Machine Learning Assignment 2 Report

- Implementation
- + phi (X, lift): Lift the input instances to feature space, where lift denotes type of lifting function.

$$\Phi_1(X) = \begin{bmatrix} x_1 \\ x_2 \\ \vdots \\ x_N \end{bmatrix}$$

$$\Phi_{2}(X) = \begin{bmatrix} x_{1}^{2} & x_{1} \\ x_{2}^{2} & x_{2} \\ \vdots & \vdots \\ x_{N}^{2} & x_{N} \end{bmatrix}$$

$$\Phi_{1}(X) = \begin{bmatrix} x_{1} \\ x_{2} \\ \vdots \\ x_{N} \end{bmatrix} \qquad \Phi_{2}(X) = \begin{bmatrix} x_{1}^{2} & x_{1} \\ x_{2}^{2} & x_{2} \\ \vdots & \vdots \\ x_{N}^{2} & x_{N} \end{bmatrix} \qquad \Phi_{3}(X) = \begin{bmatrix} x_{1}^{3} & x_{1}^{2} & x_{1} \\ x_{2}^{3} & x_{2}^{2} & x_{2} \\ \vdots & \vdots & \vdots \\ x_{N}^{3} & x_{N}^{2} & x_{N} \end{bmatrix}$$

```
function featureX = phi(X, lift)
featureX = X;
while lift>1
featureX = [X.^lift featureX];
    lift = lift-1;
```

- +train(X, r, lift):
 - Initial W, B randomly in normal distribution.
 - Set error tolerance ϵ =10⁻¹⁰
 - Iterations from 0 to 10000

For every iteration,

- Set
$$\eta = \frac{\left|\left|(r-y)\Phi(X)\right|\right|}{3}$$

- Implement

$$W^{(i+1)} = W^{(i)} + \eta \sum_{t=1}^{N} (r^{(t)} - y^{(t)}) \Phi(x^{(t)})$$

$$B^{(i+1)} = B^{(i)} + \eta \sum_{t=1}^{N} (r^{(t)} - y^{(t)})$$

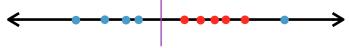
- function perceptronClassifierObj = train(X, r, lift) featureX = PerceptronClassifier.phi(X, lift); W = randn(lift, 1); B = randn;err = 1e-10: for i=0:10000 preW = W; preB = B;predictY = featureX*W + B; y = sign(predictY); eta = norm((r-y)'*featureX)/3; W = preW + eta*((r-y)'*featureX)'; B = preB + eta*sum(r-y); if norm(W-preW)<err break: perceptronClassifierObj = PerceptronClassifier(W, B);
- Halt if the convergence $||W^{(i+1)} W^{(i)}|| < \epsilon$ occurs, or the maximum iteration is reached.
- + Generalization (X, r, lift):

Generalization error is defined as the average of number of incorrect prediction.

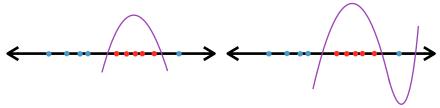
```
function generalErr = GeneralizationError(obj, X, r, N, lift)
liftX = obj.phi(X, lift);
predictY = liftX*obj.w + obj.b;
y = sign(predictY);
 generalErr = sum(y\sim=r)/N;
```

+ ConsistencyBound(N, lift):

For $\phi_1(X)$, the optimal way to shatter instances by a line is depicted as below. Since 10% of the instances will be mispredicted, the R[h*] = N/10.



For $\phi_2(X)$ and $\phi_3(X)$, it is always able to shatter instances, so the $R[h^*] = 0$;



Results

As N increases:

- (1) generalization error decreases because we have more history data to train and obtain a more accurate model.
- (2) consistency bound decreases. Observed from the formula of the bound, we can tighten the bound with larger N.

As dimension of feature space increases:

- (1) generalization error decreases. Because higher dimension model is more likely to shatter instances, thus prediction becomes more accurate.
- (2) consistency bound increases because the VC dimension increases.

