CIS 519 Homework 2: Linear Classifiers

Handed Out: October 5, 2020

Due: October 19, 2020 at 11:59pm.

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Although the solutions are my own, I consulted with the following people while working on this homework:

• TODO (if applicable): ...

Preface

- Feel free to talk to other members of the class in doing the homework. I am more concerned
 that you learn how to solve the problem than that you demonstrate that you solved it entirely
 on your own. You should, however, write down your solution yourself. Please include here
 the list of people you consulted with in the course of working on the homework:
- While we encourage discussion within and outside the class, cheating and copying code is strictly not allowed. Copied code will result in the entire assignment being discarded at the very least.
- Please use Piazza if you have questions about the homework. Also, please come to the TAs
 recitations and to the office hours.
- The homework is due at 11:59 PM on the due date. We will be using Gradescope for
 collecting the homework assignments. You should have been automatically added to
 Gradescope. If not, please ask a TA for assistance. Post on Piazza and contact the TAs if you
 are having technical difficulties in submitting the assignment.
- Here are some resources you will need for this assignment
 (https://www.seas.upenn.edu/~cis519/fall2020/assets/HW/HW2/hw2-materials.zip
 (https://www.seas.upenn.edu/~cis519/fall2020/assets/HW/HW2/hw2-materials.zip))

Overview

About Jupyter Notebooks

In this homework assignment, we will use a Jupyter notebook to implement, analyze, and discuss ML classifiers. Knowing and being comfortable with Jupyter notebooks is a must in every data scientist, ML engineer, researcher, etc. They are widely used in industry and are a standard form of communication in ML by intertwining text and code to "tell a story". There are many resources that can introduce you to Jupyter notebooks (they are pretty easy to understand!), and if you still need help any of the TAs are more than willing to help.

We will be using a local instance of Jupyter instead of Colab. You are of course free to use Colab, but you will need to understand how to hook your Colab instance with Google Drive to upload the datasets and to save images.

About the Homework

You will experiment with several different linear classifiers and analyze their performances in both real and synthetic datasets. The goal is to understand the differences and similarities between the algorithms and the impact that the dataset characteristics have on the algorithms' learning behaviors and performances.

In total, there are seven different learning algorithms which you will implement. Six are variants of the Perceptron algorithm and the seventh is a support vector machine (SVM). The details of these models is described in Section 1.

In order to evaluate the performances of these models, you will use several different datasets. The first two datasets are synthetic datasets that have features and labels that were programatically generated. They were generated using the same script but use different input parameters that produced sparse and dense variants. The second two datasets are for the task of named-entity recognition (NER), identifying the names of people, locations, and organizations within text. One comes from news text and the other from a corpus of emails. For these two datasets, you need to implement the feature extraction yourself. All of the datasets and feature extraction information are described in Section 2.

Finally, you will run two sets of experiments, one on the synthetic data and one on the NER data. The first set will analyze how the amount of training data impacts model performance. The second will look at the consequences of having training and testing data that come from different domains. The details of the experiments are described in Section 3.

Distribution of Points

The homework has 4 sections for a total of 100 points + 10 extra credit points:

- Section 0: Warmup (5 points)
- Section 1: Linear Classifiers (30 points)
- Section 2: Datasets (0 points, just text)
- Section 3: Experiments (65 points)
 - Synthetic Experiment:
 - Parameter Tuning (10 points)
 - Learning Curves(10 points)
 - Final Test Accuracies (5 points)
 - Discussion Questions (5 points)
 - Noise Experiment (10 points extra credit)
 - NER Experiment:
 - Feature Extraction (25 points)
 - Final Test Accuracies (5 points)
 - F₁ Discussion Questions (5 points)

Section 0: Warmup

Only For Colab

If you want to complete this homework in Colab, you are more than welcome to. You will need a little bit more maneuvering since you will need to upload the files of hw2 to your Google Drive and run the following two cells:

```
In [1]:  # Uncomment if you want to use Colab for this homework.
# from google.colab import drive
# drive.mount('/content/drive', force_remount=True)
In [2]:  # Uncomment if you want to use Colab for this homework.
# %cd /content/drive/My Drive/Colab Notebooks/YOUR PATH TO HW FOLDER
```

Python Version

Python 3.6 or above is required for this homework. Make sure you have it installed.

```
In [3]: # Let's check.
import sys
if sys.version_info[:2] < (3, 6):
    raise Exception("You have Python version " + str(sys.version_info))</pre>
```

Imports and Helper Functions (5 points total)

Let's import useful modules we will need throughout the homework as well as implement helper functions for our experiment. **Read and remember** what each function is doing, as you will probably need some of them down the line.

```
In [6]:
            Helper function that loads a synthetic dataset from the dataset root (e.g. "s
            You should not need to edit this method.
            def load_synthetic_data(dataset_type):
                def load jsonl(file path):
                    data = []
                    with open(file_path, 'r') as f:
                        for line in f:
                            data.append(json.loads(line))
                    return data
                def load txt(file path):
                    data = []
                    with open(file path, 'r') as f:
                        for line in f:
                            data.append(int(line.strip()))
                    return data
                def convert_to_sparse(X):
                    sparse = []
                    for x in X:
                        data = \{\}
                        for i, value in enumerate(x):
                            if value != 0:
                                data[str(i)] = value
                        sparse.append(data)
                    return sparse
                path = os.path.join(SYNTHETIC_PATH, dataset_type)
                X_train = load_jsonl(os.path.join(path, 'train.X'))
                X_dev = load_jsonl(os.path.join(path, 'dev.X'))
                X test = load jsonl(os.path.join(path, 'test.X'))
                num_features = len(X_train[0])
                features = [str(i) for i in range(num features)]
                X_train = convert_to_sparse(X_train)
                X_dev = convert_to_sparse(X_dev)
                X_test = convert_to_sparse(X_test)
                y train = load txt(os.path.join(path, 'train.y'))
                y_dev = load_txt(os.path.join(path, 'dev.y'))
                y_test = load_txt(os.path.join(path, 'test.y'))
                return X_train, y_train, X_dev, y_dev, X_test, y_test, features
```

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```
In [7]:
            Helper function that loads the NER data from a path (e.g. "ner/conll/train").
            You should not need to edit this method.
            def load_ner_data(dataset=None, dataset_type=None):
                # List of tuples for each sentence
                data = []
                path = os.path.join(os.path.join(NER_PATH, dataset), dataset_type)
                for filename in os.listdir(path):
                    with open(os.path.join(path, filename), 'r') as file:
                        sentence = []
                        for line in file:
                            if line == '\n':
                                data.append(sentence)
                                sentence = []
                            else:
                                sentence.append(tuple(line.split()))
                return data
```

```
In [8]:
            A helper function that plots the relationship between number of examples
            and accuracies for all the models.
            You should not need to edit this method.
            def plot learning curves(
                perceptron accs,
                winnow accs,
                adagrad_accs,
                avg perceptron accs,
                avg_winnow_accs,
                avg_adagrad_accs,
                svm accs,
                save dest = 'NA'
            ):
                This function will plot the learning curve for the 7 different models.
                Pass the accuracies as lists of length 11 where each item corresponds
                to a point on the learning curve.
                accuracies = [
                    ('perceptron', perceptron_accs),
                    ('winnow', winnow_accs),
                    ('adagrad', adagrad_accs),
                    ('avg-perceptron', avg perceptron accs),
                    ('avg-winnow', avg_winnow_accs),
                    ('avg-adagrad', avg_adagrad_accs),
                    ('svm', svm accs)
                1
                x = [500, 1000, 1500, 2000, 2500, 3000, 3500, 4000, 4500, 5000, 10000]
                plt.figure()
                f, (ax, ax2) = plt.subplots(1, 2, sharey=True, facecolor='w')
                for label, acc list in accuracies:
                    assert len(acc list) == 11
                    ax.plot(x, acc list, label=label)
                    ax2.plot(x, acc_list, label=label)
                ax.set xlim(0, 5500)
                ax2.set xlim(9500, 10000)
                ax2.set xticks([10000])
                # hide the spines between ax and ax2
                ax.spines['right'].set_visible(False)
                ax2.spines['left'].set_visible(False)
                ax.yaxis.tick left()
                ax.tick params(labelright='off')
                ax2.yaxis.tick right()
                ax2.legend()
                if (save_dest == 'sparse'):
                    plt.savefig('images/sparse_final.png')
                elif (save dest == 'dense'):
```

```
plt.savefig('images/dense_final.png')
plt.show()
```

F1 Score (5 points)

For some part of the homework, you will use the F1 score instead of accuracy to evaluate how well a model does. The F1 score is computed as the harmonic mean of the precision and recall of the classifier. Precision measures the number of correctly identified positive results by the total number of positive results. Recall, on the other hand, measures the number of correctly identified positive results divided by the number of all samples that should have been identified as positive. More formally, we have that

Precision =
$$\frac{TP}{TP + FP}$$
Recall =
$$\frac{TP}{TP + FN}$$

where TP is the number of true positives, FP false positives and FN false negatives. Combining these two we define F1 as

$$F1 = 2 \cdot \frac{Precision \cdot Recall}{Precision + Recall}$$

You now need to implement the calculation of F1 yourself using the provided function header. It will be unit tested on Gradescope.

