## 10-701 Cheat Sheet

### Non-Parametric

notes on chap 6 of PRML).

MaxLikelihood learning window will give you delta functions, which is a kind of over fitting. Use Leave-one-out cross validation for model selection. Idea: Use some of the data to estimate density; Use other part to evaluate how well it works. Pick the parameter that works best.

$$\begin{split} \log p(x_i|\hat{X}\backslash \left\{x_i\right\}) &= \log \frac{1}{n-1} \sum_{j\neq i} k(x_i,x_j), \text{ the sum over all} \\ \text{points is } \frac{1}{n} \sum_{i=1}^n \log \left[\frac{n}{n-1} p(x_i) - \frac{1}{n-1} k(x_i,x_i)\right] \text{ where } p(x) &= \frac{1}{n} \sum_{i=1}^n k(x_i,x). \end{split}$$

why must we not check too many parameters? that you can overfit more; for a given dataset, a few particular parameter values might happen to do well in k-fold CV by sheer chance, where if you had a new dataset they might not do so well. Checking a reasonable number of parameter values makes you less likely to hit those "lucky" spots helps mitigate this risk.

Silverman's Rule for kernel size Use average distance from k nearest neighbors  $r_i = \frac{r}{k} \sum_{x \in \text{NN}(x_i, k)} ||x_i - x||$ .

Watson Nadaraya 1. estimate p(x|y=1) and p(x|y=-1); 2. compute by Bayes rule

$$p(y|x) = \frac{p(x|y)p(y)}{p(x)} = \frac{\frac{1}{my}\sum_{y_i=y}k(x_i,x)\cdot\frac{m_y}{m}}{\frac{1}{m}\sum_ik(x_i,x)}.$$
 3. Decision boundary

 $p(y=1|x) - p(y=-1|x) = \frac{\sum_{j} y_{j} k(x_{j},x)}{\sum_{i} k(x_{i},x)} = \sum_{j} y_{j} \frac{k(x_{j},x)}{\sum_{i} k(x_{i},x)}$ 

Actually, we assume that p(x-y) is equal to  $1/m_y * \sum_{i} k(x_i, x)$ . Using this definition, we can see p(x,-1) + p(x,1) = p(x|-1)p(-1) + p(x|1)p(1) = p(x).This can be incorporated into the regression framework in chap 6 of PRML. Where we define  $f(x-x_n, t \neq t_n) = 0$ , and  $f(x-x_n,t=t_n)=f(x-x_n)$ . Using this definition, we can derive all the probabilities on this slide. (see my handwritten Regression case is the same equation.

**kNN** Let optimal error rate be p. Given unlimited **iid** data. 1NN's error rate is  $\leq 2p(1-p)$ .

### Matrix Cookbook

$$\begin{array}{l} \frac{\partial x^T a}{\partial x} = \frac{\partial a^T x}{\partial x} = a \\ \frac{\partial a^T X b}{\partial X} = ab^T, \ \frac{\partial a^T X^T b}{\partial X} = ba^T, \ \frac{\partial a^T (X^T | X) a}{\partial X} = aa^T \\ W \in S, \ \frac{\partial}{\partial s} (x - As)^T W (x - As) = -2A^T W (x - As), \\ \frac{\partial}{\partial x} (x - s)^T W (x - s) = 2W (x - s), \\ \frac{\partial}{\partial s} (x - s)^T W (x - s) = -2W (x - s), \\ \frac{\partial}{\partial s} (x - As)^T W (x - As) = 2W (x - As), \\ \frac{\partial}{\partial a} (x - As)^T W (x - As) = -2W (x - As)s^T. \\ \text{Tr}(A) = \sum_i A_{ii}. \text{ For two equal sized matrices, } \text{Tr}(A^T B) = \\ \text{Tr}(B^T A) = \text{Tr}(AB^T) = \text{Tr}(BA^T) = \sum_{i,j} A_{ij} B_{ij}. \\ \text{Tr}(A) = \text{Tr}(A^T), \text{Tr}(A + B) = \text{Tr}(A) + \text{Tr}(B). \text{ For square matricis,} \\ \text{Tr}(AB) = \text{Tr}(BA), \text{Tr}(ABC) = \text{Tr}(CAB) = \text{Tr}(BCA) \\ \text{(trace rotation)}. \end{array}$$

# Classifers and Regressors

Naive Bayes Conditionally independent:

 $P(x_1, x_2, \dots | C) = \prod_i P(x_i | C)$ . One way to avoid divide by zero: add  $(1, 1, \ldots, 1)$  and  $(0, 0, \ldots, 0)$  to both classes.

zero: add 
$$(1,1,\ldots,1)$$
 and  $(0,0,\ldots,0)$  to both classes.  
**Learns**  $P(x_i|y)$  for  $Discrete\ x_i - P(x_i|y) = \frac{\#D(X_i=x_i,Y=y)}{\#D(Y=y)}$ 

