

# 실습2: svm 후반부





## 확률적 경사 하강법

- SVM에서 손실과 그래디언트에 대한 벡터화 사용 계산 구현 완료
- 확률적 경사 하강법을 사용하여 손실을 최소화할 예정

Al1/Al/classifiers/linear\_classifier.py 파일 안의 LinearClassifier 클
 래스의 train 메서드 구현 필요

```
class LinearClassifier(object):
    def __init__(self):
        self.W = None
    def train(
        self.
       learning rate=1e-3.
       reg=1e-5,
       num_iters=100,
       batch size=200.
       verbose=False.
       Train this linear classifier using stochastic gradient descent.
        Inputs:
       - X: A numpy array of shape (N, D) containing training data; there are N
         training samples each of dimension D.
       - y: A numpy array of shape (N,) containing training labels; y[i] = c
         means that X[i] has label 0 <= c < C for C classes.
       - learning_rate: (float) learning rate for optimization.
       - reg: (float) regularization strength.
       - num iters: (integer) number of steps to take when optimizing
       - batch_size: (integer) number of training examples to use at each step.
       - verbose: (boolean) If true, print progress during optimization.
       Outputs:
        A list containing the value of the loss function at each training iteration.
```

```
# TODO:
# Sample batch size elements from the training data and their
# corresponding labels to use in this round of gradient descent.
# Store the data in X batch and their corresponding labels in
# y_batch; after sampling X_batch should have shape (batch_size, dim)
# and y_batch should have shape (batch_size,)
# Hint: Use np.random.choice to generate indices. Sampling with
# replacement is faster than sampling without replacement.
# ****START OF YOUR CODE (DO NOT DELETE/MODIFY THIS LINE)****
pass
# ****END OF YOUR CODE (DO NOT DELETE/MODIFY THIS LINE)****
# evaluate loss and gradient
loss, grad = self.loss(X batch, y batch, reg)
loss history.append(loss)
# perform parameter update
# Update the weights using the gradient and the learning rate.
# ****START OF YOUR CODE (DO NOT DELETE/MODIFY THIS LINE) ****
pass
```

• 훈련세트에서 batch\_size 크기의 배치를 선택

```
num train, dim = X.shape
num classes = (
   np.max(y) + 1
 # assume y takes values 0...K-1 where K is number of classes
if self W is None:
   # lazily initialize W
   self.W = 0.001 * np.random.randn(dim, num_classes)
# Run stochastic gradient descent to optimize W
loss history = []
for it in range(num_iters):
   X batch = None
   y batch = None
   # TODO:
   # Sample batch size elements from the training data and their
   # corresponding labels to use in this round of gradient descent.
   # Store the data in X batch and their corresponding labels in
   # y batch; after sampling X batch should have shape (batch size, dim)
   # and y_batch should have shape (batch_size,)
   # Hint: Use np.random.choice to generate indices. Sampling with
   # replacement is faster than sampling without replacement.
   # ****START OF YOUR CODE (DO NOT DELETE/MODIFY THIS LINE)****
   pass
```

- 순차적 선택이 아닌 랜덤 선택 사용
- 중복을 허용하는 복원 추출 사용
- batch\_size x D 크기의 배열 X\_batch 및 batch\_size 크기 배열
   y\_batch 생성

- loss 메서드는 이미 구현이 되어있음
- 가중치 W 업데이트 부분을 구현해야함

```
# Vanilla Minibatch Gradient Descent

while True:

data_batch = sample_training_data(data, 256) # sample 256 examples

weights_grad = evaluate_gradient(loss_fun, data_batch, weights)

weights += - step_size * weights_grad # perform parameter update
```

## 확률적 경사 하강법

## 손실 그래프 그리기

```
# A useful debugging strategy is to plot the loss as a function of
# iteration number:
plt.plot(loss_hist)
plt.xlabel('Iteration number')
plt.ylabel('Loss value')
plt.show()
```

