# DS-GA 1003: Machine Learning and Computational Statistics Homework 3: SVM and Sentiment Analysis

Due: Monday, February 29, 2016, at 6pm (Submit via NYU Classes)

Instructions: Your answers to the questions below, including plots and mathematical work, should be submitted as a single file, either HTML or PDF. You may include your code inline or submit it as a separate file. You may either scan hand-written work or, preferably, write your answers using software that typesets mathematics (e.g. L\*TFX, LyX, or MathJax via iPython).

### 1 Introduction

In this assignment, we'll be working with natural language data. In particular, we'll be doing sentiment analysis on movie reviews. This problem will give you the opportunity to try your hand at feature engineering, which is one of the most important parts of many data science problems. From a technical standpoint, this homework has two new pieces. First, you'll be implementing Pegasos. Pegasos is essentially stochastic subgradient descent for the SVM with a particular schedule for the step-size. Second, because in natural language domains we typically have huge feature spaces, we work with sparse representations of feature vectors, where only the non-zero entries are explicitly recorded. This will require coding your gradient and SGD code using hash tables (dictionaries in Python), rather than numpy arrays. We begin with some practice with subgradients and an easy problem that introduces the Perceptron algorithm.

## 2 Calculating Subgradients

Recall that a vector  $q \in \mathbf{R}^d$  is a subgradient of  $f : \mathbf{R}^d \to \mathbf{R}$  at x if for all z,

$$f(z) \ge f(x) + g^T(z - x).$$

As we noted in lecture, there may be 0, 1, or infinitely many subgradients at any point. The **subdifferential** of f at a point x, denoted  $\partial f(x)$ , is the set of all subgradients of f at x.

Just as there is a calculus for gradients, there is a calculus for subgradients<sup>1</sup>. For our purposes, we can usually get by using the definition of subgradient directly. However, in the first problem below we derive a property that will make our life easier for finding a subgradient of the hinge loss and perceptron loss.

 $<sup>^1\</sup>mathrm{A}$  good reference for subgradients are these notes https://see.stanford.edu/materials/lsocoee364b/01-subgradients\_notes.pdf.

1. [Subgradients for pointwise maximum of functions] Suppose  $f_1, \ldots, f_m : \mathbf{R}^d \to \mathbf{R}$  are convex functions, and

$$f(x) = \max_{i=1,\dots,m} f_i(x).$$

Let k be any index for which  $f_k(x) = f(x)$ , and choose  $g \in \partial f_k(x)$ . [We are using the fact that a convex function on  $\mathbf{R}^d$  has a non-empty subdifferential at all points.] Show that  $g \in \partial f(x)$ .

#### Solution:

Fix any  $z \in \mathbf{R}^d$ . Then

$$f(z) \ge f_k(z)$$
 by definition of  $f(x)$   
 $\ge f_k(x) + g^T(z - x)$  since  $g \in \partial f_k(x)$   
 $= f(x) + g^T(z - x)$  by choice of  $k$ .

Thus  $f(z) \ge f(x) + g^T(z-x)$ . Since this is true for all  $z \in \mathbf{R}^d$ , we have shown  $g \in \partial f(x)$ .

2. [Subgradient of hinge loss for linear prediction] Give a subgradient of

$$J(w) = \max\left\{0, 1 - yw^T x\right\}.$$

#### Solution:

By the previous problem, a subgradient is given by

$$g = \begin{cases} -yx & \text{for } yw^T x < 1\\ 0 & \text{for } yw^T x \ge 1. \end{cases}$$

## 3 Perceptron

The perceptron algorithm is often the first classification algorithm taught in machine learning classes. Suppose we have a labeled training set  $(x_1, y_1), \ldots, (x_n, y_n) \in \mathbf{R}^d \times \{-1, 1\}$ . In the perceptron algorithm, we are looking for a hyperplane that perfectly separates the classes. That is, we're looking for  $w \in \mathbf{R}^d$  such that

$$y_i w^T x_i > 0 \ \forall i \in \{1, \dots, n\}.$$

Visually, this would mean that all the x's for which y = 1 are on one side of the hyperplane  $\{x \mid w^T x = 0\}$ , and all the x's for which y = -1 are on the other side. When such a hyperplane exists, we say that the data are **linearly separable**. The perceptron algorithm is given in Algorithm 1.

