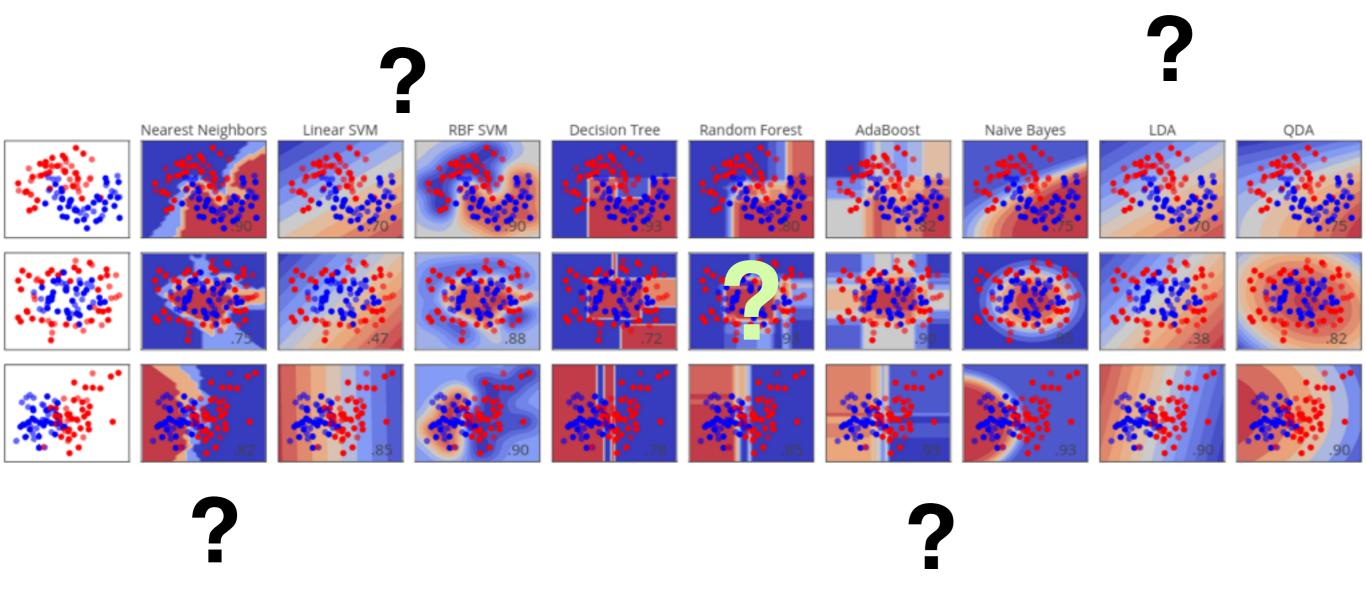


Algorithm Selection

Two-day workshop
Duke University
Adam Chekroud

Preface

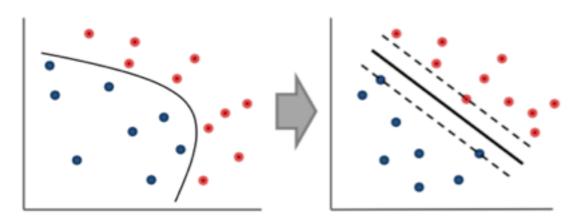


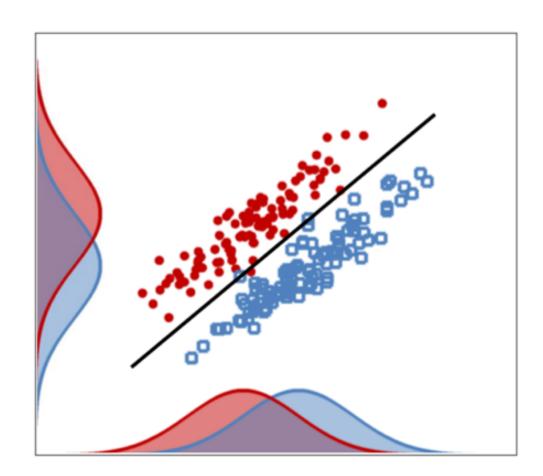
"Algorithm Selection: Try all of them*"

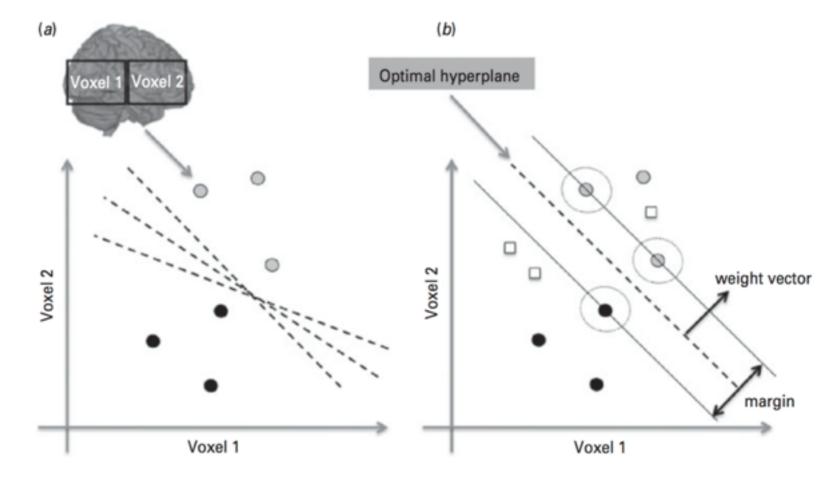
^{*} only caveat: with appropriate train-test precautions

