

## 第 12 节课习题

### 1. 使用 LM 算法估计曲线参数

#### 1) 绘制阻尼因子 $\mu$ 随着迭代变化的曲线图

在样例代码中我们使用 Pangolin 来绘制 残差平方和 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> &currentValue_vec, string name){
    // 参考 Pangolin Tutorial: https://github.com/yuntianli91/pangolin\_tutorial/blob/master/task5/main.cpp
    // Create OpenGL window in single line
    pangolin::CreateWindowAndBind( window_title: name, w: 640, h: 480);
    // Data logger object
    pangolin::DataLog log;
    // Optionally add named labels
    std::vector<std::string> labels;
    labels.push_back(std::string(name));
    log.SetLabels(labels);
    // OpenGL 'view' of data. We might have many views of the same data.
    auto max_value_iterator = max_element( first: currentValue_vec.begin(), last: currentValue_vec.end());
    pangolin::Plotter plotter( default_log: &log, left: 0, right: currentValue_vec.size() - 1, bottom: 0, top: *max_value, tickx: 1, ticky: *max_value/20);
    plotter.SetBounds( bottom: 0.0, top: 1.0, left: 0.0, right: 1.0);
    // plotter.Track("$i");//坐标轴自动滚动
    pangolin::DisplayBase().AddDisplay( &plotter);
    // Default hooks for exiting (Esc) and fullscreen (tab).
    unsigned iter = 0;
    while( !pangolin::ShouldQuit() )
    {
        glClear( mask: GL_COLOR_BUFFER_BIT | GL_DEPTH_BUFFER_BIT);

        if(iter < currentValue_vec.size()){
            log.Log( v: currentValue_vec[iter]);
        }
        iter++;

        // Render graph, Swap frames and Process Events
        pangolin::FinishFrame();
    }
}
```

并在 bool Problem::Solve(int iterations)函数内创建变量 currentChi\_vec 和 currentLambda\_vec 来存储所有迭代步骤的残差 currentChi\_和阻尼因子 currentLambda\_:

```
vector<double> currentChi_vec;
vector<double> currentLambda_vec;

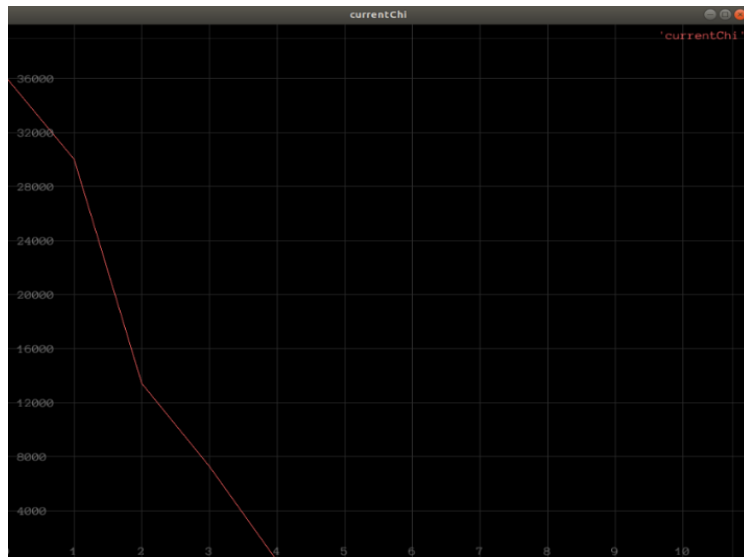
while (!stop && (iter < iterations)) {
    currentChi_vec.push_back(currentChi_);
    currentLambda_vec.push_back(currentLambda_);
}
```

并且在 bool Problem::Solve(int iterations)函数末尾调用 draw\_curve 函数来绘制残差和阻尼因子随着迭代步骤 iter 变化的曲线图:

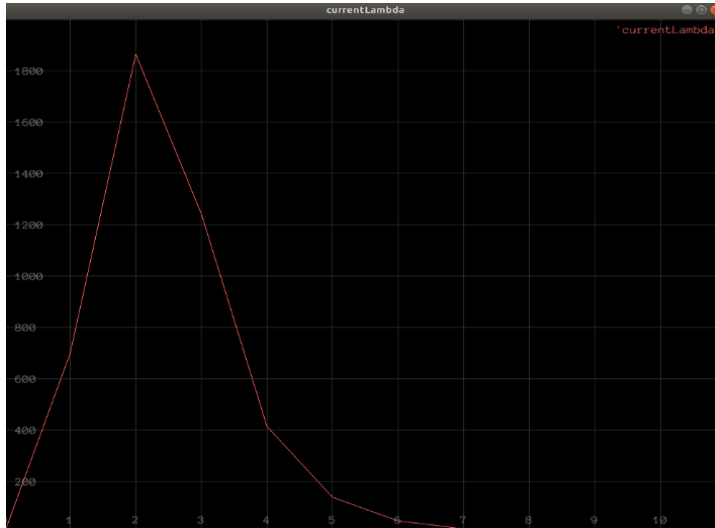
```
// draw the curve
draw_curve( currentValue_vec: currentChi_vec, name: "currentChi");
draw_curve( currentValue_vec: currentLambda_vec, name: "currentLambda");
```

曲线绘制结果如下:

残差平方和变化:



阻尼因子变化:



我们看到在总共的 11 步迭代步骤中，残差一直在下降，这是因为在算法中不满足更新策略的更新都被删除了，导致我们采用的更新都是使得残差缩小的。同时我们看到一开始阻尼因子很小，后来算法逐步增大阻尼因子，来缩小迭代步长，此时 LM 算法逐步趋近于最速下降算法。后来阻尼因子逐渐减小，LM 算法趋近于高斯牛顿。

## 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，代码更改如下：

