

Bioinformatics and Biomarker Discovery Part 1: Foundations

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Themes of Bioinformatics



Bioinformatics =
Data Mgmt +
Knowledge Discovery +
Sequence Analysis +
Physical Modeling +

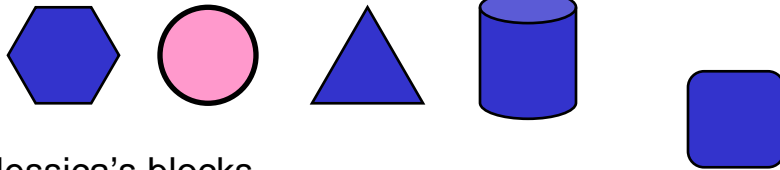
Knowledge Discovery =
Statistics + Algorithms + Databases

Applications include diagnosis, prognosis, &
treatment optimization, often thru biomarker
discovery

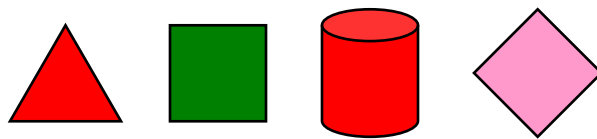
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What is Knowledge Discovery?

Jonathan's blocks



Jessica's blocks



Whose block is this?

Jonathan's rules : Blue or Circle
Jessica's rules : All the rest

What is Knowledge Discovery?



Question: Can you explain how?

Key Steps of Knowledge Discovery

- **Training data gathering**
- **Feature generation**
 - k-grams, colour, texture, domain know-how, ...
- **Feature selection**
 - Entropy, χ^2 , CFS, t-test, domain know-how...
- **Feature integration**
 - SVM, ANN, PCL, CART, C4.5, kNN, ...

What is Accuracy?

What is Accuracy?

	predicted as positive	predicted as negative
positive	TP	FN
negative	FP	TN

$$\begin{aligned} \text{Accuracy} &= \frac{\text{No. of correct predictions}}{\text{No. of predictions}} \\ &= \frac{\text{TP} + \text{TN}}{\text{TP} + \text{TN} + \text{FP} + \text{FN}} \end{aligned}$$

Examples (Balanced Population)

classifier	TP	TN	FP	FN	Accuracy
A	25	25	25	25	50%
B	50	25	25	0	75%
C	25	50	0	25	75%
D	37	37	13	13	74%

- Clearly, B, C, D are all better than A
- Is B better than C, D?
- Is C better than B, D?
- Is D better than B, C?

Accuracy may not tell the whole story

Examples (Unbalanced Population)

classifier	TP	TN	FP	FN	Accuracy
A	25	75	75	25	50%
B	0	150	0	50	75%
C	50	0	150	0	25%
D	30	100	50	20	65%

- Clearly, D is better than A
- Is B better than A, C, D?

Exercise: What is B's Prediction strategy?

What is Sensitivity (aka Recall)?

	predicted as positive	predicted as negative
positive	TP	FN
negative	FP	TN

$$\begin{aligned}
 \text{Sensitivity}_{\text{wrt positives}} &= \frac{\text{No. of correct positive predictions}}{\text{No. of positives}} \\
 &= \frac{TP}{TP + FN}
 \end{aligned}$$

Sometimes sensitivity wrt negatives is termed **specificity**

What is Precision?

	predicted as positive	predicted as negative
positive	TP	FN
negative	FP	TN

$$\begin{aligned} \text{Precision} &= \frac{\text{No. of correct positive predictions}}{\text{No. of positives predictions}} \\ &= \frac{TP}{TP + FP} \end{aligned}$$

Unbalanced Population Revisited

classifier	TP	TN	FP	FN	Accuracy	Sensitivity	Precision
A	25	75	75	25	50%	50%	25%
B	0	150	0	50	75%	0%	ND
C	50	0	150	0	25%	100%	25%
D	30	100	50	20	65%	60%	38%

- What are the sensitivity and precision of B and C?
- Is B better than A, C, D?

