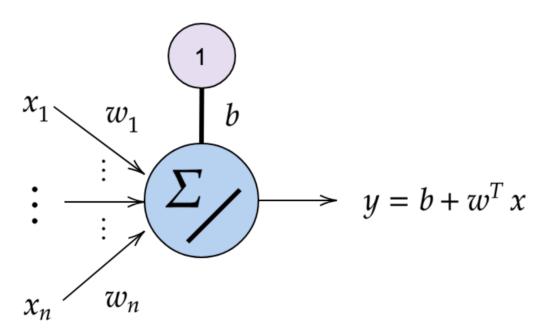


Second linear classifier:

• Fisher discriminant

First network:

• Perceptron



Linear Classifiers

Assume data $\mathbf{x}^{(i)} \in \mathbb{R}^d$ have been categorized into two classes.

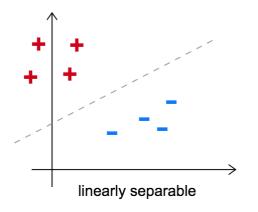
• Last time used probabilistic methods (Bayes' Rule) to separate classes.

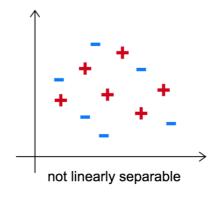
Method 1 - Fisher Discriminant

Classical statistical technique

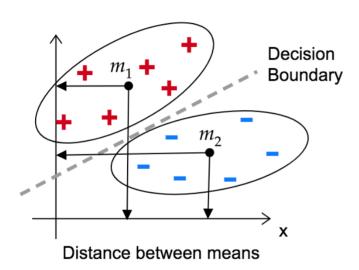
Method 2 - Perceptron

- Algorithm with many variants
- One of the few learning algorithms with a proof of convergence





Linear Classifiers



- Data sets are linearly separable but not when projected onto the x or y axis
- Is there a better line? What about line that maximizes projected distance between data sets?

Given points $\mathbf{x} \in \mathbb{R}^d$, compute class means \mathbf{m}_1 and \mathbf{m}_2

$$\mathbf{m}_1 = rac{1}{n_1} \sum_{C_1} \mathbf{x}_i \qquad \mathbf{m}_2 \qquad = rac{1}{n_2} \sum_{C_2} \mathbf{x}_i .$$

Find projection vector ${\bf w}$ that maximizes the projected distance between ${\bf m}_1$ and ${\bf m}_2$

$$\max_{\mathbf{w}} \|\mathbf{w}^T(\mathbf{m}_2 - \mathbf{m}_1)\|$$

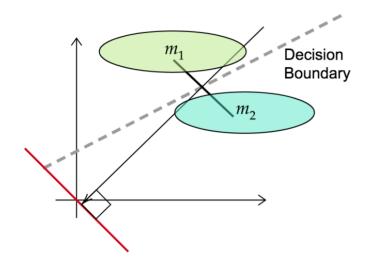
We know for unit vectors \mathbf{w} , that this is maximized when \mathbf{w} is in the same direction as $\mathbf{m}_2 - \mathbf{m}_1$

• Projecting onto direction defined by the line connecting the data set means maximizes distance between means of projections.

The projection of the mean vectors equals the mean of the projections

$$\mathbf{w}^T \mathbf{m}_i = \mathbf{w}^T \left(rac{1}{n_i} \sum_{C_i} \mathbf{x}_i
ight) = rac{1}{n_i} \sum_{C_i} \mathbf{w}^T \mathbf{x}_i$$

Linear Classifiers



- The data sets are linearly separable.
- Projecting onto a line parallel to line between means maximizes mean distance of the projected points.
- Again, data is separable but not when projected onto line connecting means.
- Is there a better line?
- Variance of the projected data must be part of the solution. This projection smears out the classes too much

Two preliminary results:

Empirical Covariance

• Will show that if X is a centered (zero mean) data matrix with samples arranged in rows, then the empirical covariance matrix is proportional to $\text{Cov} \sim X^T X$. If the data samples are arranged by column, then $\text{Cov}(X) \sim XX^T$

Re-expressing a projection using covariance

• The scalar $(\mathbf{w}^T\mathbf{x})^2$ can be computed from a vectors covariance matrix.