```
// 该模型 模板参数：观测值维度，类型，连接顶点类型
class CurveFittingEdge: public Edge
{
public:
    EIGEN_MAKE_ALIGNED_OPERATOR_NEW
    CurveFittingEdge( double x, double y ): Edge( residual_dimension: 1, num_vertices: 1, vertices_types: std::vector<std::string>{"abc"}) {
        x_ = x;
        y_ = y;
    }
    // 计算曲线模型误差
    virtual void ComputeResidual() override
    {
        Vec3 abc = vertices_[0]->Parameters(); // 估计的参数
        // residual_(0) = std::exp( abc(0)*x_ + abc(1)*x_ + abc(2) ) - y_; // 构建残差
        residual_( index: 0 ) = abc( index: 0)*x_ + abc( index: 1)*x_ + abc( index: 2) - y_; // 构建残差
    }

    // 计算残差对变量的雅可比
    virtual void ComputeJacobians() override
    {
        // Vec3 abc = vertices_[0]->Parameters();
        // double exp_y = std::exp( abc(0)*x_ + abc(1)*x_ + abc(2) );
        Eigen::Matrix<double, 1, 3> jaco_abc; // 误差为1维，状态量 3 个，所以是 1x3 的雅可比矩阵

        // jaco_abc << x_ * x_ * exp_y, x_ * exp_y, 1 * exp_y;
        jaco_abc << x_ * x_, x_, 1;
        jacobians_[0] = jaco_abc;
    }
}
```

参数估计如下：

```
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。

### 3) 实现更优秀的阻尼因子策略, 参考文献 [1] 的 4.1.1 节

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:

```
bool Problem::Solve(int iterations) {
    cout << "Please enter a LM parameter update strategy (1, 2 or 3):" << endl;
    cin >> update_strategy_;
}
```

