CS229-Cheatsheet

Supervised Learning

- Gradient Descent: to minimize $J(\theta)$, we perform $\theta_j := \theta_j \alpha \frac{\partial}{\partial \theta} J(\theta)$
- $\nabla_A AB = B^T$, $\nabla_{A^T} f(A) = (\nabla_A f(A))^T$, $\nabla_A tr ABA^T C = CAB + C^T AB^T$, $\nabla_A |A| = |A|(A^{-1})^T$
- Normal Equations and Least Squares

$$J(\theta) = \frac{1}{2} \sum_{i=1}^{m} (h_{\theta}(x^{i}) - y^{i})^{2} \to \nabla_{\theta} J(\theta) =$$

$$\nabla_{\theta} \frac{1}{2} (X\theta - y)^{T} (X\theta - y) = X^{T} X\theta - X^{T} y = 0 \to$$

$$X^{T} X\theta = X^{T} y \to \theta = (X^{T} X)^{-1} X^{T} y.$$

- Locally Weighted Regression Fit θ to minimize $\sum_{i=0}^m (y^i \theta^T x^i)^2 \text{ where } w^i = e^{-\frac{(x^i x)^2}{2\tau^2}}$
- Logistic Regression: $h_{\theta}(x) = g(\theta^T x) = \frac{1}{1 + e^{-\theta^T x}},$ $g(z) = \frac{1}{1 + e^{-z}}, g'(z) = \frac{d}{dz} \frac{1}{1 + e^{-z}} = g(z)(1 g(z)),$ $p(y|x;\theta) = (h_{\theta}(x))^y (1 h_{\theta}(x))^{1-y}.$ $l(\theta) = logL(\theta) = \sum_{i=1}^m y^i logh(x^i) + (1 y^i) log(1 h(x^i)),$ $\frac{\partial}{\partial \theta_i} l(\theta) = (y h_{\theta}(x))x_j$
- Perceptron Learning Algorithm $\theta_i := \theta_i + \alpha(y^i h_{\theta}(x^i))x_i^i$
- Newton's Method: $\theta := \theta \frac{f(\theta)}{f'(\theta)}$, we want the first derivative to be zero, then $\theta := \theta \frac{l'(\theta)}{l''(\theta)}$, if θ is a vector then $\theta := \theta H^{-1}\nabla_{\theta}l(\theta)$ where $H_{ij} = \frac{\partial^{2}l(\theta)}{\partial\theta_{i}\partial\theta_{j}}$
- Exponential Family $p(y; \eta) = b(y)exp(\eta^T T(y) a(\eta))$
- General Linear Model Assumptions: 1. $y|x; \eta \sim Exponential Family(\eta)$. 2. Given x our goal is to predict the expected value of T(y) which is usually just y, so we would like our hypothesis to satisfy h(x) = E(y|x). 3. The natural parameter η and inputs x are related linearly. $\eta = \theta^T x$.
- Canonical response function: the distribution's mean as a function of the natural parameter $g(\eta) = E(T(y); \eta)$.

Generative Learning Algorithm Gaussian Discriminant Analysis

• $y \sim Bernoulli(\phi), x|y = 0 \sim N(\mu_0, \Sigma), x|y = 1 \sim N(\mu_1, \Sigma).$

- $p(y) = \phi^y (1 \phi)^{1-y}$
- $p(x|y=0) = \frac{1}{(2\pi)^{n/2}|\Sigma|^{1/2}} exp(-\frac{1}{2}(x-\mu_0)^T \Sigma^{-1}(x-\mu_0))$
- $p(x|y=1) = \frac{1}{(2\pi)^{n/2}|\Sigma|^{1/2}} exp(-\frac{1}{2}(x-\mu_1)^T \Sigma^{-1}(x-\mu_1))$
- $\bullet \ l(\phi,\mu_0,\mu_1,\Sigma) = \log \prod_{i=1}^m p(x^i,y^i;\phi,\mu_0,\mu_1,\Sigma) =$ $\log \prod_{i=1}^m p(x^i|y^i;\phi,\mu_0,\mu_1,\Sigma) p(y^i,\phi).$
- By maximizing l with respect to the parameters, we find the maximum likelihood of the parameters to be:

$$\begin{split} \phi &= \frac{1}{m} \mathbf{1}\{y^i = 1\} \\ \mu_0 &= \frac{\sum_{i=1}^m \{y^i = 0\} x^i}{\sum_{i=1}^m \{y^i = 0\}} \\ \mu_1 &= \frac{\sum_{i=1}^m \{y^i = 1\} x^i}{\sum_{i=1}^m \{y^i = 1\}} \\ \Sigma &= \frac{1}{m} \sum_{i=1}^m (x^i - \mu_{y^i})^T (x^i - \mu_{y^i}) \end{split}$$