$$egin{aligned} (\mathbf{w}^T\mathbf{x})^2 &= (\mathbf{w}^T\mathbf{x})(\mathbf{w}^T\mathbf{x})^T \ &= (\mathbf{w}^T\mathbf{x})(\mathbf{x}^T\mathbf{w}) \ &= \mathbf{w}^T(\mathbf{x}\mathbf{x}^T)\mathbf{w} \end{aligned}$$

$$Cov[X, Y] = E[(X - E[X])(Y - E[Y])] = E[XY]$$
 when $E[X] = E[Y] = 0$

Given data vectors $\mathbf{x}^{(i)}$, $i=1,\cdots,m$ want to compute the empirical covariance of the \mathbf{x} components

Let $\tilde{\mathbf{x}}^{(i)}$ be the centered (zero mean) version of the data $\mathbf{x}^{(i)}$

$$\mathbf{ ilde{x}}^{(i)} = \mathbf{x}^{(i)} - rac{1}{m} \sum \mathbf{x}^{(i)}$$

Using that Cov(X,Y) = E[XY] when E[X] = E[Y] = 0, the (j,k) element of the (symmetric) covariance matrix is the mean of \tilde{x}_j times \tilde{x}_k

$$C_{j\,k} = rac{1}{m} \sum_{i=1}^m ilde{x}_j ilde{x}_k pprox E[X_j X_k]$$

Note: This empirical equation is biased, should divide by m-1 to get the unbiased estimate. The difference is seldom important.

The matrix outer product gives products of vector components

$$egin{bmatrix} x_1^{(i)} \ dots \ x_d^{(i)} \end{bmatrix} [x_1^{(i)}, \cdots, x_d^{(i)}] = egin{bmatrix} x_1^{(i)} x_1^{(i)} & \cdots & x_1^{(i)} x_d^{(i)} \ dots & dots \ x_d^{(i)} x_1^{(i)} & \cdots & x_d^{(i)} x_d^{(i)} \end{bmatrix}$$

Showed earlier that one way to interpret matrix multiplication AB is

$$AB = \sum_{i} col_{i}(A) \bigotimes row_{i}(B)$$

If the vectors $\mathbf{x}^{(i)}$ are arranged as the columns of X then the covariance matrix is:

$$egin{aligned} C &= rac{1}{m} \sum_i col_i(X) \, row_i(X^T) \ &= rac{1}{m} X X^T \end{aligned}$$

note 1: It is the transpose of this if data is arranged by rows.

note 2: The dimension of the covariance matrix should be $d \times d$ if the data vectors have d elements

Given m data points $\mathbf{x}^{(i)} \in \mathbb{R}^d$ write the data matrix X as

$$X = egin{bmatrix} \mathbf{x}_1^{(1)} & \cdots & \mathbf{x}_d^{(1)} \ dots & & dots \ \mathbf{x}_1^{(m)} & \cdots & \mathbf{x}_d^{(m)} \end{bmatrix}$$

In X, the samples are arranged in rows and features/coordinates in columns
The sample mean vector (column means) is

$$\mu = rac{1}{m} \sum_{i=1}^m \mathbf{x}^{(i)}$$

 $\mu \in \mathbb{R}^d$. The covariance matrix for the data has generic (j,k) entries (definition)

$$ext{Cov}(X)_{jk} = E[(\mathbf{x}_j - \mu_j)(\mathbf{x}_j - \mu_j)] pprox rac{1}{m-1} \sum_{i=1}^m (\mathbf{x}_j^{(i)} - \mu_j)(\mathbf{x}_k^{(i)} - \mu_k)$$

$$\mathrm{Cov}(X) \in d imes d$$

If we take just one row of the data, say $\mathbf{x}^{(p)}$ and think of $\mathbf{x}^{(p)} - \mu$ as a row vector, compute

$$S^{(p)} = \underbrace{(\mathbf{x}^{(p)} - \mu)^T}_{ ext{column}} \underbrace{(\mathbf{x}^{(p)} - \mu)}_{ ext{row}}$$

The generic (i, j) entry for $S^{(p)}$ is

$$S_{ij}^{(p)} = (\mathbf{x}_i^{(p)} - \mu_i)^T (\mathbf{x}_j^{(p)} - \mu_j)$$

This is the $\mathbf{x}^{(p)}$ contribution to the covariance matrix. Using this, the Cov(X) can be written as

$$egin{align} \operatorname{Cov}(X) &= rac{1}{m-1} \sum_{i}^{m} (\mathbf{x}^{(i)} - \mu)^T (\mathbf{x}^{(i)} - \mu) \ &= rac{1}{m-1} \sum_{i}^{m} S^{(i)} \end{split}$$

This is a sum of outer products.