Abstract Model of a Classifier

- Given a test sample S
- Compute scores $p(S), n(S)$
- Predict S as negative if $p(S) < t * n(s)$
- Predict S as positive if $p(S) \geq t * n(s)$

t is the decision threshold of the classifier

changing t affects the recall and precision,
and hence accuracy, of the classifier

An Example

S	$P(S)$	$N(S)$	Actual Class	Predicted Class $\emptyset \ t = 3$	Predicted Class $\emptyset \ t = 2$
2	0.961252	0.038748	P	P	P
3	0.435302	0.564698	N	N	N
6	0.691596	0.308404	P	N	P
7	0.180885	0.819115	N	N	N
8	0.814909	0.185091	P	P	P
10	0.887220	0.112780	P	P	P
			accuracy	5 / 6	6 / 6
			recall	3 / 4	4 / 4
			precision	3 / 3	4 / 4

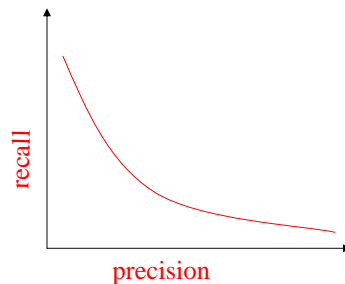
Recall that ...

- Predict S as negative if $p(S) < t * n(s)$
- Predict S as positive if $p(S) \geq t * n(s)$

Precision-Recall Trade-off



- A predicts better than B if A has better recall and precision than B
- There is a trade-off between recall and precision
- In some applications, once you reach a satisfactory precision, you optimize for recall
- In some applications, once you reach a satisfactory recall, you optimize for precision



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Comparing Prediction Performance



- **Accuracy is the obvious measure**
 - But it conveys the right intuition only when the positive and negative populations are roughly equal in size
- **Recall and precision together form a better measure**
 - But what do you do when A has better recall than B and B has better precision than A?

So let us look at some alternate measures

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Adjusted Accuracy

- Weigh by the importance of the classes

$$\text{Adjusted accuracy} = \alpha * \text{Sensitivity} + \beta * \text{Specificity}$$

$$\text{where } \alpha + \beta = 1$$

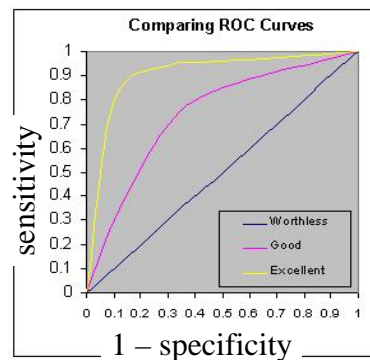
$$\text{typically, } \alpha = \beta = 0.5$$

classifier	TP	TN	FP	FN	Accuracy	Adj Accuracy
A	25	75	75	25	50%	50%
B	0	150	0	50	75%	50%
C	50	0	150	0	25%	50%
D	30	100	50	20	65%	63%

But people can't always agree on values for α , β

ROC Curves

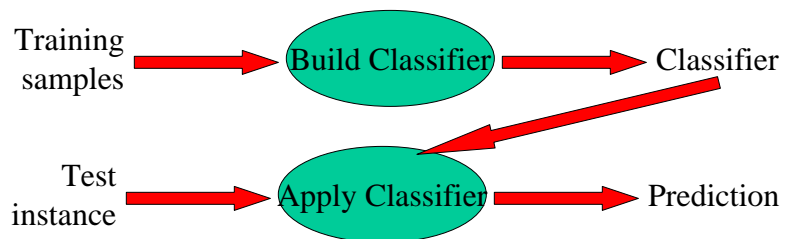
- By changing t , we get a range of sensitivities and specificities of a classifier
- A predicts better than B if A has better sensitivities than B at most specificities
- Leads to ROC curve that plots sensitivity vs. (1 - specificity)
- Then the larger the area under the ROC curve, the better



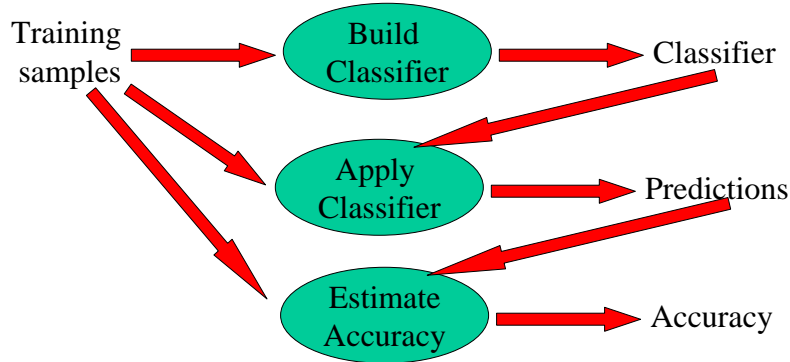
What is Cross Validation?