#### 方法一:

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

为了实现该参数更新方法, 源代码中要改变的代码如下。

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

```
void Problem::AddLambdatoHessianLM() {
    ulong size = Hessian_.cols();
    assert(Hessian_.rows() == Hessian_.cols() && "Hessian is not square");
    for (ulong i = 0; i < size; ++i) {
        if(update_strategy_ == 1)
            // Hessian_(i, i) += currentLambda_ + Hessian_(i, i);
            Hessian_(row: i, col: i) += (1.+currentLambda_);
        else if(update_strategy_ == 2 || update_strategy_ == 3)
            Hessian_(row: i, col: i) += currentLambda_;
        else{}
    }
}

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Lambda前的值, 在这里直接赋值
    for (ulong i = 0; i < size; ++i) {
        if(update_strategy_ == 1)
            // Hessian_(i, i) -= currentLambda_ + Hessian_(i, i); // false remove
            Hessian_(row: i, col: i) /= (1.+currentLambda_);
        else if(update_strategy_ == 2 || update_strategy_ == 3)
            Hessian_(row: i, col: i) -= currentLambda_;
        else{}
    }
}
```

相对应的, 我们在 `bool Problem::IsGoodStepInLM()` 中给出方法 1 的阻尼因子  $\lambda$  的更新方法, 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 rho = 0;

    // recompute residuals after update state
    // 统计所有的残差
    double tempChi = 0.0;
    for (auto edge : pair<...> : edges_) {
        edge.second->ComputeResidual();
        tempChi += edge.second->Chi2();
    }

    if(update_strategy_ == 1)
    {
        unsigned Hessian_size = Hessian_.rows();
        MatXX diag_Hessian = MatXX::Zero( rows: Hessian_size, cols: Hessian_size);
        for (ulong i = 0; i < Hessian_size; ++i) {
            diag_Hessian( row: i, col: i) = Hessian_( row: i, col: i);
        }
        scale = delta_x_.transpose() * (currentLambda_ + diag_Hessian * delta_x_ + b_);
        scale += 1e-3; // make sure it's non-zero :)
        rho = (currentChi_ - tempChi) / scale;

        double L_up = 11.0;
        double L_down = 9.0;
        if (rho > 0 && !isfinite( x tempChi)) // last step was good, 误差在下降
        {
            currentLambda_ = (std::max)(currentLambda_/L_down, 1e-7);
            currentChi_ = tempChi;
            return true;
        } else {
            currentLambda_ = (std::min)(currentLambda_ * L_up, 1e7);
            return false;
        }
    }
}
```

在函数 `void Problem::ComputeLambdaInitLM()` 中我们规定阻尼因子的初始值，三个方法的策略都是一样的：

$$\mu_0 = \tau \cdot \max \left\{ \left( \mathbf{J}^T \mathbf{J} \right)_{ii} \right\}$$

通常，按需设定  $\tau \sim [10^{-8}, 1]$ ，在原代码中  $\tau = 10^{-5}$ ：

```

// LM
void Problem::ComputeLambdaInitLM() {
    ni_ = 2.;
    currentLambda_ = -1.;
    currentChi_ = 0.0;
    // TODO: robust cost chi2
    for (auto edge : pairs_> : edges_) {
        currentChi_ += edge.second->Chi2();
    }
    if (err_prior_.rows() > 0)
        currentChi_ += err_prior_.norm();

    stopThresholdLM_ = 1e-6 * currentChi_; // 迭代条件为 误差下降 1e-6 倍

    double maxDiagonal = 0;
    ulong size = Hessian_.cols();
    assert(Hessian_.rows() == Hessian_.cols() && "Hessian is not square");
    for (ulong i = 0; i < size; ++i) {
        maxDiagonal = std::max(fabs(Hessian_(row: i, col: i)), maxDiagonal);
    }
    double tau = 1e-5;
    currentLambda_ = tau * maxDiagonal;
}