There is also something called the **perceptron loss**, given by

$$\ell(\hat{y}, y) = \max\{0, -\hat{y}y\}.$$

In this problem we will see why this loss function has this name.

### Algorithm 1: Perceptron Algorithm

```
input: Training set (x_1,y_1),\ldots,(x_n,y_n)\in\mathbf{R}^d\times\{-1,1\} w^{(0)}=(0,\ldots,0)\in\mathbf{R}^d k=0 # step number repeat all_correct = TRUE for i=1,2,\ldots,n # loop through data if (y_ix_i^Tw^{(k)}\leq 0) w^{(k+1)}=w^{(k)}+y_ix_i all_correct = FALSE else w^{(k+1)}=w^{(k)} end if k=k+1 end for until (all_correct == TRUE) return w^{(k)}
```

1. Show that if  $\{x \mid w^T x = 0\}$  is a separating hyperplane for a training set  $\mathcal{D} = ((x_1, y_1), \dots, (x_n, y_n))$ , then the average perceptron loss on  $\mathcal{D}$  is 0.

#### Solution:

If  $w^T x = 0$  is a separating hyperplane, then  $y_i w^T x_i > 0$  for all i, and  $\ell(w^T x_i, y_i) = \max\{0, -w^T x_i y_i\} = 0$  for all i. Thus the average over i is also 0.

2. Let  $\mathcal{H}$  be the linear hypothesis space consisting of functions  $x \mapsto w^T x$ . Consider running SGD to minimize the empirical risk with the perceptron loss. We'll use the version of SGD in which we cycle through the data points in each epoch. Show that if we use a fixed step size 1 and we terminate when our training loss is 0, then we are exactly doing the Perceptron algorithm.

#### Solution:

The empirical risk is given by

$$J(w) = \frac{1}{n} \sum_{i=1}^{n} \max\{0, -y_i w^T x_i\}.$$

Let  $f_i = \max\{0, -y_i w^T x_i\}$ . A subgradient of  $f_i$  at w is

$$g = \begin{cases} -y_i x_i & \text{for } y_i w^T x_i < 0\\ 0 & \text{for } y_i w^T x_i \ge 0. \end{cases}$$

With a step size of 1, the step from  $w^{(k)}$  to  $w^{(k+1)}$  is given by

$$w^{(k+1)} = \begin{cases} w^{(k)} - (-y_i x_i) & \text{for } y_i w^T x_i < 0 \\ w^{(k)} & \text{for } y_i w^T x_i \ge 0, \end{cases}$$

which is exactly implemented by the perceptron algorithm.

3. Suppose the perceptron algorithm returns w. Show that w is a linear combination of the input points. That is, we can write  $w = \sum_{i=1}^{n} \alpha_i x_i$  for some  $\alpha_1, \ldots, \alpha_n \in \mathbf{R}$ . The  $x_i$  for which  $\alpha_i \neq 0$  are called support vectors. Give a characterization of points that are support vectors and not support vectors.

#### Solution:

Note that initially w = 0, and we only ever add to it multiples of  $x_i$  for some i. Thus w is a linear combination of  $x_i$ 's. If a point  $x_i$  was never misclassified when it was examined by the algorithm, then it is not a support vector. Otherwise, it is.

### 4 The Data

We will be using the "polarity dataset v2.0", constructed by Pang and Lee (http://www.cs.cornell.edu/People/pabo/movie%2Dreview%2Ddata/). It has the full text from 2000 movies reivews: 1000 reviews are classified as "positive" and 1000 as "negative." Our goal is to predict whether a review has positive or negative sentiment from the text of the review. Each review is stored in a separate file: the positive reviews are in a folder called "pos", and the negative reviews are in "neg". We have provided some code in load.py to assist with reading these files. You can use the code, or write your own version. The code removes some special symbols from the reviews. Later you can check if this helps or hurts your results.