Naive Bayes

• Naive Assumption:

$$p(x_1, x_2, \dots | y) = p(x_1 | y) p(x_2 | y) \dots = \prod_{i=1}^{n} p(x_i | y)$$

- Laplace Smoothing $\phi_j = \frac{\sum_{i=1}^m 1\{z^i = j\}}{m} \to \frac{\sum_{i=1}^m 1\{z^i = j+1\}}{m+k}, \text{ where } k$ represent the number of possible outcomes for z.
- $\begin{aligned} \bullet & \text{ Event Driven Text Classification:} \\ \phi_{k|y=1} &= \frac{\sum_{i=1}^{m} \sum_{j=1}^{n_i} 1\{x_j^i = k \wedge y^i = 1\}}{\sum_{i=1}^{m} 1\{y^i = 1\}n_i} \\ \phi_{k|y=0} &= \frac{\sum_{i=1}^{m} \sum_{j=1}^{n_i} 1\{x_j^i = k \wedge y^i = 0\}}{\sum_{i=1}^{m} 1\{y^i = 0\}n_i} \\ \phi_y &= \frac{\sum_{i=1}^{m} 1\{y^i = 1\}}{m} \end{aligned}$

Support Vector Machines

- Classifier: $h_{w,b}(x) = g(w^T x + b)$ where g(z) = 1 if z > 0 and g(z) = -1 otherwise.
- Functional Margins: $\hat{\gamma}^i = y^i(w^Tx + b)$, the smallest functional margin in the training set is called: $\hat{\gamma} = min_{i=1,2,...,m}\hat{\gamma}^i$
- Geometric Margins: $\gamma^i = y^i((\frac{w}{||w||})^T x^i + \frac{b}{||w||})$, the smallest geometric margin in a training set is $: \gamma = min_{i=1,...,m} \gamma^i$
- Optimal Margin Classifier: $min_{\gamma,w,b} \frac{1}{2} ||w||^2$ s.t $y^i(w^T x^i + b) > 1, i = 1, 2, ..., m$

• Lagrangian

$$L(w, b, \alpha) = \frac{1}{2}||w||^2 - \sum_{i=1}^{m} \alpha_i(y^i(w^Tx^i + b) - 1)$$

• The dual problem

$$max_{\alpha}W(\alpha) = \sum_{i=1}^{m} \alpha_i - \frac{1}{2} \sum_{i,j=1}^{m} y^i y^j \alpha_i \alpha_j (x^i)^T x^j$$

s.t
$$\sum_{i=1}^{m} \alpha_i y^i = 0$$

• Observations: 1. Most of the α_i s will be zero 2.

$$w^{T}x + b = (\sum_{i=1}^{m} \alpha_{i} y^{i} x^{i})^{T} x + b$$

• KKT Conditions:

$$\frac{\partial}{\partial w_i} L(w^*, \alpha^*, \beta^*) = 0, i = 1, \dots n$$

$$\frac{\partial}{\partial \beta_i} L(w^*, \alpha^*, \beta^*) = 0, i = 1, \dots l$$

$$\alpha^* g_i(w^*) = 0, i = 1, \dots, k$$

$$g_i(w^*) \le 0, i = 1, \dots, k$$

$$\alpha^* \ge 0, i = 1, \dots, k$$

- Mercer Theorem: Let $K : \mathbb{R}^n \times \mathbb{R}^n \to \mathbb{R}$ be given, then for K to be a valid kernel, it is necessary and sufficient that for any $\{x_1, x_2, \dots, x^m\}$, the corresponding kernel matrix is symmetric positive semi-definite.
- Regularization (revised optimal margin classifier

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

s.t $y^i(w^T x^i + b) \ge 1 - \xi_i, i = 1, 2, \dots, m$
 $\xi_i \ge 0, i = 1, \dots, m$

