}}(\mathbf{p})))} & \text{if using eq'n (13) for } \mathbf{h}_{\mathsf{lm}} \ (16) \end{array}$$

其中参数 L_{\uparrow} 和 L_{\downarrow} 根据文献可设为 L_{\uparrow} = 11, L_{\downarrow} = 9.

```
| bool Problem::IsGoodStepInLM() {
     double scale = 0:
    double tempChi = 0.0:
    for (auto edge : pair<...> : edges_) {
        tempChi += edge.second->Chi2();
        unsigned Hessian_size = Hessian_.rows();
        MatXX diag_Hessian = MatXX::Zero( rows Hessian_size, cols Hessian_size);
        for (ulong i = 0; i < Hessian_size; ++i) {
        scale = delta_x_.transpose() * (currentLambda_ * diag_Hessian * delta_x_ + b_);
        if (rho > 0 && isfinite(x tempChi)) // last step was good, 误差在下降
            currentChi = tempChi:
```



1.3 实现更优秀的阻尼因子策略

方法二:

2. $\lambda_0 = \lambda_o \max \left[\text{diag}[\mathbf{J}^\mathsf{T} \mathbf{W} \mathbf{J}] \right]; \lambda_o \text{ is user-specified.}$ use eq'n (12) for \mathbf{h}_{lm} and eq'n (15) for ρ $\alpha = \left(\left(\mathbf{J}^\mathsf{T} \mathbf{W} (\mathbf{y} - \hat{\mathbf{y}}(\mathbf{p})) \right)^\mathsf{T} \mathbf{h} \right) / \left(\left(\chi^2 (\mathbf{p} + \mathbf{h}) - \chi^2 (\mathbf{p}) \right) / 2 + 2 \left(\mathbf{J}^\mathsf{T} \mathbf{W} (\mathbf{y} - \hat{\mathbf{y}}(\mathbf{p})) \right)^\mathsf{T} \mathbf{h} \right);$ if $\rho_i(\alpha \mathbf{h}) > \epsilon_4$: $\mathbf{p} \leftarrow \mathbf{p} + \alpha \mathbf{h}; \lambda_{i+1} = \max \left[\lambda_i / (1 + \alpha), 10^{-7} \right];$ otherwise: $\lambda_{i+1} = \lambda_i + |\chi^2 (\mathbf{p} + \alpha \mathbf{h}) - \chi^2 (\mathbf{p})| / (2\alpha);$

方法二相较方法一更为复杂一些,我们需要先回滚 Δx ,然后再更新 $\alpha \cdot \Delta x$ 来计算残差和,然后根据计算出的rho和alpha来更新的阻尼因子 λ .

```
for (auto edge :pair<...> : edges_) {
if (rho > 0 && isfinite( X tempChi)) // last step was good, 误差在下降
```



1.3 实现更优秀的阻尼因子策略

方法三:

3. $\lambda_0 = \lambda_o \max \left[\text{diag}[\mathbf{J}^\mathsf{T} \mathbf{W} \mathbf{J}] \right]; \ \lambda_o \text{ is user-specified [6].}$ use eq'n (12) for \mathbf{h}_{lm} and eq'n (15) for ρ if $\rho_i(\mathbf{h}) > \epsilon_4$: $\mathbf{p} \leftarrow \mathbf{p} + \mathbf{h}$; $\lambda_{i+1} = \lambda_i \max \left[1/3, 1 - (2\rho_i - 1)^3 \right]; \nu_i = 2$; otherwise: $\lambda_{i+1} = \lambda_i \nu_i$; $\nu_{i+1} = 2\nu_i$;

方法3是课上介绍的Nielsen策略,也是给出的原代码中实现的参数更新方法。

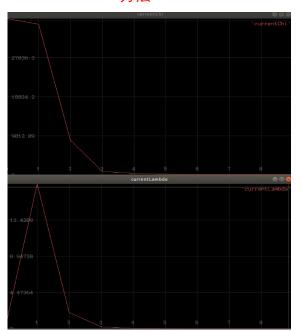
```
else if(update_strategy_ == 3) {
    scale += 1e-3; // make sure it's non-zero :)
    if (rho > 0 && isfinite(x tempChi)) // last step was good, 误差在下降
        double alpha = 1. - pow( x: (2 * rho - 1), y: 3);
        alpha = std::min(alpha, 2, / 3.):
        double scaleFactor = (std::max)(1. / 3., alpha);
        currentLambda_ *= scaleFactor;
        currentChi_ = tempChi;
    else {
```



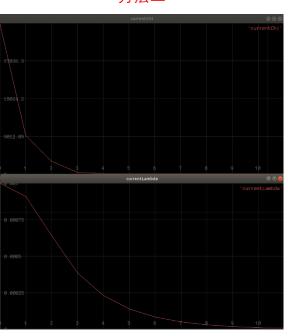
1.3 实现更优秀的阻尼因子策略

三种方法阻尼因子变化情况进行比较:

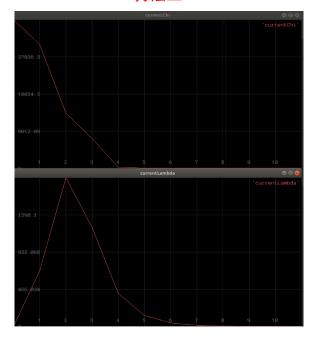
方法一



方法二



方法三



2. 公式推导雅可比F, G中的两项



2. 公式推导雅可比F, G中的两项



3. 证明式(9)



```
证明式(9):
 已知 非正益信息矩阵 丁丁 特征值 (7) 和对应特征 向量为 「约〕
 12日 LM がは中 (プリナル1) AXIm=-JT 中東新星 AXIm=- Z Vy FIT
iz:
     # F(x+0x) = L(0x) = 1 L(0x) ((0x)
                      = 1 (fx) + 14x) T (fx) + Jax)
                       = + f 7 + 0x 7 7 + + + 0x 7 7 7 0x
                       = FX) + (J'f) ax + ± ax J J ax
      3 \stackrel{*}{\underset{\sim}{}} : F'_{(x)} \approx (J^T f)^T, F''_{(x)} \approx J^T J
     接 プリニレハレブ 代入正規が経:
             (VAVT+MI) AXIm = - JTY
      由于特征相量矩阵为正交矩阵: VVT=I,特其什么,结
            (VAV + MVV ) AXIM = -J +
      => V(A+MI)V OXIM= -F'
       由于特征相景矩阵一定可逆,且以下=以一,可得:
         AXIM =-V (A+MI) -1V T FIT
            => OXIM = - Z ViTVi FIT
```



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