Fisher condition

Find

$$\max_{\mathbf{w}} J(\mathbf{w})$$

where

$$J=rac{(m_2-m_1)^2}{s_1^2+s_2^2}$$

- m_i are class means projected onto ${f w}$, so m_2-m_1 measures between class distance
- ullet s_i^2 are within-class scatter, so $s_1^2+s_2^2$ measures total within-class variation
- *J* is maximized by finding a **w** that balances increasing the distance between the projected class means while not smearing the class scatter too much

Let ${\bf w}$ define the unknown 1^d projection subspace. Assume $\|{\bf w}\|=1$

Data sets
$$x_j^{(i)} \in \mathbb{R}^d$$
, $i=1,2\dots n_j,\, j\in 1,2$

$$egin{aligned} y^{(i)} &= \mathbf{w}^T \mathbf{x}^{(i)} \ \mathbf{m}_j &= rac{1}{n_j} \sum_{C_j} \mathbf{x}^{(i)} \ & ilde{m}_j &= rac{1}{n_j} \sum_{C_j} y^{(i)} \ &= rac{1}{n_j} \sum_{C_j} \mathbf{w}^T \mathbf{x}^{(i)} \ &= \mathbf{w}^T \mathbf{m}_j \end{aligned}$$

The distance between projected means is

$$| ilde{m}_1- ilde{m}_2|=|\mathbf{w}^T(\mathbf{m}_1-\mathbf{m}_2)|$$

Already showed that the distance between projected means is maximized by the line connecting the data set means.

Define the scatter for the projected data as

$$s_j^2 = \sum_{y \in C_j} (y - ilde{m})^2$$

so $s_1^2 + s_2^2$ is an estimate of the total variance of the projected data.

The **Fisher Linear Discriminant** is the line that maximizes:

$$J(\mathbf{w}) = rac{(ilde{m}_1 - ilde{m}_2)^2}{s_1^2 + s_2^2} \, .$$

It takes some algebra to express J as a function \mathbf{w}

Consider $s_1^2 + s_2^2$

$$s_1^2 + s_2^2 = \sum_{C_1} (y - ilde{m}_1)^2 + \sum_{C_2} (y - ilde{m}_2)^2$$

$$egin{aligned} s_i^2 &= \sum_{x \in C_i} (\mathbf{w}^T x - \mathbf{w}^T \mathbf{m}_i)^2 \ &= \sum_{x \in C_i} (\mathbf{w}^T (\mathbf{x} - \mathbf{m}_i))^2 \end{aligned}$$

Now use result derived earlier $(\mathbf{w}^T x)^2 = \mathbf{w}^T (XX^T) \mathbf{w}$ and rewrite the within class scatter as

$$egin{aligned} s_i &= \sum_{x \in C_i} \mathbf{w}^T (\mathbf{x} - \mathbf{m}_i) (\mathbf{x} - \mathbf{m}_i)^T \mathbf{w} \ &= \mathbf{w}^T \Bigl(\sum_{x \in C_i} (\mathbf{x} - \mathbf{m}_i) (\mathbf{x} - \mathbf{m}_i)^T \Bigr) \mathbf{w} \end{aligned}$$

Note that summation is proportional to the data covariance matrix for class i. It measures variation within the class.

Defining

$$egin{aligned} S_i &= \sum_{x \in C_i} (\mathbf{x} - \mathbf{m}_i) (\mathbf{x} - \mathbf{m}_i)^T \ S_w &= S_1 + S_2 \end{aligned}$$

Gives

$$egin{aligned} s_1^2 + s_2^2 &= \mathbf{w}^T S_1 \mathbf{w} + \mathbf{w}^T S_2 \mathbf{w} \ &= \mathbf{w}^T S_w \mathbf{w} \end{aligned}$$

The numerator of the Fisher cost function *J* is

$$egin{aligned} (ilde{m}_1 - ilde{m}_2)^2 &= (\mathbf{w}^T \mathbf{m}_1 - \mathbf{w}^T \mathbf{m}_2)^2 \ &= (\mathbf{w}^T (\mathbf{m}_1 - \mathbf{m}_2))^2 \ &= \mathbf{w}^T \Big((\mathbf{m}_1 - \mathbf{m}_2) (\mathbf{m}_1 - \mathbf{m}_2)^T \Big) \mathbf{w} \end{aligned}$$