## 구현2: LinearClassifier 클래스의 predict 메서드

- 점수 행렬 계산:  $S = X \cdot W$
- 점수 행렬의 각 행에서 가장 큰 점수를 가진 것의 인덱스(label)을
  - 구하면 됨
- np.argmax 함수활용 가능

```
def predict(self, X):
   Use the trained weights of this linear classifier to predict labels for
   data points.
   Inputs:
   - X: A numpy array of shape (N, D) containing training data; there are N
    training samples each of dimension D.
   Returns:
   - y_pred: Predicted labels for the data in X. y_pred is a 1-dimensional
    array of length N, and each element is an integer giving the predicted
    class
   y_pred = np.zeros(X.shape[0])
   # Implement this method. Store the predicted labels in y_pred.
   # ****START OF YOUR CODE (DO NOT DELETE/MODIFY THIS LINE)****
   pass
   # ****END OF YOUR CODE (DO NOT DELETE/MODIFY THIS LINE)****
   return y_pred
```

## 구현3: 검증 과정 구현

```
# Use the validation set to tune hyperparameters (regularization strength and
# learning rate). You should experiment with different ranges for the learning
# rates and regularization strengths; if you are careful you should be able to
# get a classification accuracy of about 0.39 (> 0.385) on the validation set.
# Note: you may see runtime/overflow warnings during hyper-parameter search.
# This may be caused by extreme values, and is not a bug.
# results is dictionary mapping tuples of the form
# (learning rate, regularization strength) to tuples of the form
# (training_accuracy, validation_accuracy). The accuracy is simply the fraction
# of data points that are correctly classified.
results = {}
best val = -1 # The highest validation accuracy that we have seen so far.
best_svm = None # The LinearSVM object that achieved the highest validation rate.
# TODO:
# Write code that chooses the best hyperparameters by tuning on the validation #
# set. For each combination of hyperparameters, train a linear SVM on the
# training set, compute its accuracy on the training and validation sets, and
# store these numbers in the results dictionary. In addition, store the best
# validation accuracy in best val and the LinearSVM object that achieves this
# accuracy in best_svm.
# Hint: You should use a small value for num iters as you develop your
# validation code so that the SVMs don't take much time to train; once you are
# confident that your validation code works, you should rerun the validation
# code with a larger value for num iters.
```

## 구현3: 검증 과정 구현

- 학습률, 정규화 강도에 대한 이중 loop 사용
- train 및 predict 메서드를 사용하여 훈련 및 레이블 예측
- 훈련세트 및 검증세트 각각에 대한 정확도를 모두 구해야함

## 정확도 시각화

```
# Visualize the cross-validation results
import math
import pdb
# pdb.set_trace()
x scatter = [math.log10(x[0])] for x in results]
y_scatter = [math.log10(x[1]) for x in results]
# plot training accuracy
marker size = 100
colors = [results[x][0]] for x in results]
plt.subplot(2, 1, 1)
plt.tight_layout(pad=3)
plt.scatter(x_scatter, y_scatter, marker_size, c=colors, cmap=plt.cm.coolwarm)
plt.colorbar()
plt.xlabel('log learning rate')
plt.ylabel('log regularization strength')
plt.title('CIFAR-10 training accuracy')
# plot validation accuracy
colors = [results[x][1] for x in results] # default size of markers is 20
plt.subplot(2, 1, 2)
plt.scatter(x_scatter, y_scatter, marker_size, c=colors, cmap=plt.cm.coolwarm)
plt.colorbar()
plt.xlabel('log learning rate')
plt.ylabel('log regularization strength')
plt.title('CIFAR-10 validation accuracy')
plt.show()
```

## 테스트 세트 분류

```
# Evaluate the best svm on test set
y_test_pred = best_svm.predict(X_test)
test_accuracy = np.mean(y_test == y_test_pred)
print('linear SVM on raw pixels final test set accuracy: %f' % test_accuracy)
```

## 템플릿 그리기

• 훈련된 가중치 행렬을 이미지로 그림 (템플릿)

```
# Visualize the learned weights for each class.
# Depending on your choice of learning rate and regularization strength, these may
# or may not be nice to look at.
w = best_svm.W[:-1,:] # strip out the bias
w = w.reshape(32, 32, 3, 10)
w_min, w_max = np.min(w), np.max(w)
classes = ['plane', 'car', 'bird', 'cat', 'deer', 'dog', 'frog', 'horse', 'ship', 'truck']
for i in range(10):
   plt.subplot(2, 5, i + 1)

# Rescale the weights to be between 0 and 255
wimg = 255.0 * (w[:, :, :, i].squeeze() - w_min) / (w_max - w_min)
plt.imshow(wimg.astype('uint8'))
plt.axis('off')
plt.title(classes[i])
```