```
In [9]:

    def calculate_f1(y_gold, y_model):

                tp count = 0
                fp count = 0
                fn count = 0
                for i in range (0, len(y_gold)):
                     if(y model[i] == 1):
                         if(y gold[i] == 1):
                             tp_count+=1
                         if(y_gold[i] != 1):
                             fp count+=1
                     else:
                         if(y_gold[i] == 1):
                             fn count+=1
                precision = tp_count/(tp_count+fp_count)
                recall = tp_count/(tp_count+fn_count)
                f1 = 2*(precision*recall)/(precision+recall)
                return f1
```

Looking at the formula for the F1 score, what is the highest and lowest possible value?

Section 1. Linear Classifiers (30 points total)

This section details the seven different algorithms that you will use in the experiments. For each of the algorithms, we describe the initialization you should use to start training and the different parameter settings that you should use for the experiment on the synthetic datasets. Each of the update functions for the Perceptron, Winnow, and Perceptron with AdaGrad will be unittested on Gradescope, so please do not edit the function definitions.

1.1 Base Classifier

```
In [11]:
         DO NOT MODIFY
                 The Classifier class is the base class for all of the Perceptron-based
                 algorithms. Your class should override the "process_example" and
                 "predict single" functions. Further, the averaged models should
                 override the "finalize" method, where the final parameter values
                 should be calculated.
                 You should not need to edit this class any further.
                 ITERATIONS = 10
                 def train(self, X, y):
                     for iteration in range(self.ITERATIONS):
                         for x_i, y_i in zip(X, y):
                             self.process_example(x_i, y_i)
                     self.finalize()
                 def process_example(self, x, y):
                     Makes a prediction using the current parameter values for
                     the features x and potentially updates the parameters based
                     on the gradient. Note "x" is a dictionary which maps from the feature
                     name to the feature value and y is either 1 or -1.
                     raise NotImplementedError
                 def finalize(self):
                     """Calculates the final parameter values for the averaged models."""
                 def predict(self, X):
                     Predicts labels for all of the input examples. You should not need
                     to override this method.
                     return [self.predict_single(x) for x in X]
                 def predict_single(self, x):
                     Predicts a label, 1 or -1, for the input example. "x" is a dictionary
                     which maps from the feature name to the feature value.
                     raise NotImplementedError
```

1.2 Basic Perceptron (2 points)

We do this classifier for you, so enjoy the two free points and pay attention to the techniques and code written.

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This is the basic version of the Perceptron Algorithm. In this version, an update will be performed on the example (\mathbf{x}, y) if $y(\mathbf{w}^{\mathsf{T}}\mathbf{x} + \theta) \leq 0$. The Perceptron needs to learn both the bias term θ and the weight vector \mathbf{w} parameters. When the Perceptron makes a mistake on the example (\mathbf{x}, y) , both \mathbf{w} and θ need to be updated using the following update equations:

$$\mathbf{w}^{\text{new}} \leftarrow \mathbf{w} + \boldsymbol{\eta} \cdot \boldsymbol{y} \cdot \mathbf{x}$$
$$\theta^{\text{new}} \leftarrow \theta + \boldsymbol{\eta} \cdot \boldsymbol{y}$$

where η is the learning rate.

1.2.2 Hyperparameters

We set η to 1, so there are no hyperparameters to tune.

Note: If we assume that the order of the examples presented to the algorithm is fixed, we initialize $\mathbf{w}=\mathbf{0}$ and $\theta=0$, and train both together, then the learning rate η does not have any effect. In fact you can show that, if \mathbf{w}_1 and θ_1 are the outputs of the Perceptron algorithm with learning rate η_1 , then \mathbf{w}_1/η_1 and θ_1/η_1 will be the result of the Perceptron with learning rate 1 (note that these two hyperplanes give identical predictions).

1.2.3 Initialization

$$\mathbf{w} = [0, 0, ..., 0]$$
 and $\theta = 0$.

```
In [12]:
         DO NOT MODIFY THIS CELL
                 The Perceptron model. Note how we are subclassing `Classifier`.
                def __init__(self, features):
                    Initializes the parameters for the Perceptron model. "features"
                    is a list of all of the features of the model where each is
                    represented by a string.
                    # NOTE: Do not change the names of these 3 data members because
                    # they are used in the unit tests
                    self.eta = 1
                    self.theta = 0
                    self.w = {feature: 0.0 for feature in features}
                 def process example(self, x, y):
                    y_pred = self.predict_single(x)
                    if y != y pred:
                        for feature, value in x.items():
                            self.w[feature] += self.eta * y * value
                        self.theta += self.eta * y
                 def predict_single(self, x):
                    score = 0
                    for feature, value in x.items():
                        score += self.w[feature] * value
                    score += self.theta
                    if score <= 0:
                        return -1
                    return 1
```

For the rest of the Perceptron-based algorithms, you will have to implement the corresponding class like we have done for Perceptron . Use the Perceptron class as a guide for how to implement the functions.

1.3 Winnow (5 points)

1.3.1 Description

The Winnow algorithm is a variant of the Perceptron algorithm with multiplicative updates. Since the algorithm requires that the target function is monotonic, you will only use it on the synthetic datasets.

The Winnow algorithm only learns parameters \mathbf{w} . We will fix $\theta = -n$, where n is the number of features. When the Winnow algorithm makes a mistake on the example (\mathbf{x}, y) , the parameters are updated with the following equation:

$$w_i^{\text{new}} \leftarrow w_i \cdot \alpha^{y \cdot x_i}$$

where w_i and x_i are the *i*th components of the corresponding vectors. Here, α is a promotion/demotion hyperparameter.

1.3.2 Hyperparameters

For the experiment, choose $\alpha \in \{1.1, 1.01, 1.005, 1.0005, 1.0001\}$.

1.3.3 Initialization

```
\mathbf{w} = [1, 1, \dots, 1] and \theta = -n (constant).
```

```
In [13]:
                def init (self, alpha, features):
                    # DO NOT change the names of these 3 data members because
                    # they are used in the unit tests
                    self.alpha = alpha
                    self.w = {feature: 1.0 for feature in features}
                    self.theta = -len(features)
                def process example(self, x, y):
                    y_pred = self.predict_single(x)
                    if (y pred != y):
                        for feature, value in x.items():
                            self.w[feature] = self.w[feature] * self.alpha**(y*value)
                def predict single(self, x):
                    score = 0
                    for feature, value in x.items():
                        score += self.w[feature] * value
                    score += self.theta
                    if (score <= 0):
                        return -1
                    else:
                        return 1
```

1.4 AdaGrad (10 points)

1.4.1 Description

AdaGrad is a variant of the Perceptron algorithm that adapts the learning rate for each parameter based on historical information. The idea is that frequently changing features get smaller learning rates and stable features higher ones.

To derive the update equations for this model, we first need to start with the loss function. Instead of using the hinge loss with the elbow at 0 (like the basic Perceptron does), we will instead use the standard hinge loss with the elbow at 1:

$$\mathcal{L}(\mathbf{x}, y, \mathbf{w}, \theta) = \max\{0, 1 - y(\mathbf{w}^{\mathsf{T}}\mathbf{x} + \theta)\}\$$

Then, by taking the partial derivative of \mathcal{L} with respect to \mathbf{w} and θ , we can derive the respective graidents (make sure you understand how you could derive these gradients on your own):

$$\frac{\partial \mathcal{L}}{\partial \mathbf{w}} = \begin{cases} \mathbf{0} & \text{if } y(\mathbf{w}^{\mathsf{T}} \mathbf{x} + \theta) > 1 \\ -y \cdot \mathbf{x} & \text{otherwise} \end{cases}$$

$$\frac{\partial \mathcal{L}}{\partial \theta} = \begin{cases} 0 & \text{if } y(\mathbf{w}^{\mathsf{T}} \mathbf{x} + \theta) > 1 \\ -y & \text{otherwise} \end{cases}$$

Then for each parameter, we will keep track of the sum of the parameters' squared gradients. In the following equations, the k superscript refers to the kth non-zero gradient (i.e., the kth weight vector/misclassified example) and t is the number of mistakes seen thus far.

$$G_{j}^{t} = \sum_{k=1}^{t} \left(\frac{\partial \mathcal{L}}{\partial w_{j}^{k}}\right)^{2}$$
$$H^{t} = \sum_{k=1}^{t} \left(\frac{\partial \mathcal{L}}{\partial \theta^{k}}\right)^{2}$$

For example, on the 3rd mistake (t=3), G_j^3 is the sum of the squares of the first three non-zero gradients $(\left(\frac{\partial \mathcal{L}}{\partial w_j^1}\right)^2, \left(\frac{\partial \mathcal{L}}{\partial w_j^2}\right)^2$, and $\left(\frac{\partial \mathcal{L}}{\partial w_j^3}\right)^2$). Then \mathbf{G}^3 is used to calculate the 4th value of the weight vector as follows. On example (\mathbf{x},y) , if $y(\mathbf{w}^{\mathsf{T}}\mathbf{x}+\theta)\leq 1$, then the parameters are updated with the following equations:

$$\mathbf{w}^{t+1} \leftarrow \mathbf{w}^{t} + \eta \cdot \frac{y \cdot \mathbf{x}}{\sqrt{\mathbf{G}^{t}}}$$
$$\theta^{t+1} \leftarrow \theta^{t} + \eta \frac{y}{\sqrt{H^{t}}}$$

Note that, although we use the hinge loss with the elbow at 1 for training, you still make the prediction based on whether or not $y(\mathbf{w}^{\mathsf{T}}\mathbf{x} + \theta) \leq 0$ during testing.

1.4.2 Hyperparameters

For the experiment, choose $\eta \in \{1.5, 0.25, 0.03, 0.005, 0.001\}$

1.4.3 Initialization

$$\mathbf{w} = [0, 0, \dots, 0] \text{ and } \theta = 0.$$