Define $S_B=(\mathbf{m}_1-\mathbf{m}_2)(\mathbf{m}_1-\mathbf{m}_2)^T$. S_B measures variation between classes. Combining results:

$$J(\mathbf{w}) = rac{\mathbf{w}^T S_B \mathbf{w}}{\mathbf{w}^T S_w \mathbf{w}}$$

Want to find the w that maximizes J. Note that S_B has rank 1 and

$$S_B\mathbf{x} = (\mathbf{m}_2 - \mathbf{m}_1)(\mathbf{m}_2 - \mathbf{m}_1)^T\mathbf{x}$$

The objective function is a ratio of scalars. Use vector/matrix derivatives for before to compute the derivative of J wrt \mathbf{w}

$$\frac{\partial J}{\partial \mathbf{w}} = 0$$

gives:

$$(\mathbf{w}^T S_B \mathbf{w}) S_w \mathbf{w} = (\mathbf{w}^T S_w \mathbf{w}) S_B \mathbf{w}$$

 S_B is the outer product of $\mathbf{m}_2 - \mathbf{m}_1$ so $S_B \mathbf{w}$ is in the direction of $\mathbf{m}_2 - \mathbf{m}_1$.

We are only interested in the direction of \mathbf{w} so constant multipliers can be ignored

$$c_B \, S_w \mathbf{w} = c_w ((\mathbf{m}_2 - \mathbf{m}_1)^T \mathbf{w}) (\mathbf{m}_2 - \mathbf{m}_1)$$

With $c_m = (\mathbf{m}_2 - \mathbf{m}_1)^T \mathbf{w}$

$$\mathbf{w} = rac{c_w}{c_B} c_m S_w^{-1} (\mathbf{m}_2 - \mathbf{m}_1)$$

$$\mathbf{w} \sim S_w^{-1} (\mathbf{m}_2 - \mathbf{m}_1)$$

where

$$egin{align} \mathbf{m}_j &= rac{1}{n_j} \sum_{C_j} \mathbf{x}^{(i)} \ S_i &= \sum_{x \in C_i} (\mathbf{x} - \mathbf{m}_i) (\mathbf{x} - \mathbf{m}_i)^T \ S_w &= S_1 + S_2 \end{aligned}$$

So far have a direction which helps separate two classes when they are projected. Now need threshold which discriminates the projected data (scalars)

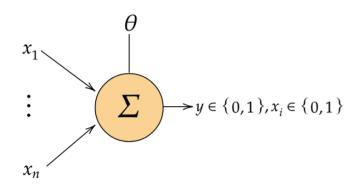
- Could assume that the projected data are Gaussian and use Bayes' rule from Week 1
- Bishop shows the following:

$$egin{aligned} \mathbf{m} &= rac{1}{n_1 + n_2} (n_1 \mathbf{m}_1 + n_2 \mathbf{m}_2) \ g(\mathbf{x}) &= \mathbf{w}^T (\mathbf{x} - \mathbf{m}) \end{aligned}$$

• The discriminant is:

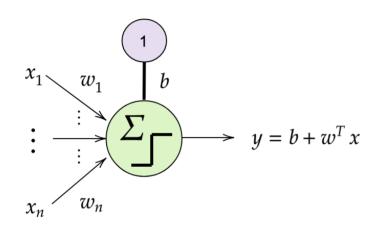
$$g(\mathbf{x}) = \left\{ egin{array}{ll} > 0 & \mathcal{C}_1 \ o. \, w. & \mathcal{C}_2 \end{array}
ight.$$

The perceptron evolved from the McCulloch-Pitts neuron (1943) where $y \in \{0,1\}$, $\mathbf{x}_i \in \{0,1\}$. There are no dynamic weights



$$egin{aligned} y &= f(g(\mathbf{x})) \ g(\mathbf{x}) &= \sum_i^n \mathbf{x}_i \ f &= egin{cases} 1 & ext{if } g(x) \geq 0 \ 0 & ext{o.w.} \end{cases}$$

- Threshold θ fixed
- Inhibitory inputs have veto power
- Excitory inputs have integer values (weights)