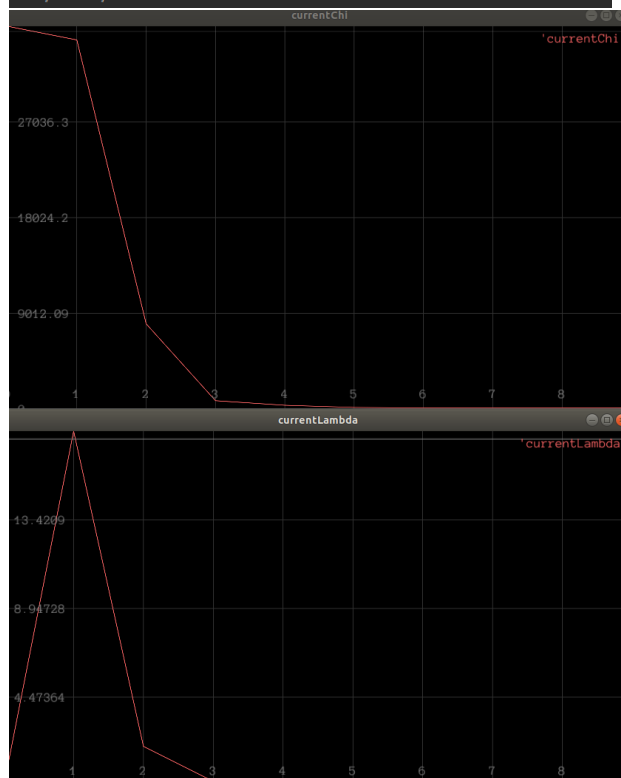
```

方法 1 计算得到的参数计算以及绘制如下：

```

Test CurveFitting start...
Please enter a LM parameter update strategy (1, 2 or 3):
1
iter: 0 , chi= 36048.3 , Lambda= 0.001
iter: 1 , chi= 34760.2 , Lambda= 17.8946
iter: 2 , chi= 8020.58 , Lambda= 1.98828
iter: 3 , chi= 779.997 , Lambda= 0.22092
iter: 4 , chi= 348.805 , Lambda= 0.0245467
iter: 5 , chi= 145.33 , Lambda= 0.00272741
iter: 6 , chi= 101 , Lambda= 0.000303046
iter: 7 , chi= 92.3181 , Lambda= 3.36718e-05
iter: 8 , chi= 91.3999 , Lambda= 3.74131e-06
iter: 9 , chi= 91.3959 , Lambda= 4.15701e-07
problem solve cost: 22.0139 ms
makeHessian cost: 14.5564 ms
-----After optimization, we got these parameters :
0.941955  2.0945  0.9656
-----ground truth:
1.0,  2.0,  1.0

```



## 方法二:

2.  $\lambda_0 = \lambda_o \max [\text{diag}[\mathbf{J}^T \mathbf{W} \mathbf{J}]]$ ;  $\lambda_o$  is user-specified.  
 use eq'n (12) for  $\mathbf{h}_{lm}$  and eq'n (15) for  $\rho$   

$$\alpha = \left( \left( \mathbf{J}^T \mathbf{W} (\mathbf{y} - \hat{\mathbf{y}}(\mathbf{p})) \right)^T \mathbf{h} \right) / \left( (\chi^2(\mathbf{p} + \mathbf{h}) - \chi^2(\mathbf{p})) / 2 + 2 \left( \mathbf{J}^T \mathbf{W} (\mathbf{y} - \hat{\mathbf{y}}(\mathbf{p})) \right)^T \mathbf{h} \right);$$
  
 if  $\rho_i(\alpha \mathbf{h}) > \epsilon_4$ :  $\mathbf{p} \leftarrow \mathbf{p} + \alpha \mathbf{h}$ ;  $\lambda_{i+1} = \max [\lambda_i / (1 + \alpha), 10^{-7}]$ ;  
 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 $\Delta x$ ，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  $\lambda$ :

```

else if(update_strategy_ == 2)
{
    double alpha = double(b_.transpose() * delta_x_) / ((tempChi - currentChi_) / 2.0 + 2.0 * b_.transpose() * delta_x_);
    alpha = std::max(alpha, 1e-1);
    // cout << "alpha_factor = " << alpha_factor << endl;