```
In [14]:
                 def __init__(self, eta, features):
                     # DO NOT change the names of these 3 data members because
                     # they are used in the unit tests
                     self.eta = eta
                     self.w = {feature: 0.0 for feature in features}
                     self.theta = 0
                     self.G = {feature: 1e-5 for feature in features} # 1e-5 prevents div
                     self.H = 0
                 def process example(self, x, y):
                     score = self.predict_single_train(x)
                     if (score*y <= 1):
                         for feature, value in x.items():
                             dw = -y*value
                             self.G[feature]+=(dw**2)
                             self.w[feature] = self.w[feature] + (self.eta*y*value)/(np.sq
                        dt = -y
                         self.H += dt**2
                         self.theta = self.theta + (self.eta*y)/(np.sqrt(self.H))
                 def predict_single_train(self, x):
                     score = 0
                     for feature, value in x.items():
                         score += self.w[feature] * value
                     score += self.theta
                     return score
                 def predict_single(self, x):
                     score = 0
                     for feature, value in x.items():
                         score += self.w[feature] * value
                     score += self.theta
                     if (score <= 0):
                         return -1
                     else:
                         return 1
```

1.5 Averaged Models (15 points)

You will also implement the averaged version of the previous three algorithms.

During the course of training, each of the above algorithms will have K+1 different parameter settings for the K different updates it will make during training. The regular implementation of these algorithms uses the parameter values after the Kth update as the final ones. Instead, the averaged version use the weighted average of the K+1 parameter values as the final parameter values. Let m_k denote the number of correctly classified examples by the kth parameter values and M the total number of correctly classified examples. The final parameter values are

$$M = \sum_{k=1}^{K+1} m_k$$

$$\mathbf{w} \leftarrow \frac{1}{M} \sum_{k=1}^{K+1} m_k \cdot \mathbf{w}^k$$

$$\theta \leftarrow \frac{1}{M} \sum_{k=1}^{K+1} m_k \cdot \theta^k$$

For each of the averaged versions of Perceptron, Winnow, and AdaGrad, use the same hyperparameters and initialization as before.

1.5.1 Implementation Note

Implementing the averaged variants of these algorithms can be tricky. While the final parameter values are based on the sum of K different vectors, there is no need to maintain all of these parameters. Instead, you should implement these algorithms by keeping only two vectors, one which maintains the cumulative sum and the current one.

Additionally, there are two ways of keeping track of these two vectors. One is more straightforward but prohibitively slow. The second requires some algebra to derive but is significantly faster to run. Try to analyze how the final weight vector is a function of the intermediate updates and their corresponding weights. It should take less than a minute or two for ten iterations for any of the averaged algorithms. You need to think about how to efficiently implement the averaged algorithms yourself.

Further, the implementation for Winnow is slightly more complicated than the other two, so if you consistently have low accuracy for the averaged Winnow, take a closer look at the derivation.

```
In [15]:
         def __init__(self, features):
                     self.eta = 1
                     self.w = {feature: 0.0 for feature in features}
                     self.theta = 0
                     self.cumul_w = {feature: 0.0 for feature in features}
                     self.cumul theta = 0
                     self.classified = 0
                     self.final m = 0
                     self.update_x = [feature for feature in features]
                     self.last update = {feature: 0 for feature in features}
                 def process_example(self, x, y):
                     self.final m+=1
                     y pred = self.predict single(x)
                     if (y pred == y):
                         self.classified+=1
                     else:
                         for feature in self.update x:
                             self.cumul_w[feature] += self.w[feature] * (self.final_m - 1
                             self.last update[feature] = self.final m-1
                        for feature in x:
                             self.w[feature] += self.eta * y * x[feature]
                         self.cumul_theta += self.theta * self.classified
                         self.theta += self.eta * y
                         self.classified=1
                         self.update_x = [feature for feature in x]
                 def predict single(self, x):
                     score = 0
                     for feature, value in x.items():
                         score += self.w[feature] * value
                     score += self.theta
                     if score <= 0:</pre>
                         return -1
                     return 1
                 def finalize(self):
                     for feature in self.cumul w:
                             self.cumul w[feature] += self.w[feature] * (self.final m - se
                     self.cumul theta += self.theta * self.classified
                     final_w = {feature: self.cumul_w[feature]/self.final_m for feature in
                     final theta = self.cumul theta/self.final m
                     self.w = final w
                     self.theta = final theta
```

```
In [16]:
                           def __init__(self, alpha, features):
                                                          self.alpha = alpha
                                                          self.w = {feature: 1.0 for feature in features}
                                                          self.theta = -len(features)
                                                          self.cumul w = {feature: 1.0 for feature in features}
                                                          self.classified = 0
                                                          self.final m = 0
                                                          self.update x = [feature for feature in features]
                                                          self.last_update = {feature: 0 for feature in features}
                                               def process_example(self, x, y):
                                                          self.final m+=1
                                                          y pred = self.predict single(x)
                                                          if (y_pred == y):
                                                                     self.classified+=1
                                                          else:
                                                                     for feature in self.update x:
                                                                                self.cumul_w[feature] += self.w[feature] * (self.final_m - 1
                                                                                self.last_update[feature] = self.final_m-1
                                                                     for feature in x:
                                                                                self.w[feature] = self.w[feature] * self.alpha**(y*x[feature]
                                                                     self.classified=1
                                                                     self.update_x = [feature for feature in x]
                                               def predict_single(self, x):
                                                          score = 0
                                                          for feature, value in x.items():
                                                                      score += self.w[feature] * value
                                                          score += self.theta
                                                          if (score <= 0):
                                                                      return -1
                                                          else:
                                                                     return 1
                                               def finalize(self):
                                                          for feature in self.cumul w:
                                                                                self.cumul w[feature] += self.w[feature] * (self.final m - self.w[feature] * (self.w[feature] m - self.w[feature] * (self.w[feature] m - self.w[feature] * (self.w[feature] m - self.w[feature] m - self.w[feature] * (self.w[feature] m - self.w[feature] 
                                                          final_w = {feature: self.cumul_w[feature]/self.final_m for feature in
                                                          self.w = final w
```

```
In [17]:
         class AveragedAdaGrad(Classifier):
                 def __init__(self, eta, features):
                     self.eta = eta
                     self.w = {feature: 0.0 for feature in features}
                     self.theta = 0
                     self.G = {feature: 1e-5 for feature in features}
                     self.H = 0
                     self.cumul w = {feature: 0.0 for feature in features}
                     self.cumul theta = 0
                     self.classified = 0
                     self.final m = 0
                     self.update_x = [feature for feature in features]
                     self.last_update = {feature: 0 for feature in features}
                 def process_example(self, x, y):
                     self.final m+=1
                     score = self.predict single train(x)
                     if (score*y > 1):
                         self.classified+=1
                     else:
                         for feature in self.update x:
                             self.cumul_w[feature] += self.w[feature] * (self.final_m - 1
                             self.last update[feature] = self.final m-1
                         for feature in x:
                             dw=-y*x[feature]
                             self.G[feature]+=(dw**2)
                             self.w[feature]=self.w[feature] + (self.eta*y*x[feature])/(np
                         dt = -y
                         self.H += dt**2
                         self.cumul_theta += self.theta * self.classified
                         self.theta = self.theta + (self.eta*y)/(np.sqrt(self.H))
                         self.classified=1
                         self.update_x = [feature for feature in x]
                 def predict_single_train(self, x):
                     score = 0
                     for feature, value in x.items():
                          score += self.w[feature] * value
                     score += self.theta
                     return score
                 def predict_single(self, x):
                     score = 0
                     for feature, value in x.items():
                         score += self.w[feature] * value
                     score += self.theta
                     if (score <= 0):
                         return -1
                     else:
                         return 1
```

1.6 Support Vector Machines

Although we have not yet covered SVMs in class, you can still train them using the sklearn library. We will use a soft margin SVM for non-linearly separable data. You should use the sklearn implementation as follows:

```
from sklearn.svm import LinearSVC
classifier = LinearSVC(loss='hinge')
classifier.fit(X, y)
```

sklearn requires a different feature representation than what we use for the Perceptron models. The provided Python template code demonstrates how to convert to the require representation.

Given training samples $S = \{(\mathbf{x}^1, y^1), (\mathbf{x}^2, y^2), \dots, (\mathbf{x}^m, y^m)\}$, the objective for the SVM is the following:

$$\min_{\mathbf{w},b,\xi} \frac{1}{2} ||\mathbf{w}||_2^2 + C \sum_{i=1}^m \xi_i$$

subject to the following constraints:

$$y^{i}(\mathbf{w}^{\mathsf{T}}\mathbf{x}^{i}+b) \geq 1-\xi_{i} \text{ for } i=1,2,\ldots,m$$

 $\xi_{i} \geq 0 \text{ for } i=1,2,\ldots,m$

Section 2. Datasets

In this section, we describe the synthetic and NER datasets that you will use for your experiments. For the NER datasets, there is also an explanation for the features which you need to extract from the data.

2.1 Synthetic Data

2.1.1 Introduction

The synthetic datasets have features and labels which are automatically generated from a python script. Each instance will have n binary features and are labeled according to a l-of-m-of-n Boolean function. Specifically, there is a set of m features such that an example if positive if and

only if at least l of these m features are active. The set of m features is the same for the dataset (i.e., it is not a separate set of m features for each individual instance).

We provide two versions of the synthetic dataset called sparse and dense. For both datasets, we set l=10 and m=20. We set n=200 for the sparse data and n=40 for the dense data. Additionally, we add noise to the data as follows: With probability 0.05 the label assigned by the function is changed and with probability 0.001 each feature value is changed. Consequently, the data is not linearly separable.

We have provided you with three data splits for both sparse and dense with 10,000 training, 2,000 development, and 2,000 testing examples. Section 3 describes the experiments that you need to run on these datasets.

2.1.2 Feature Representation

The features of the synthetic data provided are vectors of 0s and 1s. Storing these large matrices requires lots of memory so we use a sparse representation that stores them as dictionaries instead. For example, the vector [0,1,0,0,0,1] can be stored as $\{"x2": 1,"x6": 1\}$ (using 1-based indexing). We have provided you with the code for parsing and converting the data to this format. You can use these for the all algorithms you develop except the SVM. Since you will be using the implementation of SVM from sklearn, you will need to provide a vector to it. You can use sklearn.feature_extraction.DictVectorizer for converting feature-value dictionaries to vectors.

2.2 NER Data

In addition to the synthetic data, we have provided you two datasets for the task of named-entity recognition (NER). The goal is to identify whether strings in text represent names of people, organizations, or locations. An example instance looks like the following:

```
[PER Wolff] , currently a journalist in [LOC Argentina] , played wit
h
[PER del Bosque] in the final years of the seventies in
[ORG Real Madrid] .
```

In this problem, we will simplify the task to identifying whether a string is named entity or not (that is, you don't have to say which type of entity it is). For each token in the input, we will use the tag I to denote that token is an entity and 0 otherwise. For example, the full tagging for the above instace is as follows:

```
[I Wolff] [0 ,] [0 currently] [a] [0 journalist] [0 in] [I Argentin
a]
    [0 ,] [0 played] [0 with] [I del] [I Bosque] [0 in] [0 the] [0 fina
l]
    [0 years] [0 of] [0 the] [0 seventies] [0 in] [I Real] [I Madrid] .
```

Given a sentence $S=w_1,w_2,\ldots,w_n$, you need to predict the I , 0 tag for each word in the sentence. That is, you will produce the sequence $Y=y_1,y_2,\ldots,y_n$ where $y_i\in\{$ I , 0 $\}$.

2.2.1 Datasets: CoNLL and Enron

We have provided two datasets, the CoNLL dataset which is text from news articles, and Enron, a corpus of emails. The files contain one word and one tag per line. For CoNLL, there are training, development, and testing files, whereas Enron only has a test dataset. There are 14,987 training sentences (204,567 words), 336 development sentences (3,779 words), and 303 testing sentences (3,880 words) in CoNLL. For Enron there are 368 sentences (11,852 words).

Please note that the CoNLL dataset is available only for the purposes of this assignment. It is copyrighted, and you are granted access because you are a Penn student, but please delete it when you are done with the homework.

2.2.2 Feature Extraction

The NER data is provided as raw text, and you are required to extract features for the classifier. In this assignment, we will only consider binary features based on the context of the word that is supposed to be tagged.

Assume that there are V unique words in the dataset and each word has been assigned a unique ID which is a number $\{1, 2, \dots, V\}$. Further, w_{-k} and w_{+k} indicate the kth word before and after the target word. The feature templates that you should use to generate features are as follows:

Template	Number of Features
w_{-3}	V
w_{-2}	V
w_{-1}	V
w_{+1}	V
w_{+2}	V
w_{+3}	V
w_{-1} & w_{-2}	$V \times V$
w_{+1} & w_{+2}	$V \times V$
w_{-1} & w_{+1}	$V \times V$

Each feature template corresponds to a set of features that you will compute (similar to the features you generated in problem 2 from the first homework assignment). The w_{-3} feature template corresponds to V features where the ith feature is 1 if the third word to the left of the target word has ID i. The $w_{-1}\&w_{+1}$ feature template corresponds to $V\times V$ features where there is one feature for every unique pair words. For example, feature $(i-1)\times V+j$ is a binary feature that is 1 if the word 1 to the left of the target has ID i and the first word to the right of the target has ID j. In practice, you will not need to keep track of the feature IDs. Instead, each feature will be given a name such as " $w_{-1}= {\rm the}\&w_{+1}= {\rm cat}$ ".

In total, all of the above feature templates correspond to a very large number of features. However, for each word, there will be exactly 9 features which are active (non-zero), so the feature vector is quite sparse. You will represent this as a dictionary which maps from the feature name to the value. In the provided Python template, we have implemented a couple of the features for you to demonstrate how to compute them and what the naming scheme should look like.

In order to deal with the first two words and the last two words in a sentence, we will add special symbol "SSS" and "EEE" to the vocabulary to represent the words before the first word and the words after the last word. Notice that in the test data you may encounter a word that was not observed in training, and therefore is not in your dictionary. In this case, you cannot generate a feature for it, resulting in less than 7 active features in some of the test examples.

Section 3. Experiments (65 points total)

You will run two sets of experiments, one using the synthetic data and one using the NER data.

3.1 Synthetic Experiment (30 + 10 extra credit points)

This experiment will explore the impact that the amount of training data has on model performance. First, you will do hyperparameter tuning for Winnow and Perceptron with AdaGrad (both standard and averaged versions). Then you will generate learning curves that will plot the size of the training data against the performance. Finally, for each of the models trained on all of the training data, you will find the test score. You should use accuracy to compute the performance of the model.

In summary, the experiment consists of three parts

- 1. Parameter Tuning
- 2. Learning Curves
- 3. Final Evaluation

3.1.1 Parameter Tuning (10 points)

For both the Winnow and Perceptron with AdaGrad (standard and averaged), there are hyperparameters that you need to choose. (The same is true for SVM, but you should only use the default settings.) Similarly to cross-validation from Homework 1, we will estimate how well each model will do on the true test data using the development dataset (we will not run cross-validation), and choose the hyperparameter settings based on these results.

For each hyperparameter value in Section 1, train a model using that value on the training data and compute the accuracy on the development dataset. Each model should be trained for 10 iterations (i.e., 10 passes over the entire dataset).

TODO: Fill in the table with the best hyperparameter values and the corresponding validation accuracies. Repeat this for both the sparse and dense data.

Note:

Hyperparameter accuracies were calculated in the final evaluation section. Code used to generate accuracies is built into the run_synthetic_experiment function. Accuracies are listed in the tables, but code generating them is in section below tables.

Winnow Sweep

α	Sparse	Dense
1.1	.8935	.8995
1.01	.927	.9215
1.005	.9195	.908
1.0005	.563	.8615
1.0001	.5205	.614

Averaged Winnow Sweep

α	Sparse	Dense
1.1	.939	.9445
1.01	.8875	.9315
1.005	.8145	.913
1.0005	.5275	.682
1.0001	.5055	.5465

AdaGrad Sweep

η	Sparse	Dense
1.5	.913	.9445
0.25	.7695	.7825
0.03	.6635	.5795
0.005	.5	.5085
0.001	.4985	.5085

Averaged AdaGrad Sweep

η	Sparse	Dense
1.5	.852	.9445
0.25	.762	.7055
0.03	.643	.5885
0.005	.5055	.5085
0.001	.559	.5085

3.1.2 Learning Curves (10 points)

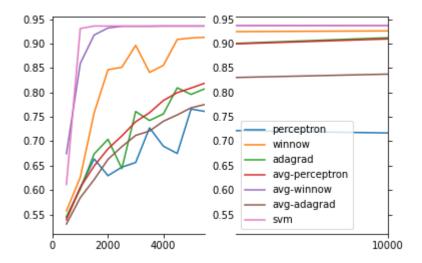
Next, you will train all 7 models with different amounts of training data. For Winnow and Perceptron with AdaGrad (standard and averaged), use the best hyperparameters from the parameter tuning experiment.

Each of the datasets contains 10,000 training examples. You will train each model 11 times on varying amounts of training data. The first 10 will increase by 500 examples: 500, 1k, 1.5k, 2k, ..., 5k. The 11th model should use all 10k examples. Each Perceptron-based model should be trained for 10 iterations (e.g., 10 passes over the total number of training examples available to that model). The SVM can be run until convergence with the default parameters.

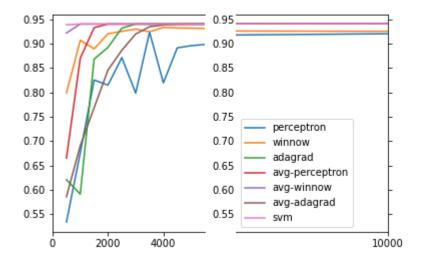
For each model, compute the accuracy on the development dataset and plot the results using the provided code. There should be a separate plot for the sparse and dense data.

Note how we have included an image in markdown. You should do the same for both plots and include them in the output below by running your experiment, saving your plots to the images folder, and linking it to this cell.

Sparse Plot



Dense Plot



3.1.3 Final Evaluation (5 points)

Finally, for each of the 7 models, train the models on all of the training data and compute the test accuracy. For Winnow and Perceptron with AdaGrad, use the best hyperparameter settings you found. Report these accuracies in a table.

We will run our models with [500, 1000, 1500, 2000, 2500, 3000, 3500, 4000, 4500, 5000, 10000] examples.

```
In [26]:
         # Load in data sets
                X_train, y_train, X_dev, y_dev, X_test, y_test, features = load_synthetiq
                # Select model hyperparameters
                winnow hp choices = [1.1, 1.01, 1.005, 1.0005, 1.0001]
                 ada_hp_choices = [1.5, .25, .03, .005, .001]
                winnow_hp_accs = []
                 ada_hp_accs = []
                winnow avg hp accs = []
                 ada_avg_hp_accs = []
                 for hp in winnow hp choices:
                    w = Winnow(hp, features)
                    w.train(X_train, y_train)
                    w pred = w.predict(X dev)
                    w acc = accuracy score(y dev, w pred)
                    winnow_hp_accs.append(w_acc)
                    print('Winnow Accuracy for Hyperparameter: ' + str(hp) + ' = ' + str(
                    w_avg = AveragedWinnow(hp, features)
                    w_avg.train(X_train, y_train)
                    w avg pred = w avg.predict(X dev)
                    w_avg_acc = accuracy_score(y_dev, w_avg_pred)
                    winnow avg hp accs.append(w avg acc)
                    print('AveragedWinnow Accuracy for Hyperparameter: ' + str(hp) + ' =
                    print()
                 for hp in ada hp choices:
                    a = AdaGrad(hp, features)
                    a.train(X_train, y_train)
                    a pred = a.predict(X dev)
                     a_acc = accuracy_score(y_dev, a_pred)
                     ada hp accs.append(a acc)
                    print('AdaGrad Accuracy for Hyperparameter: ' + str(hp) + ' = ' + str
                    a avg = AveragedAdaGrad(hp, features)
                    a_avg.train(X_train, y_train)
                    a_avg_pred = a_avg.predict(X_dev)
                    a avg acc = accuracy score(y dev, a avg pred)
                    ada_avg_hp_accs.append(a_avg_acc)
                    print('AveragedAdaGrad Accuracy for Hyperparameter: ' + str(hp) + ' =
                    print()
                winnow hp selection = winnow hp choices[winnow hp accs.index(max(winnow h
                 ada hp selection = ada hp choices[ada hp accs.index(max(ada hp accs))]
                winnow_avgd_hp_selection = winnow_hp_choices[winnow_avg_hp_accs.index(max
                 ada_avgd_hp_selection = ada_hp_choices[ada_avg_hp_accs.index(max(ada_avg_
                 print('Ideal Winnow Hyperparameter: ' + str(winnow hp selection) + ', Acd
```

```
print('Ideal AdaGrad Hyperparameter: ' + str(ada_hp_selection) + ', Accur
print('Ideal AveragedWinnow Hyperparameter: ' + str(winnow_avgd_hp_select
print('Ideal AveragedAdaGrad Hyperparameter: ' + str(ada_avgd_hp_selection)
print()
# Train models, compute test accuracy
per_accs = []
avg_per_accs = []
winnow accs = []
avg winnow accs = []
ada_accs = []
avg ada accs = []
svc_accs = []
models = {0: Perceptron, 1: AveragedPerceptron,
         2: Winnow, 3: AveragedWinnow,
         4: AdaGrad, 5: AveragedAdaGrad}
model_names = ['Perceptron', 'AveragedPerceptron', 'Winnow', 'AveragedWin
              'AdaGrad', 'AveragedAdaGrad']
accs = [per accs, avg per accs, winnow accs, avg winnow accs, ada accs, a
dv = DictVectorizer(sparse=False, dtype=int)
for size in sample sizes:
    x_train_slice = X_train[0:size]
    y train slice = y train[0:size]
    for i in range(0, len(models)):
        if (i==2):
            classifier=models[i](winnow hp selection, features)
        elif (i==3):
            classifier=models[i](winnow avgd hp selection, features)
        elif (i==4):
            classifier=models[i](ada_hp_selection, features)
        elif (i==5):
            classifier=models[i](ada avgd hp selection, features)
        else:
            classifier=models[i](features)
        classifier.train(x_train_slice, y_train_slice)
        y pred = classifier.predict(X test)
        acc = accuracy_score(y_test, y_pred)
        accs[i].append(acc)
        print(model_names[i] + ' accuracy with sample size ' + str(size)
    svc_x_train_slice = dv.fit_transform(x_train_slice)
    svc_x_test = dv.fit_transform(X_test)
    svc = LinearSVC(loss='hinge')
    svc.fit(svc_x_train_slice, y_train_slice)
    svc pred = svc.predict(svc x test)
    svc_acc = accuracy_score(y_test, svc_pred)
    accs[6].append(svc_acc)
    print('SVC accuracy with sample size ' + str(size) + ' is ' + str(svd
```

```
print()

# Plot Accuracies

save_destination = 'NA'