1. Load all the data and randomly split it into 1500 training examples and 500 test examples.

## 5 Sparse Representations

The most basic way to represent text documents for machine learning is with a "bag-of-words" representation. Here every possible word is a feature, and the value of a word feature is the number of times that word appears in the document. Of course, most words will not appear in any particular document, and those counts will be zero. Rather than store a huge number of zeros, we use a sparse representation, in which we only store the counts that are nonzero. The counts are stored in a key/value store (such as a dictionary in Python). For example, "Harry Potter and Harry Potter II" would be represented as the following Python dict:  $x=\{\text{`Harry':2, `Potter':2, `and':1, 'II':1}\}$ . We will be using linear classifiers of the form  $f(x)=w^Tx$ , and we can store the w vector in a sparse format as well, such as  $w=\{\text{`minimal':1.3,`Harry':-1.1,`viable':-4.2,`and':2.2,`product':9.1}\}$ . The inner product between w and x would only involve the features that appear in both x and y since whatever doesn't appear is assumed to be zero. For this example, the inner product would be x[Harry] \* w[Harry] + x[and] \* w[and] = 2\*(-1.1) + 1\*(2.2). Although we hate to spoil the fun, to help you along, we've included two functions for working with sparse vectors: 1) a dot product

between two vectors represented as dict's and 2) a function that increments one sparse vector by a scaled multiple of another vector, which is a very common operation. These functions are located in util.py.

1. Write a function that converts an example (e.g. a list of words) into a sparse bag-of-words representation. You may find Python's Counter class to be useful here: https://docs.python.org/2/library/collections.html. Note that a Counter is also a dict.

#### Solution:

```
class review_instance:
    """ Build the review class """
   def __init__(self):
        self.word_list = []
        self.label = 0
        self.word_dict = Counter()
        self.word_tfidf = Counter()
    def set_label(self, label_val):
        self.label = label_val
    def read_file(self, file_name):
        Read each file into a list of strings. And set the list to self.
           word_list
        f = open(file_name)
        lines = f.read().split(' ')
        symbols = '${}()[].,:;+-*/&|<>=~" '
       words = map(lambda Element: Element.translate(None, symbols).strip()
          , lines)
        words = filter(None, words)
        self.word_list = words
        self.construct_word_dict(stop_word=stopwords.words('english'))
    def construct_word_dict(self, stop_word=None, count_words=None):
        count the words in word_list, transform to a dict of (word:
           word_count)
        :param stop_word: a list of stop words, The words you hope to
           filter, not included in dict, default set to None
        :param count_words: a list, the words you hope to keep in the
           count_dict, if set to None, will keep all the words
        :return:
        ....
        c = Counter()
        c = Counter(self.word_list)
```

```
if count_words:
    c = Counter({i: c[i] for i in count_words})

if stop_word:
    for i in stop_word:
        del c[i]
    self.word_dict = c

def transform_to_tfidf(self, idf_dict):
    """

    Take a idf dict as input, construct the {word: tfidf} vector tfidf = tf*(idf+1)
    """

for word in self.word_dict:
    self.word_tfidf[word] = self.word_dict.get(word) * idf_dict.get (word, 1)
```

2. Write a version of generic\_gradient\_checker from Homework 1 that works with sparse vectors represented as dict types. See Homework 1 solutions if you didn't do that part. Since we'll be using it for stochastic methods, it should take a single (x, y) pair, rather than the entire dataset. Be sure to use the dotProduct and increment primitives we provide, or make your own.

Solution:

```
def gradient_checker_for_pegasos(review_X, review_y, weight, reg_lambda,
   objective_func=pegasos_sgd_loss, gradient_func=pegasos_sgd_gradient,
   epsilon=0.01, tolerance=1e-4):
   true_gradient = gradient_func(review_X, review_y, weight, reg_lambda)
    weight_to_check = weight.copy()
    #To make weight to check have all the words appear in w and review
       stored in dict
    increment(weight_to_check, 0, review_X)
    approx_grad = weight_to_check.copy()
    for word in weight_to_check:
        weight_plus = weight_to_check.copy()
        weight_plus[word] += epsilon
        loss_plus = objective_func(review_X, review_y, weight_plus,
           reg_lambda)
        weight_minus = weight_to_check.copy()
        weight_minus[word] -= epsilon
        loss_minus = objective_func(review_X, review_y, weight_minus,
           reg_lambda)
        approx_grad[word] = (loss_plus - loss_minus)/(2*epsilon)
    distance = np.sum([(true_gradient[i] - approx_grad[i])**2 for i in
        approx_grad])
    return distance < tolerance</pre>
```