Objectives

Nonlinear SVM



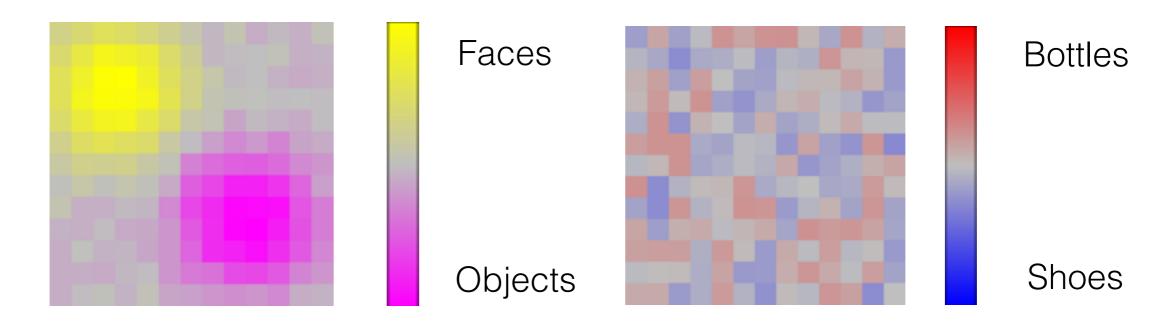




(Supervised) Machine Learning

Working definition: using computer algorithms to identify patterns in data that help us **predict** things we care about

Most accurate mapping from observables to an outcome



Neuroscience has readily adopted the notion of "distributed processing"

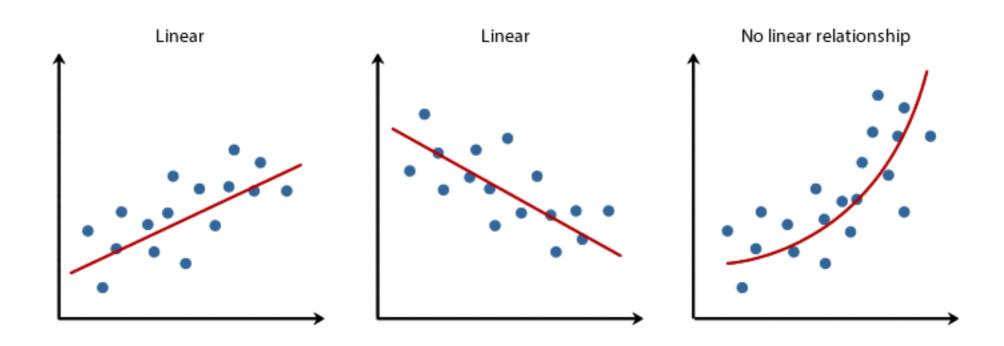
—> Multivariate approach is more sensitive in capturing complex signal

Supervised Learning

Very different to traditional inference!

Often, we must leave interpretability at the door...

While regression was all about estimating beta...