plot_learning_curves(per_accs, winnow_accs, ada_accs, avg_per_accs, avg_w
```

```
In [20]: ► | %%time
```

Run the synthetic experiment on the sparse dataset. For reference, "synthetic/sparse" is the path to where the data is located.

Note: This experiment takes substantial time (around 15 minutes), so don't worry if it's taking a long time to finish.

"""

run synthetic experiment('sparse')

Winnow Accuracy for Hyperparameter: 1.1 = 0.8935 AveragedWinnow Accuracy for Hyperparameter: 1.1 = 0.939

Winnow Accuracy for Hyperparameter: 1.01 = 0.927 AveragedWinnow Accuracy for Hyperparameter: 1.01 = 0.8875

Winnow Accuracy for Hyperparameter: 1.005 = 0.9195 AveragedWinnow Accuracy for Hyperparameter: 1.005 = 0.8145

Winnow Accuracy for Hyperparameter: 1.0005 = 0.563 AveragedWinnow Accuracy for Hyperparameter: 1.0005 = 0.5275

Winnow Accuracy for Hyperparameter: 1.0001 = 0.5205 AveragedWinnow Accuracy for Hyperparameter: 1.0001 = 0.5055

AdaGrad Accuracy for Hyperparameter: 1.5 = 0.913 AveragedAdaGrad Accuracy for Hyperparameter: 1.5 = 0.852

AdaGrad Accuracy for Hyperparameter: 0.25 = 0.7695 AveragedAdaGrad Accuracy for Hyperparameter: 0.25 = 0.762

AdaGrad Accuracy for Hyperparameter: 0.03 = 0.6635 AveragedAdaGrad Accuracy for Hyperparameter: 0.03 = 0.643

AdaGrad Accuracy for Hyperparameter: 0.005 = 0.5
AveragedAdaGrad Accuracy for Hyperparameter: 0.005 = 0.5055

AdaGrad Accuracy for Hyperparameter: 0.001 = 0.4985 AveragedAdaGrad Accuracy for Hyperparameter: 0.001 = 0.559

Ideal Winnow Hyperparameter: 1.01, Accuracy: 0.927
Ideal AdaGrad Hyperparameter: 1.5, Accuracy: 0.913
Ideal AveragedWinnow Hyperparameter: 1.1, Accuracy: 0.939
Ideal AveragedAdaGrad Hyperparameter: 1.5, Accuracy: 0.852

Perceptron accuracy with sample size 500 is 0.544
AveragedPerceptron accuracy with sample size 500 is 0.539
Winnow accuracy with sample size 500 is 0.5575
AveragedWinnow accuracy with sample size 500 is 0.6745
AdaGrad accuracy with sample size 500 is 0.5455
AveragedAdaGrad accuracy with sample size 500 is 0.5305
SVC accuracy with sample size 500 is 0.6115

/home/ian/anaconda3/lib/python3.8/site-packages/sklearn/svm/_base.py:976: C onvergenceWarning: Liblinear failed to converge, increase the number of ite rations.

warnings.warn("Liblinear failed to converge, increase "

Perceptron accuracy with sample size 1000 is 0.605
AveragedPerceptron accuracy with sample size 1000 is 0.6055
Winnow accuracy with sample size 1000 is 0.626
AveragedWinnow accuracy with sample size 1000 is 0.8595
AdaGrad accuracy with sample size 1000 is 0.602
AveragedAdaGrad accuracy with sample size 1000 is 0.5845
SVC accuracy with sample size 1000 is 0.931

/home/ian/anaconda3/lib/python3.8/site-packages/sklearn/svm/_base.py:976: C onvergenceWarning: Liblinear failed to converge, increase the number of ite rations.

warnings.warn("Liblinear failed to converge, increase "

Perceptron accuracy with sample size 1500 is 0.664
AveragedPerceptron accuracy with sample size 1500 is 0.6495
Winnow accuracy with sample size 1500 is 0.758
AveragedWinnow accuracy with sample size 1500 is 0.9175
AdaGrad accuracy with sample size 1500 is 0.674
AveragedAdaGrad accuracy with sample size 1500 is 0.6215
SVC accuracy with sample size 1500 is 0.936

Perceptron accuracy with sample size 2000 is 0.6295
AveragedPerceptron accuracy with sample size 2000 is 0.6845
Winnow accuracy with sample size 2000 is 0.8465
AveragedWinnow accuracy with sample size 2000 is 0.932
AdaGrad accuracy with sample size 2000 is 0.704
AveragedAdaGrad accuracy with sample size 2000 is 0.663
SVC accuracy with sample size 2000 is 0.936

Perceptron accuracy with sample size 2500 is 0.647
AveragedPerceptron accuracy with sample size 2500 is 0.711
Winnow accuracy with sample size 2500 is 0.8515
AveragedWinnow accuracy with sample size 2500 is 0.9355
AdaGrad accuracy with sample size 2500 is 0.644
AveragedAdaGrad accuracy with sample size 2500 is 0.689
SVC accuracy with sample size 2500 is 0.936

Perceptron accuracy with sample size 3000 is 0.6565 AveragedPerceptron accuracy with sample size 3000 is 0.74 Winnow accuracy with sample size 3000 is 0.8965 AveragedWinnow accuracy with sample size 3000 is 0.9355 AdaGrad accuracy with sample size 3000 is 0.761 AveragedAdaGrad accuracy with sample size 3000 is 0.712 SVC accuracy with sample size 3000 is 0.936

Perceptron accuracy with sample size 3500 is 0.7275 AveragedPerceptron accuracy with sample size 3500 is 0.7585 Winnow accuracy with sample size 3500 is 0.841 AveragedWinnow accuracy with sample size 3500 is 0.9355 AdaGrad accuracy with sample size 3500 is 0.7425 AveragedAdaGrad accuracy with sample size 3500 is 0.721 SVC accuracy with sample size 3500 is 0.936

Perceptron accuracy with sample size 4000 is 0.69 AveragedPerceptron accuracy with sample size 4000 is 0.7835 Winnow accuracy with sample size 4000 is 0.8555 AveragedWinnow accuracy with sample size 4000 is 0.936 AdaGrad accuracy with sample size 4000 is 0.756 AveragedAdaGrad accuracy with sample size 4000 is 0.741 SVC accuracy with sample size 4000 is 0.936

Perceptron accuracy with sample size 4500 is 0.675
AveragedPerceptron accuracy with sample size 4500 is 0.7995
Winnow accuracy with sample size 4500 is 0.9085
AveragedWinnow accuracy with sample size 4500 is 0.936
AdaGrad accuracy with sample size 4500 is 0.8095
AveragedAdaGrad accuracy with sample size 4500 is 0.754
SVC accuracy with sample size 4500 is 0.936

Perceptron accuracy with sample size 5000 is 0.766
AveragedPerceptron accuracy with sample size 5000 is 0.809
Winnow accuracy with sample size 5000 is 0.9115
AveragedWinnow accuracy with sample size 5000 is 0.936
AdaGrad accuracy with sample size 5000 is 0.796
AveragedAdaGrad accuracy with sample size 5000 is 0.7685
SVC accuracy with sample size 5000 is 0.936

Perceptron accuracy with sample size 10000 is 0.717

AveragedPerceptron accuracy with sample size 10000 is 0.9095

Winnow accuracy with sample size 10000 is 0.926

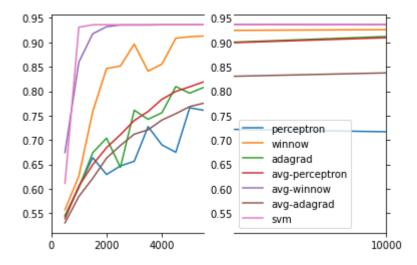
AveragedWinnow accuracy with sample size 10000 is 0.936

AdaGrad accuracy with sample size 10000 is 0.912

AveragedAdaGrad accuracy with sample size 10000 is 0.8375

SVC accuracy with sample size 10000 is 0.936

<Figure size 432x288 with 0 Axes>



CPU times: user 5min 47s, sys: 154 ms, total: 5min 47s

Wall time: 5min 47s

In [21]: ► %%time

.

Run the synthetic experiment on the dense dataset. For reference, "synthetic/dense" is the path to where the data is located. Note: this experiment should take much less time.

run_synthetic_experiment('dense')

Winnow Accuracy for Hyperparameter: 1.1 = 0.8995 AveragedWinnow Accuracy for Hyperparameter: 1.1 = 0.9445

Winnow Accuracy for Hyperparameter: 1.01 = 0.9215 AveragedWinnow Accuracy for Hyperparameter: 1.01 = 0.9315

Winnow Accuracy for Hyperparameter: 1.005 = 0.908 AveragedWinnow Accuracy for Hyperparameter: 1.005 = 0.913

Winnow Accuracy for Hyperparameter: 1.0005 = 0.8615 AveragedWinnow Accuracy for Hyperparameter: 1.0005 = 0.682

Winnow Accuracy for Hyperparameter: 1.0001 = 0.614 AveragedWinnow Accuracy for Hyperparameter: 1.0001 = 0.5465

AdaGrad Accuracy for Hyperparameter: 1.5 = 0.9445 AveragedAdaGrad Accuracy for Hyperparameter: 1.5 = 0.9445

AdaGrad Accuracy for Hyperparameter: 0.25 = 0.7825 AveragedAdaGrad Accuracy for Hyperparameter: 0.25 = 0.7055

AdaGrad Accuracy for Hyperparameter: 0.03 = 0.5795 AveragedAdaGrad Accuracy for Hyperparameter: 0.03 = 0.5885

AdaGrad Accuracy for Hyperparameter: 0.005 = 0.5085 AveragedAdaGrad Accuracy for Hyperparameter: 0.005 = 0.5085

AdaGrad Accuracy for Hyperparameter: 0.001 = 0.5085 AveragedAdaGrad Accuracy for Hyperparameter: 0.001 = 0.5085

Ideal Winnow Hyperparameter: 1.01, Accuracy: 0.9215
Ideal AdaGrad Hyperparameter: 1.5, Accuracy: 0.9445
Ideal AveragedWinnow Hyperparameter: 1.1, Accuracy: 0.9445
Ideal AveragedAdaGrad Hyperparameter: 1.5, Accuracy: 0.9445