## 6 Support Vector Machine via Pegasos

In this question you will build an SVM using the Pegasos algorithm. To align with the notation used in the Pegasos paper<sup>2</sup>, we're considering the following formulation of the SVM objective function:

$$\min_{w \in \mathbf{R}^n} \frac{\lambda}{2} ||w||^2 + \frac{1}{m} \sum_{i=1}^m \max \{0, 1 - y_i w^T x_i\}.$$

Note that, for simplicity, we are leaving off the unregularized bias term b, and the expression with "max" is just another way to write  $(1 - y_i w^T x)_+$ . Pegasos is stochastic subgradient descent using a step size rule  $\eta_t = 1/(\lambda t)$ . The pseudocode is given below:

```
Input: \lambda > 0. Choose w_1 = 0, t = 0

While epoch \leq \max_{-} = 0

For j = 1, ..., m (assumes data is randomly permuted)

t = t + 1

\eta_t = 1/(t\lambda);

If y_j w_t^T x_j < 1

w_{t+1} = (1 - \eta_t \lambda) w_t + \eta_t y_j x_j

Else

w_{t+1} = (1 - \eta_t \lambda) w_t
```

1. [Written] Compute a subgradient for the "stochastic" SVM objective, which assumes a single training point. Show that if your step size rule is  $\eta_t = 1/(\lambda t)$ , then the the corresponding SGD update is the same as given in the pseudocode. [You may use the following facts without proof: 1) If  $f_1, \ldots, f_m : \mathbf{R}^d \to \mathbf{R}$  are convex functions and  $f = f_1 + \cdots + f_m$ , then  $\partial f(x) = \partial f_1(x) + \cdots + \partial f_m(x)$ . 2) For  $\alpha \geq 0$ ,  $\partial (\alpha f)(x) = \alpha \partial f(x)$ . ]

#### Solution:

We need to find a subgradient of  $\frac{\lambda}{2}||w||^2 + \max\{0, 1 - y_i w^T x_i\}$ . Using the rules provided and the calculation in the first problem, a subgradient at w is given by

$$g = \begin{cases} \lambda w - y_i x_i & \text{for } y_i w^T x_i < 1\\ \lambda w & \text{for } y_i w^T x_i \ge 1. \end{cases}$$

So a stochastic subgradient descent update would be

$$w \leftarrow w - \eta g$$
.

If  $y_I w^T x_i < 1$ , the updated w value is

$$w - \eta (\lambda w - y_i x_i) = (1 - \eta \lambda) w + \eta y_i x_i,$$

and otherwise is simply

$$w - \eta \lambda w = w (1 - \eta \lambda),$$

which are the updates given in the algorithm.

<sup>&</sup>lt;sup>2</sup>Shalev-Shwartz et al.'s "Pegasos: Primal Estimated sub-GrAdient SOlver for SVM" http://ttic.uchicago.edu/~nati/Publications/PegasosMPB.pdf

2. Implement the Pegasos algorithm to run on a sparse data representation. The output should be a sparse weight vector w. Note that our Pegasos algorithm starts at w = 0. In a sparse representation, this corresponds to an empty dictionary. Solution:

```
def pegasos_sgd_loss(review_X, review_y, w, reg_lambda):
    reg_term = np.sum([(w[i])**2 for i in w])*reg_lambda/2.0
    hinge_term = max(0, 1-review_y*dotProduct(w, review_X))
    return reg_term+hinge_term

def pegasos_sgd_gradient(review_X, review_y, w, reg_lambda):
    weight_grad = w.copy()
    for key in weight_grad:
        weight_grad[key] *= reg_lambda
    if review_y*dotProduct(w, review_X)<1:
        increment(weight_grad, -review_y, review_X)
    else:
        pass
    return weight_grad</pre>
```

3. Note that in every step of the Pegasos algorithm, we rescale every entry of  $w_t$  by the factor  $(1 - \eta_t \lambda)$ . Implementing this directly with dictionaries is relatively slow. We can make things significantly faster by representing w as w = sW, where  $s \in \mathbf{R}$  and  $W \in \mathbf{R}^d$ . You can start with s = 1 and W all zeros (i.e. an empty dictionary). Note that both updates start with rescaling  $w_t$ , which we can do simply by setting  $s_{t+1} = (1 - \eta_t \lambda) s_t$ . If the update is  $w_{t+1} = (1 - \eta_t \lambda) w_t + \eta_t y_j x_j$ , then **verify that the update is equivalent to**:

$$s_{t+1} = (1 - \eta_t \lambda) s_t$$

$$W_{t+1} = W_t + \frac{1}{s_{t+1}} \eta_t y_j x_j.$$

There is one subtle issue with the approach described above: if we ever have  $1 - \eta_t \lambda = 0$ , then  $s_{t+1} = 0$ , and we'll have a divide by 0 in the calculation for  $W_{t+1}$ . This only happens when  $\eta_t = 1/\lambda$ . With our step-size rule of  $\eta_t = 1/(\lambda t)$ , it happens exactly when t = 1. So one approach is to just start at t = 2. More generically, note that if  $s_{t+1} = 0$ , then  $w_{t+1} = 0$ . Thus an equivalent representation is  $s_{t+1} = 1$  and W = 0. Thus if we ever get  $s_{t+1} = 0$ , simply set it back to 1 and reset  $W_{t+1}$  to zero, which is an empty dictionary in a sparse representation. Implement the Pegasos algorithm with this change. [Section 5.1 of http://leon.bottou.org/papers/bottou-tricks-2012 for a more generic version of this trick, and many other useful tricks.]

```
def pegasos_fast(review_list, max_epoch, lam, watch_list=None, tfidf= False
):
    """
A faster version of pegasos tfidf: whether use tfidf features
```

```
if lam < 0:
    sys.exit("Lam must be greater than 0")
print "lambda = %r, use tfidf = %r" %(lam, tfidf)
#Initialization
weight = Counter()
epoch = 0
t = 1.
review_number = len(review_list)
s = 1.
while epoch < max_epoch:</pre>
    start_time = time.time()
    epoch += 1
    for j in xrange(review_number):
        t += 1
        eta = 1./(t*lam)
        s = (1 - eta*lam)*s
        review = review_list[j]
        label = review.label
        if tfidf:
            feature = review.word_tfidf
        else:
            feature = review.word_dict
        if label*dotProduct(feature, weight) < 1:</pre>
            temp = eta*label/s
            increment (weight, temp, feature)
            #print weight
    end_time = time.time()
    epoch weight = Counter()
    increment (epoch_weight, s, weight)
    print "Epoch %r: in %.3f seconds. Training accuracy: %.3f, Test
        accuracy: %.3f" \ %(epoch, end_time-start_time,
        accuracy_percent(review_list, epoch_weight, tfidf=tfidf)
        accuracy_percent(watch_list, epoch_weight, tfidf=tfidf))
return epoch_weight
```

- 4. Run both implementations of Pegasos to train an SVM on the training data (using the bag-of-words feature representation described above). Make sure your implementations are correct by verifying that the two approaches give essentially the same result. Report on the time taken to run each approach.
- 5. Write a function that takes a sparse weight vector w and a collection of (x, y) pairs, and returns the percent error when predicting y using  $\operatorname{sign}(w^T x)$ . In other words, the function reports the 0-1 loss of the linear predictor  $x \mapsto w^T x$ . Solution:

```
def svm_predict(review_X, weight):
```

6. Using the bag-of-words feature representation described above, search for the regularization parameter that gives the minimal percent error on your test set. A good search strategy is to start with a set of lambdas spanning a broad range of orders of magnitude. Then, continue to zoom in until you're convinced that additional search will not significantly improve your test performance. Once you have a sense of the general range of regularization parameters that give good results, you do not have to search over orders of magnitude every time you change something (such as adding new feature).