For smoothing, use  $P(x_i|y) = \frac{\#D(X_i=x_i,Y=y)+\#D(Y=y)}{\#D(Y=y)+n_ik}$ , where  $n_i$  is the number of different possible values for  $X_i$  (In practice problem set, Jing Xiang used k = 1?) Continuous  $x_i$  – Can use any PDF, but usually use Gaussian

 $P(x_i|y) = \mathcal{N}(\mu_{X_i|y}, \sigma_{X_i|y}^2)$ , where  $\mu_{X_i|y}$  and  $\sigma_{X_i|y}$  are, respectively, the average and variance of  $X_i$  for all data points where Y = y. The Gaussian distribution already provides

Perceptron Produces linear decision boundaries. Classifies using  $\hat{y} = X_{test} w + b$  Learns w and b by updating w whenever Copyright © 2013 Yimeng Zhang.

 $y_i(w^Tx_i+b) \leq 0$  (i.e. incorrectly classified). Updates as  $w \leftarrow w + x_i y_i, b \leftarrow b + y_i$  Repeat until all examples are correctly classified. w is some linear combination  $\sum_{i} \alpha_{i} x_{i} (y_{i} * x_{i})$  of data points, and decision boundary is the linear hyperplane  $f(x) = w^T x + b$ . Note that the perceptron is the same as stochastic gradient descent with a hinge loss function of  $max(0, 1 - y_i [< w, x_i > +b])$  (we can't remove 1 in the loss function; otherwise we can set w, b = 0). Convergence of perceptron proof 1 Here we use a perceptron without b. Assume we have  $w^*$  that has margin  $\gamma$  $(\min(\boldsymbol{w}^*)^T y_i x_i = \gamma)$ , and  $\|\boldsymbol{w}^*\| = 1, \|x_i\| = 1$ . We start from  $\mathbf{w}_0 = 0$ . Assume that we have made M mistakes. We have 1)  $\boldsymbol{w}_{M} \cdot \boldsymbol{w}^{*} = (\boldsymbol{w}_{M-1} + y_{i}x_{i}) \cdot \boldsymbol{w}^{*} > \boldsymbol{w}_{M-1} \cdot \boldsymbol{w}^{*} + \gamma$ . So we have  $\boldsymbol{w}_M \cdot \boldsymbol{w}^* \ge M \gamma$ . 2)  $\mathbf{w}_{M} \cdot \mathbf{w}_{M} = (\mathbf{w}_{M-1} + y_{i}x_{i}) \cdot (\mathbf{w}_{M-1} + y_{i}x_{i}) = \mathbf{w}_{M-1}$  $\mathbf{w}_{M-1} + 2y_i x_i \cdot \mathbf{w}_{M-1} + (y_i x_i) \cdot (y_i x_i) \le \mathbf{w}_{M-1} \cdot \mathbf{w}_{M-1} + 1.$ So we have  $\boldsymbol{w}_M \cdot \boldsymbol{w}_M < M$ .

Combining them, using Cauchy-Schwarz, we have

Cauchy-Schwarz. So we have  $\sqrt{M} - M\gamma \ge 0$ .

### Kernel

Kernel function  $k(\boldsymbol{x}, \boldsymbol{x}') = \phi(\boldsymbol{x})^T \phi(\boldsymbol{x}')$  for some  $\phi(\cdot)$ . For a set of data points  $\{x_i\}$ , we have Gram matrix (kernel matrix)  $K_{ij} = k(\boldsymbol{x}_i, \boldsymbol{x}_j)$ . A necessary and sufficient condition for being a valid kernel function: K always positive semidefinite. Proof:  $\alpha^T K \alpha = \sum_{ij} \alpha_i \alpha_j K_{ij} = \sum_{ij} \alpha_i \alpha_j \langle \phi(\mathbf{x}_i), \phi(\mathbf{x}_j) \rangle =$  $\langle \sum_{i} \alpha_{i} \phi(\boldsymbol{x}_{i}), \sum_{j} \alpha_{j} \phi(\boldsymbol{x}_{j}) \rangle \geq 0.$ 

 $M\gamma < w_M \cdot w^* < ||w_M|| ||w^*|| < \sqrt{M}$ . So  $M < 1/\gamma^2$ . proof 2

Let potential function  $Q_i = ||\boldsymbol{w}_i|| - \boldsymbol{w}_i \cdot \boldsymbol{w}^*$ , where i is the

mistakes, so we have  $Q_i < \sqrt{M} - M\gamma$ . Clearly  $Q_i > 0$  by

number of iterations. Assuming up to iteration i, we have M

Mercer's Theorem for any symmtric function  $k: \mathcal{X} \times \mathcal{X} \to \mathbb{R}$  which is square integrable and satisfying  $\int_{\mathcal{X}\times\mathcal{X}} k(x,x')f(x)f(x')\mathrm{d}x\mathrm{d}x' \geq 0$  for  $f\in L_2(\mathcal{X})$ , we have a feature space  $\Phi(x)$  and  $\lambda > 0$  that  $k(x, x') = \sum_i \lambda_i \phi_i(x) \phi_i(x')$ .