• Dual of Regularization

$$\max_{\alpha} W(\alpha) = \sum_{i=1}^{m} \alpha_i - \frac{1}{2} \sum_{i,j=1}^{m} y^i y^j \alpha_i \alpha_j (x^i)^T x^j$$
s.t
$$\sum_{i=1}^{m} \alpha_i y^i = 0$$

$$C \ge \alpha_i \ge 0$$
for $i = 1, \dots, m$

Learning Theory

• Union Bound: Let A_1, A_2, \ldots, A_k be k different events (that may not be independent). Then

$$P(A_1 \cup A_2 \dots A_k) \le P(A_1) + P(A_2) + \dots P(A_k)$$

• Hoeffding inequality Let $Z_1, Z_2, \dots Z_m$ be m independent and identically distributed random variables drawn from Bernoulli(ϕ) distribution. Let

$$\hat{\phi} = \frac{1}{m} \sum_{i=1}^{m} Z_i$$
 be the mean of these random variables

and let any $\gamma > 0$ be fixed. Then

$$P(|\phi - \hat{\phi} > \gamma) \leq 2exp(-2\gamma m)$$

• Generational Error Bound: $P(|\epsilon(h_i) - \hat{\epsilon}(h_i)| > \gamma) \le 2exp(-2\gamma^2 m)$

• Uniform Convergence:

$$P(\neg \exists h \in H. | \epsilon(h_i) - \hat{\epsilon}(h_i) | > \gamma)$$

$$= P(\forall h \in H. | \epsilon(h_i) - \hat{\epsilon}(h_i) | \leq \gamma)$$

$$\geq 1 - 2kexp(-2\gamma^2 m)$$

- Solving m, γ, δ We just need to use the equation $\delta = 2kexp(-2\gamma^2m)$ to solve for one variable given the other two.
- Let |H| = k, and let any m, δ , be fixed. Then with probability at least 1δ , we have that

$$\epsilon(\hat{h}) \le (\min_{h \in H} \epsilon(h)) + 2\sqrt{\frac{1}{2m}log(\frac{2k}{\delta})}$$

• Let H be given, and let d = VC(H), then with probability at least $1 - \delta$, we have that for all $h \in H$,

$$|\epsilon(h) - \hat{\epsilon}(h)| \le O(\sqrt{\frac{d}{m}log\frac{m}{d} + \frac{1}{m}log\frac{1}{d}})$$

we also have

$$|\hat{\epsilon}(h) - \epsilon(h^*)| \le O(\sqrt{\frac{d}{m}log\frac{m}{d} + \frac{1}{m}log\frac{1}{d}})$$

Regularization and Model Selection Cross Validation

- Simple Cross Validation: 1. Randomly split S into S_{train} (say 70 percent of the data and S_{cv} . Here, S_{cv} is called the hold-out cross validation set. 2. Train each model M_i on S_{train} only, to get some hypothesis h_i . 3. Select and output the hypothesis h_i that had the smallest error $\hat{\epsilon}(h_i)$ on the hold out crosss validation set.
- k-fold cross validation 1. Randomly split S into k disjoint subsets of m/k training examples each. Lets call these subsets $S_1, ..., S_k$. 3. For each model M_i , we evaluate it as follows: For j = 1, ..., k Train the model M_i on $S_1 \cup \cup S_{j1} \cup S_{j+1} \cup S_k$ (i.e., train on all the data except S_j) to get some hypothesis h_{ij} . Test the hypothesis h_{ij} on S_j , to get $\epsilon_{S_j}(h_{ij})$. The estimated generalization error of model M_i is then calculated as the average of the $\epsilon_{S_j}(h_{ij})$ s (averaged over j).

Feature Selection

• Forward Search:1. Initialize F = ∅. 2. Repeat (a) For i = 1,..., nif i ∉ F, let F_i = F ∪ {i}, and use some version of cross validation to evaluate features F_i. (I.e., train your learning algorithm using only the features in F_i, and estimate its generalization error.) (b) Set F to be the best feature subset found on step (a). 3. Select and output the best feature subset that was evaluated during the entire search procedure.

Bayesian Statistics

•
$$P(\theta|S) = \frac{P(S|\theta)p(\theta)}{p(S)} = \frac{(\prod_{i=1}^{m} p(y^i|x^i,\theta))p(\theta)}{\int_{\theta}(\prod_{i=1}^{m} p(y^i|x^i,\theta)p(\theta))d\theta}$$

•
$$P(y|x,S) = \int_{\theta} p(y|x,\theta)p(\theta|S)d\theta$$

•
$$E(y|x,S) = \int_{\mathcal{Y}} yp(y|x,S)dy$$
.

•
$$\theta_{MAP} = arg \max_{\alpha}$$