#### 과제2

#### · 인공지능 과제2

위에서는 미니배치를 선택할 때 "랜덤 선택" 방법을 사용하였다. 즉, 훈련세트에서 배치를 매번 랜덤으로 선택하였다. 이는 데이터의 다양성을 확보하지만 일부 이미지가 누락될 수도 있는 방법이다. 과제2에서는 각 반복에서 훈련세트를 랜덤하게 셔플링한 후 작은 배치들로 나누고, 순차적으로 각 배치를 사용하여 파라미터를 업데이트하는 방식을 사용한다. svm.train\_sequential이라는 함수가 바로 이렇게 셔플링+순차적선택 방식으로 훈련하는 함수이다. 이 svm.train\_sequential 함수 내용을 채워야한다.

```
[] # linear_classifier.py파일메서 LinearClassifier.train_sequential() 함수를 구현한 후 이 코드블럭을 실행시켜야함.
from Al.classifiers import LinearSVM
svm = LinearSVM()
tic = time.time()
loss_hist = svm.train_sequential(X_train, y_train, learning_rate=1e-7, reg=2.5e4, verbose=True)
toc = time.time()
print('That took %fs' % (toc - tic))
```

#### 과제2

```
def train_sequential(
  self.
  Χ,
  У.
  learning rate=1e-3.
  reg=1e-5,
  num_epochs=10,
  batch_size=200,
  verbose=False.
  num_train, dim = X.shape
  num classes = (
     np.max(y) + 1
  ) # assume v takes values 0...K-1 where K is number of classes
  if self.W is None:
     # lazily initialize W
     self.W = 0.001 * np.random.randn(dim, num_classes)
  # Run stochastic gradient descent to optimize W
  loss_history = []
   for epoch in range(num_epochs):
     # TODO:
     # 한 epoch이란 미니배치 경사하강법에서 훈련세트를 여러 배치로 나눈다음
     # 순차적으로 모든 배치에 파라미터 업데이트를 하는 것을 의미한다. 즉기매 #
     # epoch에서 모든 훈련세트 이미지를 한번씩 학습에 활용하게 된다. 여러
     # epoch은 훈련세트 이미지들을 한번씩 학습하는 과정을 여러번 반복하는
     # 것이다. 각 epoch에서 먼저, 훈련세트 이미지들을 랜덤으로 셔플링해야한다#
     # 이를 위해 random.shuffle 함수를 활용할 수 있다. 그런 다음, 각 epoch #
     # 메서는 훈련세트를 batch_size 크기의 배치들로 나눈다. 그리고, 순차적
     # 으로 각 배치에 대해 파라미터 업데이트를 수행한다. X_batch, y_batch는 #
     # 각각의 배치의 이미지 데이터 및 정답 레이블이다.
```

```
# Random shuffling
    pass
    for i in range(int(num_train/batch_size)):
     X batch = None
     v batch = None
     # sampling
      pass
     # evaluate loss and gradient
      loss, grad = self.loss(X_batch, y_batch, reg)
      loss_history.append(loss)
     # print("loss, grad: ", loss, grad)
     self. # += -learning_rate * grad
      if verbose and i%100 == 0:
        print("epoch %d, loss %f" % (epoch, loss))
return loss_history
```

## 과제2

#### Question 1

위에서는 학습률(learning rate)로 1e-7, 1e-6를 사용하였다. 만약 이보다 더 큰 학습률 (e.g., e-5, 5e-5)를 사용하면 성능이 어떤지 쓰시오. 직접 학습률을 바꿔서 돌려보되, 그 상태로 과제제출을 하지 말고 이후에 다시 작은 학습률(1e-7, 1e-6)에 대한 실험을 하여 그 결과가 있는 채로 과제를 제출해야한다.

답: