

## Training Process

#### **Common Process of NN**

My LSTM

for each epoch:

for each X, Y in data iter:

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⑥藏状态初始化

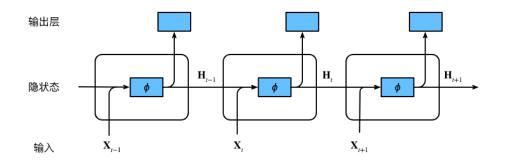
正向传播: y\_hat = net(X)

母 损失计算: I = loss(y\_hat, y)

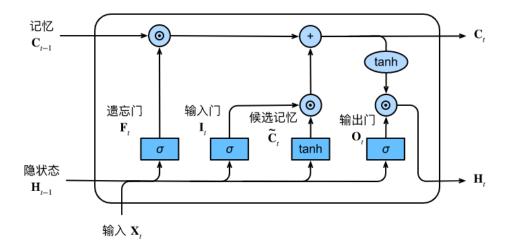
梯度归零: updater.zero\_grad()

参数更新: updater.step()

# $y_hat = net(x)$



PRNN model: Each neuro utilizes the information from the prior neuro



#### For each neuro,

Inner structure is packaged neatly

```
self.layer = nn.LSTM(feature_size, hidden_size, num_layers)
self.linear = nn.Linear(hidden_size, output_size)
```

• What we need:  $C_0$ ,  $H_0$ ,  $X_t$ 

```
output, state = self.layer(X, state)
y_hat = self.linear(output)
```

## $C_0$ , $H_0$

#### Formula

$$\left\{ \begin{array}{l} \mathbf{I}_t = \sigma(\mathbf{X}_t \mathbf{W}_{xi} + \mathbf{H}_{t-1} \mathbf{W}_{hi} + \mathbf{b}_i), \\ \mathbf{F}_t = \sigma(\mathbf{X}_t \mathbf{W}_{xf} + \mathbf{H}_{t-1} \mathbf{W}_{hf} + \mathbf{b}_f), \\ \mathbf{O}_t = \sigma(\mathbf{X}_t \mathbf{W}_{xo} + \mathbf{H}_{t-1} \mathbf{W}_{ho} + \mathbf{b}_o), \\ \tilde{\mathbf{C}}_t = \tanh(\mathbf{X}_t \mathbf{W}_{xc} + \mathbf{H}_{t-1} \mathbf{W}_{hc} + \mathbf{b}_c), \end{array} \right. \left\{ \begin{array}{l} \mathbf{C}_t = \mathbf{F}_t \odot \mathbf{C}_{t-1} + \mathbf{I}_t \odot \tilde{\mathbf{C}}_t. \\ \mathbf{H}_t = \mathbf{O}_t \odot \tanh(\mathbf{C}_t). \\ \mathbf{\hat{Y}}_t = \mathbf{H}_t \mathbf{W}_{hq} + \mathbf{b}_q. \end{array} \right.$$

# 记忆 $\mathbf{C}_{t-1}$ $\mathbf{G}_{t-1}$ $\mathbf{G}_{t-1}$

#### Shape

 $W_{\chi}$  defined in nn.LSTM: [feature\_size, hidden\_size]  $W_h$  defined in nn.LSTM: [hidden\_size, hidden\_size]  $X_t$ : [batch\_size, feature\_size]  $X_t$ : [time\_steps, batch\_size, feature\_size]  $X_t$ : [time\_steps, batch\_size, feature\_size]  $X_t$ : [batch\_size, hidden\_size, output\_size]  $X_t$ : [batch\_size, hidden\_size]

## • 隐藏状态初始化 $C_0, H_0$

## $MyDataLoader - X_t$

## X: [time\_steps, batch\_size, feature\_size]

- times\_steps为超参数,表示每次预测要用过去多长时间的数据
- feature\_size代表特征数,这是由因子数量决定的
- batch\_size是可变的,这是小批量梯度下降中用到的样本数量

## 实现过程中有遇到很多问题——

假定当前有3000支股票,100期数据,300个因子,根据经验,time\_steps=5是个不错的选择,现在需要根据当前数据生成X[5, batch\_size, 300]

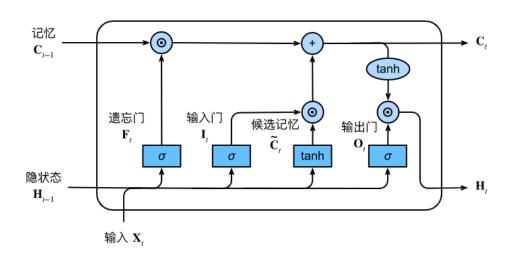
- Q1: X在时间上是连续的还是重叠的?
  - t1-t5, t6-t10, ..., t96-t100 / t1-t5, t2-t6, t3-t7, ..., t95-t99, t96-t100
- Q2: 在某些截面,一些股票会被禁止交易,如停牌、ST
  - batch\_size可变 / batch\_size不变, 丢弃这些股票
  - 如果股票总数只有10,是否batch\_size ≤10
- Q3: 在同一个X中,是否需要维持每只股票的位置一致?
  - 在每个time\_step中,第一条样本为股票A,第二条样本为股票B,...