    // 回滚delta_x_
    RollbackStates();
    // 回滚delta_x_之后，更新状态量 X = X + alpha * delta_x
    delta_x_ *= alpha;
    UpdateStates();

    tempChi = 0.0;
    for (auto edge : pairs<> : edges_) {
        edge.second->ComputeResidual();
        tempChi += edge.second->Chi2();
    }

    scale = delta_x_.transpose() * (currentLambda_ * delta_x_ + b_);
    scale += 1e-3; // make sure it's non-zero :)

    rho = (currentChi_ - tempChi) / scale;

    if (rho > 0 && !isfinite((double) tempChi)) // last step was good, 误差在下降
    {
        currentLambda_ = (std::max)(currentLambda_ / (1 + alpha), 1e-7);
        currentChi_ = tempChi;
        return true;
    }
    else {
        currentLambda_ += abs((double) tempChi - currentChi_) / (2 * alpha);
        return false;
    }
}

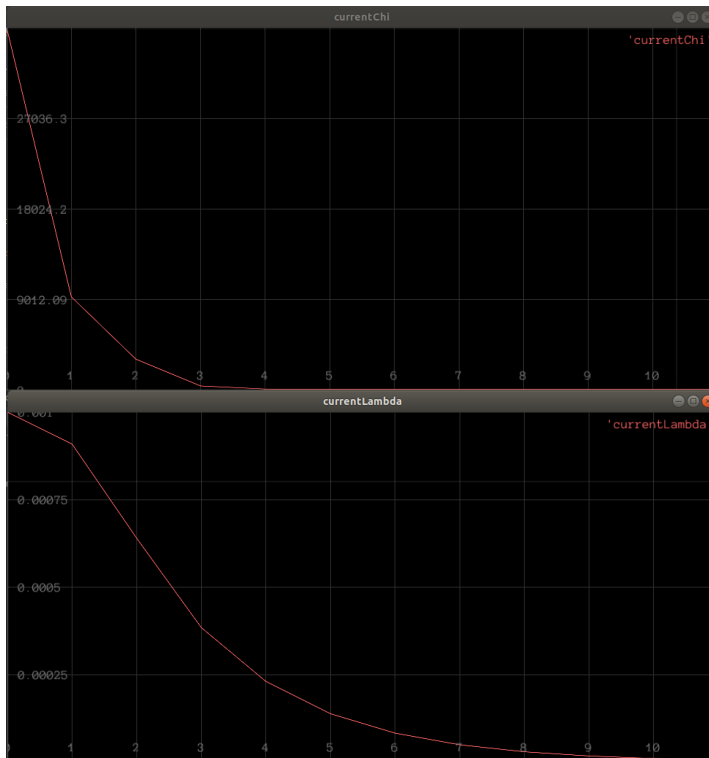
```

方法 2 计算得到的参数计算以及绘制如下:

```

Test CurveFitting start...
Please enter a LM parameter update strategy (1, 2 or 3):
2
iter: 0 , chi= 36048.3 , Lambda= 0.001
iter: 1 , chi= 9294.65 , Lambda= 0.000909091
iter: 2 , chi= 3097.71 , Lambda= 0.000640061
iter: 3 , chi= 436.783 , Lambda= 0.000384749
iter: 4 , chi= 132.226 , Lambda= 0.000230966
iter: 5 , chi= 96.0453 , Lambda= 0.000138626
iter: 6 , chi= 91.9154 , Lambda= 8.32001e-05
iter: 7 , chi= 91.4535 , Lambda= 4.99341e-05
iter: 8 , chi= 91.4022 , Lambda= 2.99687e-05
iter: 9 , chi= 91.3966 , Lambda= 1.7986e-05
iter: 10 , chi= 91.3959 , Lambda= 1.07944e-05
iter: 11 , chi= 91.3959 , Lambda= 6.47825e-06
problem solve cost: 21.1229 ms
makeHessian cost: 14.1695 ms
-----After optimization, we got these parameters :
0.942115 2.09435 0.965614
-----ground truth:
1.0, 2.0, 1.0

```



### 方法三:

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

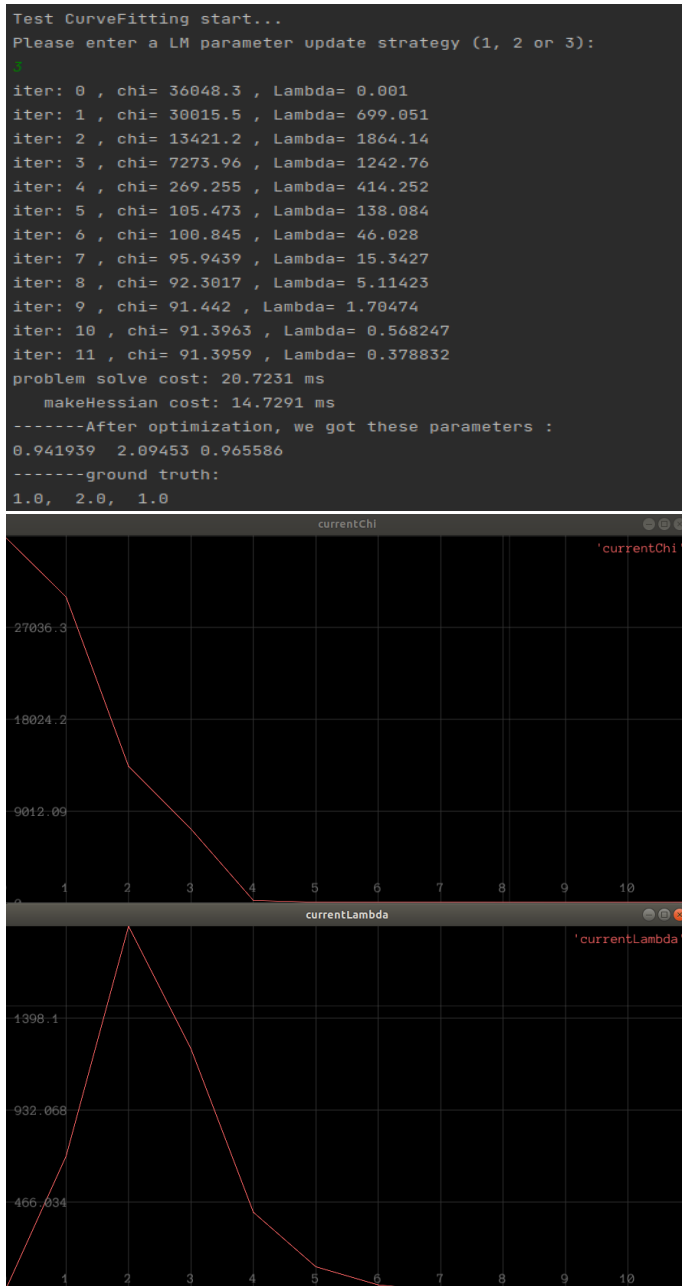
```

else if(update_strategy_ == 3) {
    scale = delta_x_.transpose() * (currentLambda_ * delta_x_ + b_);
    scale += 1e-3; // make sure it's non-zero :)
    rho = (currentChi_ - tempChi) / scale;

    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;
        ni_ = 2;
        currentChi_ = tempChi;
        return true;
    }
    else {
        currentLambda_ *= ni_;
        ni_ *= 2;
        return false;
    }
}
else{
    cout << "Can not find suitable LM update methods!" << endl;
    return false;
}

```

我们同样给出参数计算以及绘制情况:



### 总结:

比较以上三种方法，仅从我们有限的实验中，我们发现三种方法估计出的曲线参数相差不大，最后的残差和都比较接近，且迭代步骤都在 10 步左右。三种方法都是可行的 LM 参数更新方法。其中方法二最为复杂，它的特别之处在于根据比例因子来计算缩放因子 $\alpha$ ，改变更新步长为 $\alpha \cdot \Delta x$ 。从绘制的阻尼因子变化曲线来看，方法二的阻尼因子逐渐减小，变化最为平滑，方法三阻尼因子变化最为剧烈，所以从这个角度来看，方法二好于方法一，方法一好于方法三。当然可能在不同拟合参数的情况下，情况又会有所不同。



## 2. 公式推导雅可比 F, G 中的两项

位置预积分量  $\alpha$  的递推计算形式:

$$\begin{aligned}\alpha_{b|k+1} &= \alpha_{b|k} + \beta_{b|k} \delta t + \frac{1}{2} a \delta t^2 \\ &= \alpha_{b|k} + \beta_{b|k} \delta t + \frac{1}{4} \left( \mathbf{g}_{b|k} (a^{bk} - b_k^a) + \mathbf{g}_{b|k+1} (a^{bk+1} - b_k^a) \right) \delta t^2\end{aligned}$$

位置预积分量对  $k$  时刻角速度 bias 的 Jacobian:

$$\begin{aligned}\partial \alpha_{b|k+1} &= \partial \alpha_{b|k} + \partial \beta_{b|k} \delta t + \frac{1}{4} \left( \mathbf{g}_{b|k} (a^{bk} - b_k^a) + \mathbf{g}_{b|k} \otimes \left[ \frac{1}{2} \omega \delta t \right] (a^{bk+1} - b_k^a) \right) \delta t^2 \\ f_{15} = \frac{\partial \alpha_{b|k+1}}{\partial \delta b_k^g} &= \frac{1}{4} \frac{\partial \mathbf{g}_{b|k} \otimes \left[ \frac{1}{2} (\omega - \delta b_k^g) \delta t \right] (a^{bk+1} - b_k^a) \delta t^2}{\partial \delta b_k^g} \\ &= \frac{1}{4} \frac{\partial R_{b|k} \exp \left( \left[ (\omega - \delta b_k^g) \delta t \right]_{\times} \right) (a^{bk+1} - b_k^a) \delta t^2}{\partial \delta b_k^g} \\ &\approx \frac{1}{4} \frac{\partial R_{b|k} (1 + \left[ (\omega - \delta b_k^g) \delta t \right]_{\times}) (a^{bk+1} - b_k^a) \delta t^2}{\partial \delta b_k^g} \\ &= \frac{1}{4} \frac{-\partial R_{b|k} \left[ (a^{bk+1} - b_k^a) \right]_{\times} \delta t^2 \cdot [-\delta b_k^g \delta t]}{\partial \delta b_k^g} \\ &= -\frac{1}{4} \left( R_{b|k} \left[ (a^{bk+1} - b_k^a) \right]_{\times} \delta t^2 \right) (-\delta t)\end{aligned}$$

计算位置预积分量对  $k$  时刻角速度高斯白噪声的 Jacobian:

$$\begin{aligned}\omega &= \frac{1}{2} \left( (\bar{\omega}^{bk} + \mathbf{n}_k^g - b_k^g) + (\bar{\omega}^{bk+1} + \mathbf{n}_{k+1}^g - b_k^g) \right) = \frac{1}{2} \underbrace{(\bar{\omega}^{bk} + \bar{\omega}^{bk+1})}_{=\bar{\omega}} + \frac{1}{2} \mathbf{n}_k^g + \frac{1}{2} \mathbf{n}_{k+1}^g - b_k^g \\ \alpha_{b|k+1} &= \alpha_{b|k} + \beta_{b|k} \delta t + \frac{1}{4} \left( \mathbf{g}_{b|k} (a^{bk} - b_k^a) + \mathbf{g}_{b|k} \otimes \left[ \frac{1}{2} (\bar{\omega} + \frac{1}{2} \mathbf{n}_k^g + \frac{1}{2} \mathbf{n}_{k+1}^g - b_k^g) \right] (a^{bk+1} - b_k^a) \right) \delta t^2 \\ g_{12} = \frac{\partial \alpha_{b|k+1}}{\partial \delta \mathbf{n}_k^g} &= \frac{1}{4} \frac{\partial \mathbf{g}_{b|k} \otimes \left[ \frac{1}{2} (\bar{\omega} + \frac{1}{2} \mathbf{n}_k^g + \frac{1}{2} \mathbf{n}_{k+1}^g - b_k^g) \right] (a^{bk+1} - b_k^a) \delta t^2}{\partial \delta \mathbf{n}_k^g} \\ &= \frac{1}{4} \frac{\partial R_{b|k} \exp \left( \left[ \frac{1}{2} \delta \mathbf{n}_k^g \delta t \right]_{\times} \right) (a^{bk+1} - b_k^a) \delta t^2}{\partial \delta \mathbf{n}_k^g} \\ &= \frac{1}{4} \frac{\partial R_{b|k} (1 + \left[ \frac{1}{2} \delta \mathbf{n}_k^g \delta t \right]_{\times}) (a^{bk+1} - b_k^a) \delta t^2}{\partial \delta \mathbf{n}_k^g} \\ &= \frac{1}{4} \frac{\partial R_{b|k} \left[ (a^{bk+1} - b_k^a) \right]_{\times} \delta t^2 \cdot \left( \frac{1}{2} \delta \mathbf{n}_k^g \delta t \right)}{\partial \delta \mathbf{n}_k^g} \\ &= \frac{1}{4} R_{b|k} \left[ (a^{bk+1} - b_k^a) \right]_{\times} \delta t^2 \left( \frac{1}{2} \delta t \right)\end{aligned}$$

同理, 可计算位置预积分对  $k+1$  时刻角速度高斯白噪声的 Jacobian:

$$g_{14} = \frac{\partial \alpha_{b|k+1}}{\partial \delta \mathbf{n}_{k+1}^g} = g_{12}$$



### 3. 证明式(9)

证明式(9):

已知非负信息矩阵  $J^T J$  特征值  $\{\lambda_j\}$  和对应特征向量为  $\{V_j\}$

证明 LM 方法中  $(J^T J + \mu I) \Delta X_{lm} = -J^T f$  中更新量  $\Delta X_{lm} = -\sum_{j=1}^n \frac{V_j^T F'^T}{\lambda_j + \mu} V_j$

证:

$$\begin{aligned} \text{由 } F(x+\Delta x) &\approx L(\Delta x) = \frac{1}{2} L(\Delta x)^T L(\Delta x) \\ &= \frac{1}{2} (f(x) + J\Delta x)^T (f(x) + J\Delta x) \\ &= \frac{1}{2} f^T f + \Delta x^T J^T f + \frac{1}{2} \Delta x^T J^T J \Delta x \\ &= F(x) + (J^T f)^T \Delta x + \frac{1}{2} \Delta x^T J^T J \Delta x \end{aligned}$$

$$\text{可知: } F'(x) \approx (J^T f)^T, \quad F''(x) \approx J^T J$$

将  $J^T J = V\Lambda V^T$  代入正规方程:

$$(V\Lambda V^T + \mu I) \Delta X_{lm} = -J^T f$$

由于特征相量矩阵为正交矩阵:  $VV^T = I$ , 将其代入, 得

$$(V\Lambda V^T + \mu VV^T) \Delta X_{lm} = -J^T f$$

$$\Rightarrow V(\Lambda + \mu I) V^T \Delta X_{lm} = -F'^T$$

由于特征相量矩阵一定可逆, 且  $V^T = V^{-1}$ , 可得:

$$\Delta X_{lm} = -V(\Lambda + \mu I)^{-1} V^T F'^T$$

$$\Rightarrow \Delta X_{lm} = -\sum_{j=1}^n \frac{V_j^T V_j}{\lambda_j + \mu} F'^T$$

#### 参考文献:

[1] The Levenberg-Marquardt method for nonlinear least squares curve-fitting problems, Henri P. Gavin, Department of Civil and Environmental Engineering, Duke University, 2016