Construction of a Classifier



Estimate Accuracy: Wrong Way



Exercise: Why is this way of estimating accuracy wrong?

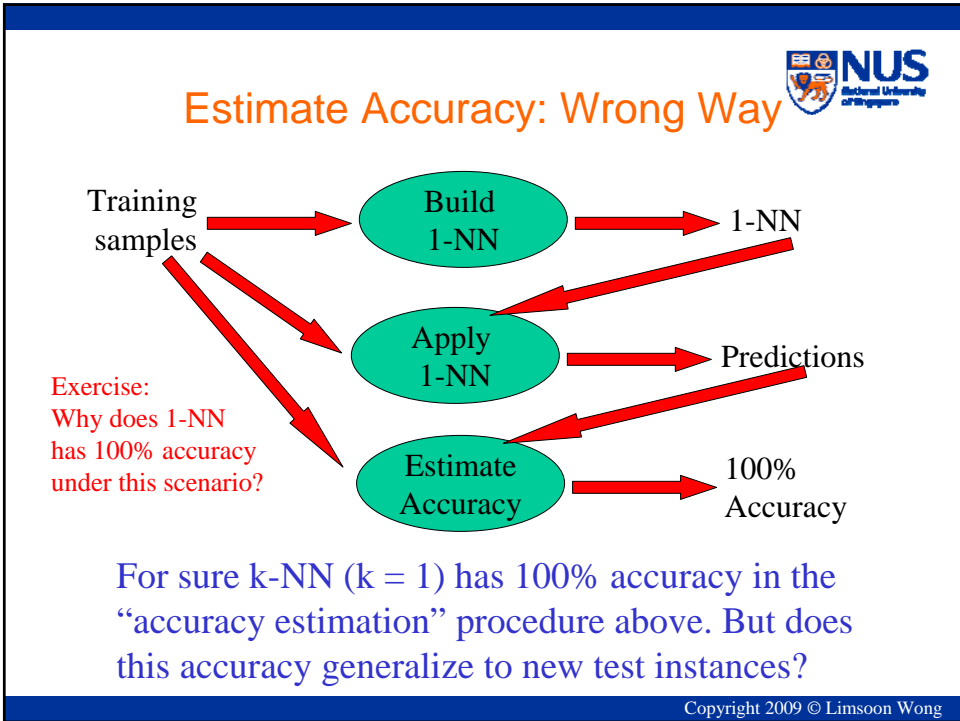
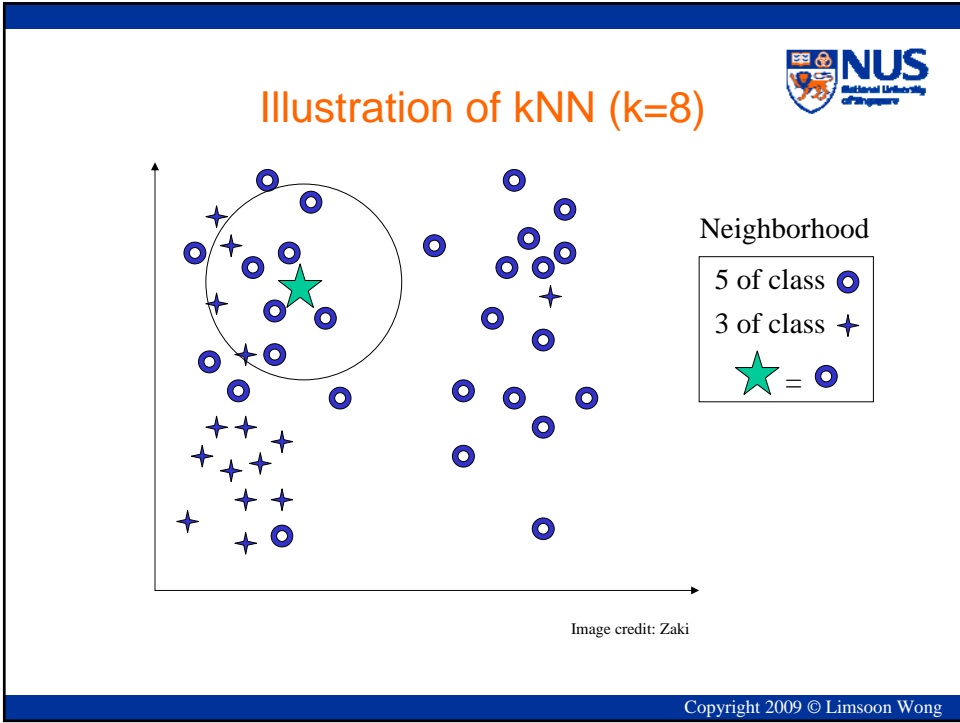
K-Nearest Neighbour Classifier (k-NN)