Perceptron accuracy with sample size 500 is 0.5335 AveragedPerceptron accuracy with sample size 500 is 0.665 Winnow accuracy with sample size 500 is 0.799 AveragedWinnow accuracy with sample size 500 is 0.922 AdaGrad accuracy with sample size 500 is 0.6205 AveragedAdaGrad accuracy with sample size 500 is 0.5855 SVC accuracy with sample size 500 is 0.939

Perceptron accuracy with sample size 1000 is 0.676 AveragedPerceptron accuracy with sample size 1000 is 0.8705 Winnow accuracy with sample size 1000 is 0.907 /home/ian/anaconda3/lib/python3.8/site-packages/sklearn/svm/_base.py:976: ConvergenceWarning: Liblinear failed to converge, increase the number of iterations.

warnings.warn("Liblinear failed to converge, increase "

AveragedWinnow accuracy with sample size 1000 is 0.9405 AdaGrad accuracy with sample size 1000 is 0.591 AveragedAdaGrad accuracy with sample size 1000 is 0.6905 SVC accuracy with sample size 1000 is 0.9405

Perceptron accuracy with sample size 1500 is 0.8255 AveragedPerceptron accuracy with sample size 1500 is 0.933 Winnow accuracy with sample size 1500 is 0.8895 AveragedWinnow accuracy with sample size 1500 is 0.9405 AdaGrad accuracy with sample size 1500 is 0.8685 AveragedAdaGrad accuracy with sample size 1500 is 0.7685 SVC accuracy with sample size 1500 is 0.9405

Perceptron accuracy with sample size 2000 is 0.815
AveragedPerceptron accuracy with sample size 2000 is 0.9405
Winnow accuracy with sample size 2000 is 0.9205
AveragedWinnow accuracy with sample size 2000 is 0.9405
AdaGrad accuracy with sample size 2000 is 0.8925
AveragedAdaGrad accuracy with sample size 2000 is 0.846
SVC accuracy with sample size 2000 is 0.9405

Perceptron accuracy with sample size 2500 is 0.8715
AveragedPerceptron accuracy with sample size 2500 is 0.9405
Winnow accuracy with sample size 2500 is 0.9255
AveragedWinnow accuracy with sample size 2500 is 0.9405
AdaGrad accuracy with sample size 2500 is 0.932
AveragedAdaGrad accuracy with sample size 2500 is 0.8865
SVC accuracy with sample size 2500 is 0.9405

Perceptron accuracy with sample size 3000 is 0.7985
AveragedPerceptron accuracy with sample size 3000 is 0.9405
Winnow accuracy with sample size 3000 is 0.93
AveragedWinnow accuracy with sample size 3000 is 0.9405
AdaGrad accuracy with sample size 3000 is 0.9405
AveragedAdaGrad accuracy with sample size 3000 is 0.92
SVC accuracy with sample size 3000 is 0.9405

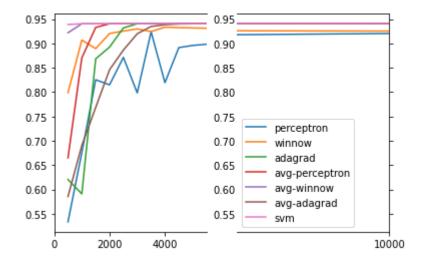
Perceptron accuracy with sample size 3500 is 0.924
AveragedPerceptron accuracy with sample size 3500 is 0.9405
Winnow accuracy with sample size 3500 is 0.9245
AveragedWinnow accuracy with sample size 3500 is 0.9405
AdaGrad accuracy with sample size 3500 is 0.9405
AveragedAdaGrad accuracy with sample size 3500 is 0.935
SVC accuracy with sample size 3500 is 0.9405

Perceptron accuracy with sample size 4000 is 0.8195 AveragedPerceptron accuracy with sample size 4000 is 0.9405 Winnow accuracy with sample size 4000 is 0.9335 AveragedWinnow accuracy with sample size 4000 is 0.9405 AdaGrad accuracy with sample size 4000 is 0.9405 AveragedAdaGrad accuracy with sample size 4000 is 0.9385 SVC accuracy with sample size 4000 is 0.9405 Perceptron accuracy with sample size 4500 is 0.8915
AveragedPerceptron accuracy with sample size 4500 is 0.9405
Winnow accuracy with sample size 4500 is 0.9325
AveragedWinnow accuracy with sample size 4500 is 0.9405
AdaGrad accuracy with sample size 4500 is 0.9405
AveragedAdaGrad accuracy with sample size 4500 is 0.94
SVC accuracy with sample size 4500 is 0.9405

Perceptron accuracy with sample size 5000 is 0.896
AveragedPerceptron accuracy with sample size 5000 is 0.9405
Winnow accuracy with sample size 5000 is 0.932
AveragedWinnow accuracy with sample size 5000 is 0.9405
AdaGrad accuracy with sample size 5000 is 0.9405
AveragedAdaGrad accuracy with sample size 5000 is 0.9405
SVC accuracy with sample size 5000 is 0.9405

Perceptron accuracy with sample size 10000 is 0.9205
AveragedPerceptron accuracy with sample size 10000 is 0.9405
Winnow accuracy with sample size 10000 is 0.9255
AveragedWinnow accuracy with sample size 10000 is 0.9405
AdaGrad accuracy with sample size 10000 is 0.9405
AveragedAdaGrad accuracy with sample size 10000 is 0.9405
SVC accuracy with sample size 10000 is 0.9405

<Figure size 432x288 with 0 Axes>



CPU times: user 1min 20s, sys: 80 ms, total: 1min 20s

Wall time: 1min 20s

Questions (5 points)

Answer the following questions:

- 1. Discuss the trends that you see when comparing the standard version of an algorithm to the averaged version (e.g., Winnow versus Averaged Winnow). Is there an observable trend?
 - The averaged version of each algorithm typically performs better than the regular (unaveraged) version, for both the sparse and dense datasets. This typically holds across all of the different sample sizes for each algorithm when the optimal hyperparameter is selected. For some hyperparameters, the averaged version did not do as well as the regular version during the portion of the experiment where we were testing on the dev data, but after selecting the optimal hyperparameters the average algorithm typically performed better. However, the exception to this rule is the Averaged AdaGrad algorithm, which would often perform worse than the regular AdaGrad algorithm.
- 2. We provided you with 10,000 training examples. Were all 10,000 necessary to achieve the best performance for each classifier? If not, how many were necessary? (Rough estimates, no exact numbers required)
 - All 10,000 results were sometimes not necessary to receive the best performance for each classifier. For the sparse dataset, the svm needed about 1500 examples, the AveragedWinnow about 2500, Winnow was pretty close at 5000, while the others actually did need all 10000 to reach optimal performance. For the dense dataset, the svm and AveragedWinnow only needed about 500, AveragedPerceptron about 1500, AdaGrad about 2500, AveragedAdaGrad and Winnow around 4000, with Perceptron the only one needing all 10000.
- 3. Report your Final Test Accuracies

Model	Sparse	Dense
Perceptron	.717	.9205
Winnow	.926	.9255
AdaGrad	.912	.9405
Averaged Perceptron	.9095	.9405
Averaged Winnow	.936	.9405
Averaged AdaGrad	.8375	.9405
SVM	.936	.9405

3.1.5 Extra Credit (10 points)

Included in the resources for this homework assignment is the code that we used to generate the synthetic data. We used a small amount of noise to create the dataset which you ran the experiments on. For extra credit, vary the amount of noise in either/both of the label and features. Then, plot the models' performances as a function of the amount of noise. Discuss your observations.

TODO: Extra Credit observations

3.2 NER Experiment: Welcome to the Real World (35 points)

The experiment with the NER data will analyze how changing the domain of the training and testing data can impact the performance of a model.

Instead of accuracy, you will use your F_1 score implementation in Section 0 to evaluate how well a model does. Recall measures how many of the actual entities the model successfully tagged as an entity.

$$\begin{split} \text{Precision} &= \frac{\#(\text{Actually Entity \& Model Predicted Entity})}{\#(\text{Model Predicted Entity})} \\ \text{Recall} &= \frac{\#(\text{Actually Entity \& Model Predicted Entity})}{\#(\text{Actually Entity})} \\ F_1 &= 2 \cdot \frac{\text{Precision} \times \text{Recall}}{\text{Precision} + \text{Recall}} \end{split}$$

For this experiment, you will only use the averaged basic Perceptron and SVM. Hence, no parameter tuning is necessary. Train both models on the CoNLL training data then compute the F_1 on the development and testing data of both CoNLL and Enron. Note that the model which is used to predict labels for Enron is trained on CoNLL data, not Enron data. Report the F_1 scores in a table.

3.2.1 Extracting NER Features (25 points)

Reread Section 2.2.2 to understand how to extract the features required to train the models and translate it to the code below.