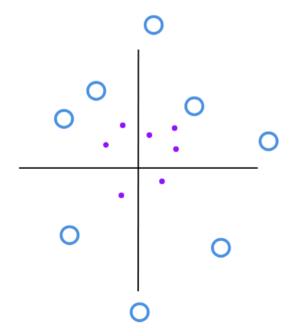
No cost function

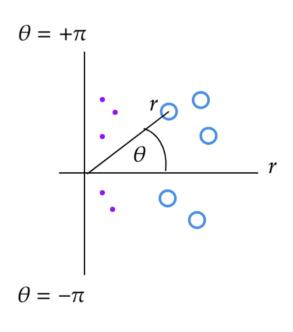
The perceptron modifies the M-P neuron to be more flexible

- Real valued inputs (positive or negative)
- Threshold becomes an adjustable parameter
- Modified output to $\{-1,1\}$
- Drops inhibitory inputs
- Has Learning rule (justly famous)
- Convergence proof (for linearly seperable data)

The original Perceptron algorithm is only useful for linearly separable data.

• Preprocessing, like transformation to polar coordinates, can change unseparable to to separable



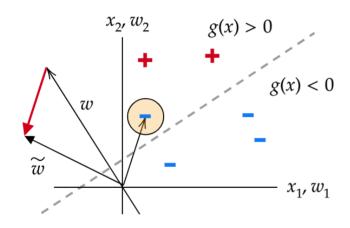


Perceptron Algorithm

```
Data \mathcal{D}=(x^{(i)},t^{(i)}), i=1,\cdots,m
x^{(i)} \in \mathbb{R}^d, t^{(i)} \in \{-1, 1\}
b, w \leftarrow 0
{Scale Data}
For i = 1, \dots, \max iterations do
   for each pair (x^{(i)}, t^{(i)}) \in \mathcal{D} do:
          z = b + w^T x^{(i)}
      if t^{(i)}z \leq 0) then
          w \leftarrow w + t^{(i)} x^{(i)}
         b \leftarrow b + t^{(i)}
       end
   end
end
```

Perceptron Algorithm

Heuristic View



Highlighted point is misclassified

say
$$b + \mathbf{w}^T \mathbf{x}^{(i)}) > 0$$
 but $t^{(i)} = -1$

Update rule:

$$\mathbf{ ilde{w}} = \mathbf{w} + t^{(i)}\mathbf{x} = \mathbf{w} - \mathbf{x}^{(i)}$$
 $\tilde{b} = b + t^{(i)} \cdot 1 = b - 1$

Update b,w and apply to $x^{(i)}$ again:

$$egin{aligned} b + \mathbf{w}^T \mathbf{x}^{(i)} &= z^{(i)} > 0 \ ilde{b} + ilde{\mathbf{w}}^T \mathbf{x}^{(i)} &= b - 1 + (\mathbf{w} - \mathbf{x}^{(i)})^T \mathbf{x}^{(i)} \ &= z^{(i)} - 1 - (\mathbf{x}^{(i)})^T \mathbf{x}_i \ &= z^{(i)} - 1 - \|\mathbf{x}^{(i)}\|^2 \end{aligned}$$

So $z^{(i)}$ is reduced

Helps this point, does it hurt other points?

Two ways to write equations

- with explicit bias
- · implicit bias and augmented data

Explicit bias

$$egin{aligned} y^{(i)} &= b + \mathbf{w}^T \mathbf{x}^{(i)} \ \mathbf{w} &= (w_1, \cdots, w_d) \in \mathbb{R}^d \ \mathbf{x}^{(i)} &= (x_1^{(i)}), \cdots, x_d^{(i)}) \in \mathbb{R}^d \end{aligned}$$

Implicit bias (homogeneous form)

$$egin{aligned} y^{(i)} &= \mathbf{ ilde{w}}^T \mathbf{ ilde{x}}^{(i)} \ \mathbf{ ilde{w}} &= (w_0, w_1, \cdots, w_d) \in \mathbb{R}^{d+1} \ \mathbf{ ilde{x}}^{(i)} &= (1, x_1^{(i)}, \cdots, x_d^{(i)}) \in \mathbb{R}^{d+1} \end{aligned}$$

identify w_0 with b

- typically, the augmented vectors $\tilde{\mathbf{w}}$ and $\tilde{\mathbf{x}}^{(i)}$ are written without tildes and are made clear by context
- Implicit bias almost always used for analysis
- Will use explicit bias in homeworks to match later neural network equations