- Assume S is well approximated by its neighbours
- Then, given a sample S , find the k observations $S_1 \dots S_k$ in the known data that are “closest” to it, and average their responses

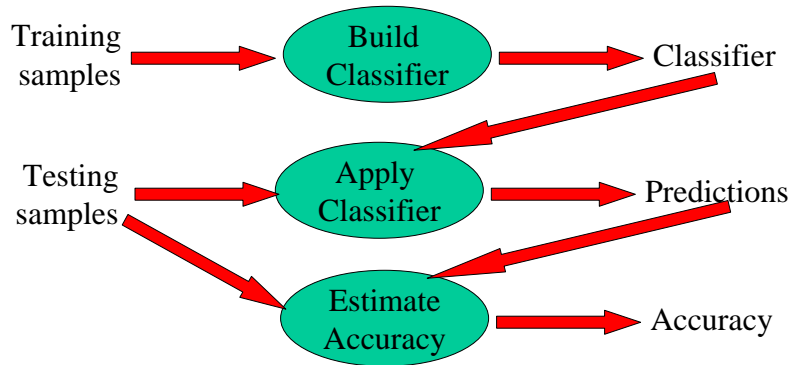
$$p(S) = \sum_{S_i \in N_k(S) \cap D^p} 1 \quad n(S) = \sum_{S_i \in N_k(S) \cap D^N} 1$$

where $N_k(S)$ is the neighbourhood of S defined by the k nearest samples to it.

Assume distance between samples is Euclidean distance for now



Estimate Accuracy: Right Way



Testing samples are NOT to be used during “Build Classifier”

How Many Training and Testing Samples?

- No fixed ratio between training and testing samples; but typically 2:1 ratio
- Proportion of instances of different classes in testing samples should be similar to proportion in training samples
- What if there are insufficient samples to reserve 1/3 for testing?
- Ans: Cross validation

Cross Validation

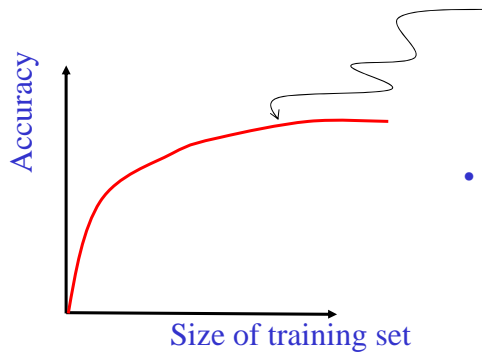


- Divide samples into k roughly equal parts
- Each part has similar proportion of samples from different classes
- Use each part to test other parts

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How Many Fold?

- If samples are divided into k parts, we call this k -fold cross validation
- Choose k so that
 - the k -fold cross validation accuracy does not change much from $k-1$ fold
 - each part within the k -fold cross validation has similar accuracy
- $k = 5$ or 10 are popular choices for k



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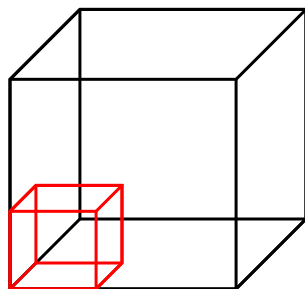
Curse of Dimensionality



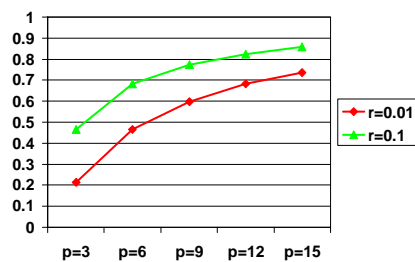
Curse of Dimensionality



- How much of each dimension is needed to cover a proportion r of total sample space?



- Calculate by $e_p(r) = r^{1/p}$
- So, to cover 10% of a 15-D space, need to sample $(0.1)^{1/15} = 85\%$ of each dimension!