```
In [22]:

    def extract ner features train(train):

                 Extracts feature dictionaries and labels from the data in "train"
                 Additionally creates a list of all of the features which were created.
                 We have implemented the w-1 and w+1 features for you to show you how
                 to create them.
                 TODO: You should add your additional featurization code here.
                 (which might require adding and/or changing existing code)
                 y = []
                 X = []
                 features = set()
                 for sentence in train:
                     padded = [('SSS', None)] * 3 + sentence[:] + [('EEE', None)] * 3
                     for i in range(3, len(padded) - 3):
                          y.append(1 if padded[i][1] == 'I' else -1)
                          feat1 = 'w-3=' + str(padded[i - 3][0])
                          feat2 = 'w-2=' + str(padded[i - 2][0])
                          feat3 = 'w-1=' + str(padded[i - 1][0])
                          feat4 = 'w+1=' + str(padded[i + 1][0])
                          feat5 = 'w+2=' + str(padded[i + 2][0])
                          feat6 = 'w+3=' + str(padded[i + 3][0])
                          feat7 = 'w-1=' + str(padded[i - 1][0]) + '&w-2=' + str(padded[i - 1][0])
                          feat8 = 'w+1=' + str(padded[i + 1][0]) + '&w+2=' + str(padded[i + 1][0])
                          feat9 = 'w-1=' + str(padded[i - 1][0]) + '&w+1=' + str(padded[i + 1][0])
                          feats = [feat1, feat2, feat3, feat4, feat5, feat6, feat7, feat8,
                          features.update(feats)
                          feats = {feature: 1 for feature in feats}
                          X.append(feats)
                 return features, X, y
```

Now, repeat the process of extracting features from the test data. What is the difference between the code above and below?

```
In [23]:
          ▶ def extract features dev or test(data, features):
                 Extracts feature dictionaries and labels from "data". The only
                 features which should be computed are those in "features". You
                 should add your additional featurization code here.
                 TODO: You should add your additional featurization code here.
                 y = []
                 X = []
                 for sentence in data:
                     padded = [('SSS', None)] * 3 + sentence[:] + [('EEE', None)] * 3
                     for i in range(3, len(padded) - 3):
                          y.append(1 if padded[i][1] == 'I' else -1)
                          feat1 = 'w-3=' + str(padded[i - 3][0])
                          feat2 = 'w-2=' + str(padded[i - 2][0])
                          feat3 = 'w-1=' + str(padded[i - 1][0])
                          feat4 = 'w+1=' + str(padded[i + 1][0])
                          feat5 = 'w+2=' + str(padded[i + 2][0])
                          feat6 = 'w+3=' + str(padded[i + 3][0])
                          feat7 = 'w-1=' + str(padded[i - 1][0]) + '&w-2=' + str(padded[i - 1][0])
                          feat8 = 'w+1=' + str(padded[i + 1][0]) + '&w+2=' + str(padded[i + 1][0])
                          feat9 = 'w-1=' + str(padded[i - 1][0]) + '&w+1=' + str(padded[i + 1][0])
                          feats = [feat1, feat2, feat3, feat4, feat5, feat6, feat7, feat8,
                          feats = {feature: 1 for feature in feats if feature in features}
                          X.append(feats)
                 return X, y
```

3.2.2 Running the NER Experiment

As stated previously, train both models on the CoNLL training data then compute the F_1 on the development and testing data of both CoNLL and Enron. Note that the model which is used to predict labels for Enron is trained on CoNLL data, not Enron data.

```
In [24]:

    def run ner experiment(data path):

                 Runs the NER experiment using the path to the ner data
                 (e.g. "ner" from the released resources). We have implemented
                 the standard Perceptron below. You should do the same for
                 the averaged version and the SVM.
                 The SVM requires transforming the features into a different
                 format. See the end of this function for how to do that.
                 train = load ner data(dataset='conll', dataset type='train')
                 conll_test = load_ner_data(dataset='conll', dataset_type='test')
                 enron_test = load_ner_data(dataset='enron', dataset_type='test')
                 features, X train, y train = extract ner features train(train)
                 X_conll_test, y_conll_test = extract_features_dev_or_test(conll_test, fea
                 X enron test, y enron test = extract features dev or test(enron test, features)
                 # TODO: We show you how to do this for Perceptron.
                 # You should do this for the Averaged Perceptron and SVM
                 classifier = Perceptron(features)
                 classifier.train(X_train, y_train)
                 y pred = classifier.predict(X conll test)
                 conll_f1 = calculate_f1(y_conll_test, y_pred)
                 y pred = classifier.predict(X enron test)
                 enron_f1 = calculate_f1(y_enron_test, y_pred)
                 print('Perceptron')
                 print('CoNLL Test Accuracy: ', conll f1)
                 print('Enron Test Accuracy: ', enron_f1)
                 print()
                 classifier = AveragedPerceptron(features)
                 classifier.train(X_train, y_train)
                 y pred = classifier.predict(X conll test)
                 conll_f1 = calculate_f1(y_conll_test, y_pred)
                 y pred = classifier.predict(X enron test)
                 enron_f1 = calculate_f1(y_enron_test, y_pred)
                 print('Averaged Perceptron')
                 print('CoNLL Test Accuracy: ', conll_f1)
                 print('Enron Test Accuracy: ', enron f1)
                 print()
                 # This is how you convert from the way we represent features in the
                 # Perceptron code to how you need to represent features for the SVM.
                 # You can then train with (X train dict, y train) and test with
                 # (X_conll_test_dict, y_conll_test) and (X_enron_test_dict, y_enron_test)
                 vectorizer = DictVectorizer()
                 X train dict = vectorizer.fit transform(X train)
                 X conll test dict = vectorizer.transform(X conll test)
                 X_enron_test_dict = vectorizer.transform(X_enron_test)
                 svc = LinearSVC(loss='hinge')
```

```
svc.fit(X_train_dict, y_train)
y_pred = svc.predict(X_conll_test_dict)
conll_f1 = calculate_f1(y_conll_test, y_pred)

y_pred = svc.predict(X_enron_test_dict)
enron_f1 = calculate_f1(y_enron_test, y_pred)
print('SVC')
print('Conll Test Accuracy: ', conll_f1)
print('Enron Test Accuracy: ', enron_f1)
print()
```

In [25]: # Run the NER experiment. "ner" is the path to where the data is located.
run_ner_experiment('ner')

Perceptron

CoNLL Test Accuracy: 0.7523076923076922 Enron Test Accuracy: 0.14016172506738545

Averaged Perceptron

CoNLL Test Accuracy: 0.8052126200274349 Enron Test Accuracy: 0.23608445297504793

SVC

CoNLL Test Accuracy: 0.827823691460055 Enron Test Accuracy: 0.24119530416221988

/home/ian/anaconda3/lib/python3.8/site-packages/sklearn/svm/_base.py:976: C onvergenceWarning: Liblinear failed to converge, increase the number of ite rations.

warnings.warn("Liblinear failed to converge, increase "

F1 Scores Table

Model	CoNLL Test F1	Enron Test F1
Averaged Perceptron	.80521	.23608
SVM	.82782	.24119

Questions (5 points)

Comment on the results:

1. Are the F₁ scores on CoNLL and Enron similar?

The F1 scores for the CoNLL and Enron datasets are pretty different, with one being about .8 and the other about .23.

2. If they are dissimilar, explain why you think the F₁ score increased/decreased on the Enron data.

I believe the F1 score decreased on the Enron data since the models were trained entirely on the CoNLL data. If the Enron data and CoNLL data are not similar, then there could be a lot of problems in trying to fit the model to the Enron data. One such problem could be features that were encountered in Enron but were not encountered in CoNLL and thus not factored into the model at all. Another problem is that the CoNLL data was pulled from newspaper articles, while the Enron data was pulled from emails. While both are text sources, the writing done in an email is typically different than writing done in a newspaper article, which could have also factored into the low Enron F1 score. Ultimately, I think the model would have performed better on a data set also pulled from a newspaper-type source, while the model prediction on Enron would have performed better had there been a training set that was from an email-like source.

Submission Instructions

We will be using Gradescope to turn in both the Python code. You should have been automatically added to Gradescope. If you do not have access, please ask the TA staff on Piazza.

There are three parts to the submission on Gradescope:

- ipynb file: Submit this notebook (.ipynb). If you are using Google Colab, you will have to download it as an ipynb file.
- PDF: This notebook saved as a PDF. We will use this to see the overall structure of your homework and code, and to check manual questions like the tables, plots, and your discussions. How should I convert this notebook to PDF, you may ask? There are a few ways, but for simplicity print the Jupyter notebook and Save as PDF.
- Code: A hw2.py file. We will use this to unit test and automate grading some parts of the homework. We only need a few functions/classes from the whole notebook. In particular, to unit test we only need:
 - calculate f1
 - highest and lowest f1 score
 - Perceptron, Winnow, and AdaGrad classes.

There are two ways to submit these pieces of code. You can either manually copy and paste these to a Python file, or you can use Jupyter's _Download as .py_, and delete all unnecessary code.