Solution:

```
def stepsize_search(X, y, theta, loss_func, grad_func, epsilon=1e-6):
    alpha = 1.0
    gamma = 0.5
    loss = loss_func(X, y, theta)
    gradient = grad_func(X, y, theta)
    while True:
        theta_next = theta - alpha*grad_func(X,y, theta)
        loss_next = loss_func(X, y, theta_next)
        if loss_next > loss-epsilon:
            alpha = alpha*gamma
        else:
            return alpha
```

7. [Optional] Recall that the "score" is the value of the prediction  $f(x) = w^T x$ . We like to think that the magnitude of the score represents the confidence of the prediction. This is something we can directly verify or refute. Break the predictions into groups based on the score (you can play with the size of the groups to get a result you think is informative). For each group, examine the percentage error. You can make a table or graph. Summarize the results. Is there a correlation between higher magnitude scores and accuracy?

## 7 Error Analysis

The natural language processing domain is particularly nice in that often one can often interpret why a model has performed well or poorly on a specific example, and sometimes it is not very difficult to come up with ideas for new features that might help fix a problem. The first step in this process is to look closely at the errors that our model makes.

1. Choose some examples that the model got wrong. List the features that contributed most heavily to the descision (e.g. rank them by  $|w_ix_i|$ ), along with  $x_i, w_i, xw_i$ . Do you understand why the model was incorrect? Can you think of a new feature that might be able to fix the issue? Include a short analysis for at least 2 incorrect examples.

### 8 Features

For a problem like this, the features you use are far more important than the learning model you choose. Whenever you enter a new problem domain, one of your first orders of business is to beg, borrow, or steal the best features you can find. This means looking at any relevant published work and seeing what they've used. Maybe it means asking a colleague what features they use. But eventually you'll need to engineer new features that help in your particular situation. To get ideas for this dataset, you might check the discussion board on this Kaggle competition, which is using a very similar dataset https://www.kaggle.com/c/sentiment-analysis-on-movie-reviews. There are also a very large number of academic research papers on sentiment analysis that you can look at for ideas.

- 1. Based on your error analysis, or on some idea you have, construct a new feature (or group of features) that you hope will improve your test performance. Describe the features and what kind of improvement they give. At this point, it's important to consider the standard errors  $(\sqrt{p(1-p)/n})$  on your performance estimates, to know whether the improvement is statistically significant.
- 2. [Optional] Try to get the best performance possible by generating lots of new features, changing the pre-processing, or any other method you want, so long as you are using the same core SVM model. Describe what you tried, and how much improvement each thing brought to the model. To get you thinking on features, here are some basic ideas of varying quality: 1) how many words are in the review? 2) How many "negative" words are there? (You'd have to construct or find a list of negative words.) 3) Word n-gram features: Instead of single-word features, you can make every pair of consecutive words a feature. 4) Character n-gram features: Ignore word boundaries and make every sequence of n characters into a feature (this will be a lot). 5) Adding an extra feature whenever a word is preceded by "not". For example "not amazing" becomes its own feature. 6) Do we really need to eliminate those funny characters in the data loading phase? Might there be useful signal there? 7) Use tf-idf instead of raw word counts. The tf-idf is calculated as

$$tfidf(f_i) = \frac{FF_i}{\log(DF_i)} \tag{1}$$

where  $FF_i$  is the feature frequency of feature  $f_i$  and  $DF_i$  is the number of document containing  $f_i$ . In this way we increase the weight of rare words. Sometimes this scheme helps,

sometimes it makes things worse. You could try using both! [Extra credit points will be awarded in proportion to how much improvement you achieve.] Solution:

- 1. Length of the text
- 2. N-gram (N;1)
- 3. Word Vector: See Mikolov's word2vec(https://code.google.com/p/word2vec/) or Stanford's GloVe(http://nlp.stanford.edu/projects/glove/). For sentences/documents, see Quoc Le's paragaph vector(http://cs.stanford.edu/~quocle/paragraph\_vector. pdf)
- 4. Topic-Model features: eg LDA
- 5. ...