Exercise: Why $e_p(r) = r^{1/p}$?

Consequence of the Curse



- Suppose the number of samples given to us in the total sample space is fixed
- Let the dimension increase
- Then the distance of the k nearest neighbours of any point increases
- Then the k nearest neighbours are less and less useful for prediction, and can confuse the k -NN classifier (and other types of classifiers as well)

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What is Feature Selection?

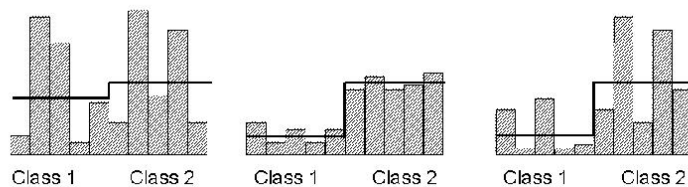


Tackling the Curse

- Given a sample space of p dimensions
- It is possible that some dimensions are irrelevant
- Need to find ways to separate those dimensions (aka features) that are relevant (aka signals) from those that are irrelevant (aka noise)

Signal Selection (Basic Idea)

- Choose a feature w/ low intra-class distance
- Choose a feature w/ high inter-class distance



Signal Selection (e.g., t-statistics)



The t-stats of a signal is defined as

$$t = \frac{|\mu_1 - \mu_2|}{\sqrt{(\sigma_1^2/n_1) + (\sigma_2^2/n_2)}}$$

where σ_i^2 is the variance of that signal in class i , μ_i is the mean of that signal in class i , and n_i is the size of class i .

Suggestion a modification to t-stats when n_1 and n_2 are small.

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Self-fulfilling Oracle



- Construct artificial dataset with 100 samples, each with 100,000 randomly generated features and randomly assigned class labels
- Select 20 features with the best t-statistics (or other methods)
- Evaluate accuracy by cross validation using only the 20 selected features
- The resultant estimated accuracy can be ~90%
- But the true accuracy should be 50%, as the data were derived randomly

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What Went Wrong?



- The 20 features were selected from the whole dataset
- Information in the held-out testing samples has thus been “leaked” to the training process
- The correct way is to re-select the 20 features at each fold; better still, use a totally new set of samples for testing

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Confounding Factors



A True Story (Charig et al, 1986)



Treatment A	Treatment B
78% (273/350)	83% (289/350)

	Treatment A	Treatment B
Small Stones	Group 1 93% (81/87)	Group 2 87% (234/270)
Large Stones	Group 3 73% (192/263)	Group 4 69% (55/80)
Both	78% (273/350)	83% (289/350)

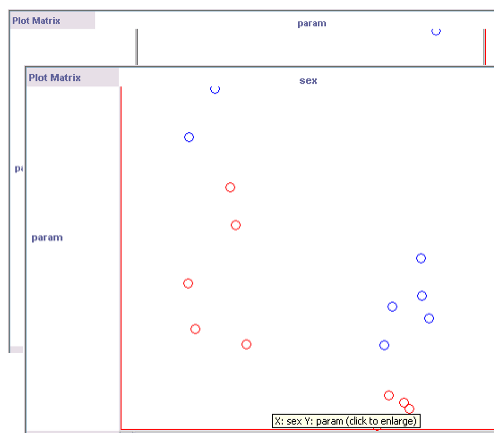
- Treatment B seems more effective than Treatment A for kidney stone
- Now Treatment A seems more effective than Treatment B
- **Case of Simpson Paradox: But we won't know this if we don't capture stone size info**

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A Made-Up Example



param	class	sex
100	alive	m
99	alive	m
98	alive	m
95	alive	m
90	alive	m
71	dead	m
80	dead	m
85	dead	m
73	dead	m
75	dead	m
71	alive	f
75	alive	f
73	alive	f
76	alive	f
72	alive	f
60	dead	f
65	dead	f
62	dead	f
63	dead	f
64	dead	f



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Concluding Remarks



What have we learned?



- **Methodology of data mining**
 - Feature generation, feature selection, feature integration
- **Evaluation of classifiers**
 - Accuracy, sensitivity, precision
 - Cross validation
- **Curse of dimensionality**
 - Feature selection concept
 - Self-fulfilling oracle

Any Questions?



Acknowledgements



- The first two slides were shown to me 10+ years ago by Tan Ah Hwee

References



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