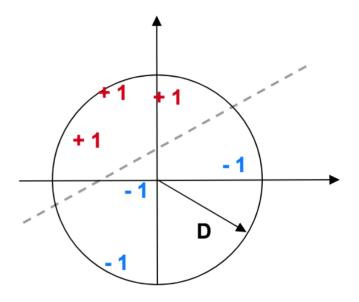
Will use homogeneous form where $x_1^{(i)}=1$

• Given data $(\mathbf{x}^{(i)}, t^{(i)})$ $i = 1, \cdots, m$

$$\mathbf{x}^{(i)} \in \mathbb{R}^{d+1}$$

$$t^{(i)} \in \{-1,1\}$$

- Assume $\|\mathbf{x}^{(i)}\| \leq D \quad \forall \quad i$
- Assume classes are linearly separable.



Theorem:

$$\exists \ \mathrm{vector} \ \mathbf{u}, \ |\mathbf{u}| = 1 \ \mathrm{such} \ \mathrm{that} \ t^{(i)} \mathbf{u}^T \mathbf{x}^{(i)} > 0 \ \forall \ i$$

This implies there is a margin $\gamma > 0 \ni t^{(i)}\mathbf{u}^T\mathbf{x}^{(i)} > \gamma$ and the total number of mistakes is bounded above by $(\frac{D}{\gamma})^2$ where $\|\mathbf{x}\| \leq D$

Let the adjustable parameters be $\mathbf{w} \in \mathbb{R}^{d+1}$ and let w=0 initially

- Theorem says that a solution (separator) exists
- Puts an upper bound on how much work is needed to find u
- Will show that \mathbf{w} will iteratively get closer to \mathbf{u} ($\mathbf{w} \cdot \mathbf{u}$ grows)
- The growth in $(\mathbf{w} \cdot \mathbf{u})$ is not dominated by the growth in \mathbf{w}

Let \mathbf{w}^k be parameter values when k-th mistake is made

•
$$\mathbf{w} = 0$$
 initially, so $\mathbf{w}^{(1)} = 0$

It follows that

$$egin{align} \operatorname{sign}(\mathbf{x^{(j)}} \cdot \mathbf{w}^{(k)})
eq t^{(j)} & \operatorname{some} j \ t^{(j)} & \operatorname{sign}(\mathbf{x^{(j)}} \cdot \mathbf{w}^{(k)}) < 0 \ \mathbf{w}^{k+1} & = \mathbf{w}^{(k)} + t^{(j)} \mathbf{x}^{(j)} \ \end{array}$$

$$\mathbf{u}^T \mathbf{w}^{k+1} = \mathbf{u}^T \mathbf{w}^{(k)} + t^{(j)} \mathbf{u}^T \mathbf{x}^{(j)} \\ > \mathbf{u}^T \mathbf{w}^{(k)} + \gamma$$

Telescope, using $\mathbf{w}^{(1)} = 0$

$$egin{aligned} \mathbf{u}^T\mathbf{w}^{(2)} &\geq \gamma \ \mathbf{u}^T\mathbf{w}^{(3)} &\geq \mathbf{u}^T\mathbf{w}^{(2)} + \gamma \ &\geq 2\gamma \end{aligned}$$

$$\mathbf{u}^T\mathbf{w}^{k+1} \geq k\gamma$$

shows projection of w on u grows

Now need to show that the growth is not all due to growth in $\mathbf{w}^{(k)}$

$$\begin{aligned} \|\mathbf{w}^{k+1}\|^2 &= \|\mathbf{w}^{(k)} + t^{(j)}x^{(j)}\|^2 \\ &= \|\mathbf{w}^k\|^2 + \|x^{(j)}\|^2 + \underbrace{2t^{(j)}(\mathbf{x^{(j)}})^T}_{must\ be\ <\ 0} \mathbf{w^{(k)}} \\ &\leq \|\mathbf{w}^k\|^2 + \|x^{(j)}\|^2 \\ &\leq \|\mathbf{w}^k\|^2 + D^2 \end{aligned}$$