## 损失计算: I = loss(y hat, y)

#### Formula

$$\left\{ \begin{array}{l} \mathbf{I}_t = \sigma(\mathbf{X}_t \mathbf{W}_{xi} + \mathbf{H}_{t-1} \mathbf{W}_{hi} + \mathbf{b}_i), \\ \mathbf{F}_t = \sigma(\mathbf{X}_t \mathbf{W}_{xf} + \mathbf{H}_{t-1} \mathbf{W}_{hf} + \mathbf{b}_f), \\ \mathbf{O}_t = \sigma(\mathbf{X}_t \mathbf{W}_{xo} + \mathbf{H}_{t-1} \mathbf{W}_{ho} + \mathbf{b}_o), \\ \tilde{\mathbf{C}}_t = \tanh(\mathbf{X}_t \mathbf{W}_{xc} + \mathbf{H}_{t-1} \mathbf{W}_{hc} + \mathbf{b}_c), \end{array} \right. \left\{ \begin{array}{l} \mathbf{C}_t = \mathbf{F}_t \odot \mathbf{C}_{t-1} + \mathbf{I}_t \odot \mathbf{C}_t \\ \mathbf{H}_t = \mathbf{O}_t \odot \tanh(\mathbf{C}_t), \\ \mathbf{\hat{Y}}_t = \mathbf{H}_t \mathbf{W}_{hq} + \mathbf{b}_q. \end{array} \right.$$

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• 只需要最后的 $\hat{y}_T$ , 前 $time_steps$ 的数据都是炮灰

```
y = (Y[-1].reshape(-1)).to(device) # 最后一期y拉成向量
y_hat = y_hat[-1].reshape(-1) # [批量大小,输出大小]
```

回归问题

```
loss = nn.MSELoss()
I = loss(y hat, y.float())
```

• 分类问题

```
loss = nn.CrossEntropyLoss()
# 好像不需要 y_hat = nn.functional.softmax(y_hat, dim=1)
I = loss(y hat, y.long()) # label一定要≥0
```

## 梯度裁剪

## RNN模型的梯度爆炸和梯度消失问题:

- 在反向传播过程中产生长度为O(T)的矩阵乘法链
- 当T较大时,它可能导致数值不稳定
- RNN往往需要额外的方式来支持稳定训练

当梯度很大或者学习率很大时,都会出现模型反复跳跃,训练无法收敛的问题:

$$|f(\mathbf{x}) - f(\mathbf{x} - \eta \mathbf{g})| \le L\eta \|\mathbf{g}\|$$

## 解决梯度爆炸问题:

- 限定一个小的学习率是可行的
- 但如果g的过大只是偶发性的 → 梯度裁剪

$$\mathbf{g} \leftarrow \min\left(1, \frac{\theta}{\|\mathbf{g}\|}\right) \mathbf{g}$$

```
@staticmethod
def grad_clipping(net, theta):
    # 所有层的参数统一裁剪
    params = [p for p in net.parameters() if p.requires_grad]
    norm = torch.sqrt(sum(torch.sum((p.grad ** 2)) for p in params))
    if norm > theta:
        for param in params:
            param.grad[:] *= theta / norm
```

## Predict

X: [time\_steps, batch\_size, feature\_size]

- times\_steps为超参数,表示每次预测要用过去多长时间的数据
- feature\_size代表特征数,这是由因子数量决定的
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假定当前有3000支股票,100期数据,300个因子,根据经验,time\_steps=5是个不错的选择,现在需要根据当前数据生成X[5, batch\_size, 300]

- 每次预测都重置隐藏状态
- 需要time\_steps-1期的预热数据
- X在时间上是重叠的
  - t1-t5, t2-t6, t3-t7, ..., t95-t99, t96-t100
- 如果某只股票不够连续的5期则不能预测
- batch\_size可以随意变动
- 在一个X中,需要维持每只股票的位置一致

```
for inputs, _, days, seclDs in test_datalter:
   batch_size = inputs.shape[1]
   state = net.begin_state(batch_size=batch_size, device=device)
   inputs = inputs.to(device)
   outputs, state = net(inputs, state)
   outputs = outputs.reshape((len(days), len(seclDs)))[-1]
```