By induction

$$egin{aligned} \|\mathbf{w}^{k+1}\|^2 & \leq kD^2 \ \sqrt{k}D & \geq \|\mathbf{w}^{k+1}\| \ & \geq (\mathbf{w}^{(k+1)})^T\mathbf{u} \ & \geq k\gamma \quad ext{by previous slide} \end{aligned}$$

$$k \leq (D/\gamma)^2$$

$$egin{aligned} \mathbf{u}^T\mathbf{w}^{k+1} &\geq k\gamma \ \|\mathbf{w}^{k+1}\|^2 \leq kD^2 \ &k \leq (D/\gamma)^2 \ &\mathbf{u}^T\mathbf{w}^{k+1} = \|\mathbf{w}^{k+1}\|\cos heta \ \cos heta &= rac{\mathbf{u}^T\mathbf{w}^{k+1}}{\|\mathbf{w}^{k+1}\|} \geq rac{k\gamma}{\sqrt{k}D} = \sqrt{k}rac{\gamma}{D} \end{aligned}$$

It follows that θ , the angle between **u** and **w** decreases as each mistake is corrected

Learning Rules

Error correcting rules

(no cost function):

- Perceptron (vanilla)
- Pocket
- Voted
- Averaged
- Perceptron with margin

Cost reduction rules

(gradient descent, MSE)

- Widrow-Hoff (LMS)
- Delta rule
- Ho Kashyap

Performance

- The error correcting algorithm will perfectly separate separable data.
- The delta rule minimizes squared error but may not be perfect on separable data.

Ensembles

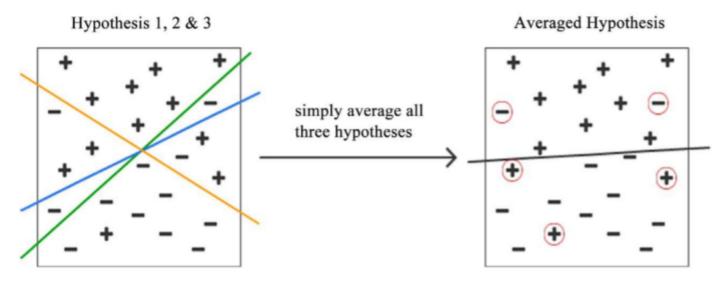


Figure 4. Ensemble model by Averaging.

Averaged Perceptron Algorithm

All hyperplane solutions $b^{(k)}$, $\mathbf{w}^{(k)}$ and all survival times $c^{(k)}$ are stored. The prediction rule uses the survival time weighting of the saved solutions:

$$egin{align} \hat{y}(x) &= ext{sign}(\sum_{k=1}^K c^{(k)}(b^{(k)} + \mathbf{w}^{(k)} \cdot x)) \ &= ext{sign}\left(\sum_{k=1}^K \left(c^{(k)}\mathbf{w}^{(k)}
ight) \cdot x + \sum_{k=1}^K c^{(k)}b^{(k)}
ight) \end{aligned}$$

Voted Perceptron Algorithm

All hyperplane solutions are remembered and weighted by how long they survive.

Run the baseline perceptron algorithm and while training store

- bias and weights at each update
- the number of samples the bias/weights survived
- The last update will survive for the entire data set

The voted perceptron prediction rule is

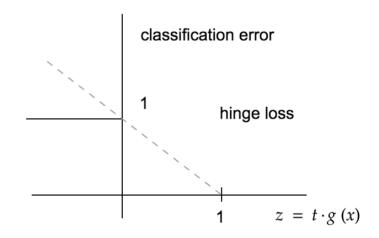
$$\hat{y}(x) = \operatorname{sign} \sum_{1}^{K} c^{(k)} \operatorname{sign}(b^{(k)} + \mathbf{w}^{(k)} \cdot \mathbf{x})$$

Perceptron Cost Function

Cost Function

$$t^{(i)} \in \{-1, +1\}$$
 $y^{(i)} = \left\{egin{array}{l} +1 ext{ if } g(x^{(i)}) \geq 0 \ -1 & o. \, w. \end{array}
ight.$

$t^{(i)}$	$y^{(i)}$	$t^{(i)} \cdot g(x)$	
- 1	- 1	> 0	correct
- 1	+ 1	≥ 0	incorrect
+ 1	- 1	< 0	incorrect
+ 1	- 1	> 0	correct



$$Cost = \max(0, 1-z) = \max(0, 1-t \cdot g(x))$$
 $rac{\partial C}{\partial W_j} = -t \cdot x_j$

Update w in opposite direction of gradient if misclassified