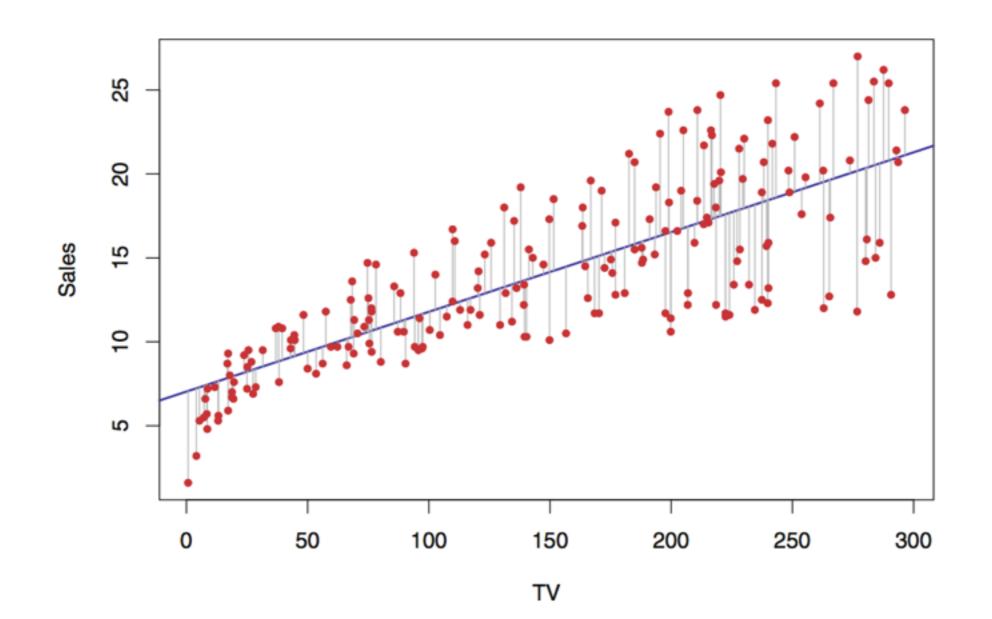


....ML is all about y-hat

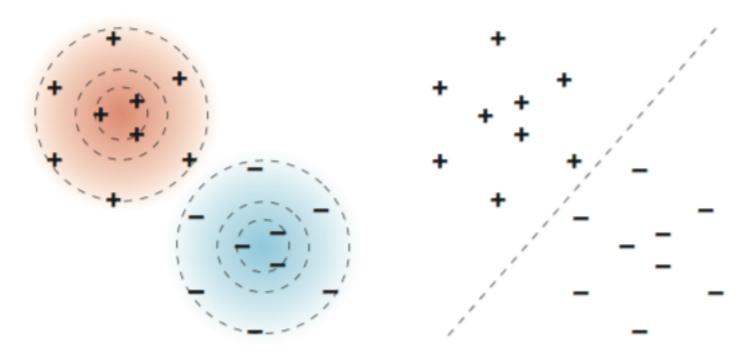
Supervised Learning

Algorithms are, for the most part, data agnostic:

 Medical diagnosis, computer vision, spam filtering, neuroscience, genetics, tailoring ads, selling you stuff, filling ur amazon wishlist.



How do we choose an algorithm?



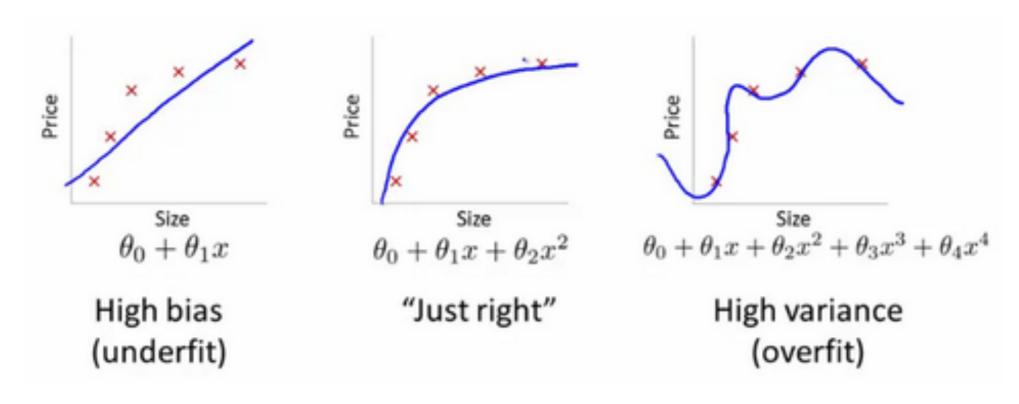
ML algorithms come with different "flavours"

Three broad aspects to consider (at least)

- Accuracy
- Feasibility
- Linearity (or, the nature of the problem)

<u>Accuracy</u>

- More accuracy is not always better (!!)
- Is an approximation sufficient?
- Is an approximation better?!
 - Simpler models less likely to over fit

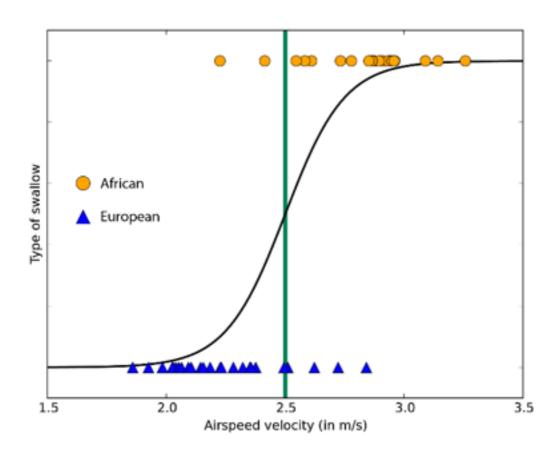


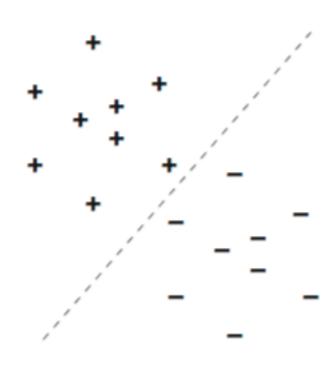
Feasibility

- What can your data support? (realistic!)
 - 10 subjects? 10k? 100k?
 - Powerful algorithms are usually data hungry
- What can you afford, computationally?
 - Not just one model
 - Tuning parameters, cross validation...
 - Implementation?
 - Laptop is easy but slow
 - HPC/AWS, fast, but hard

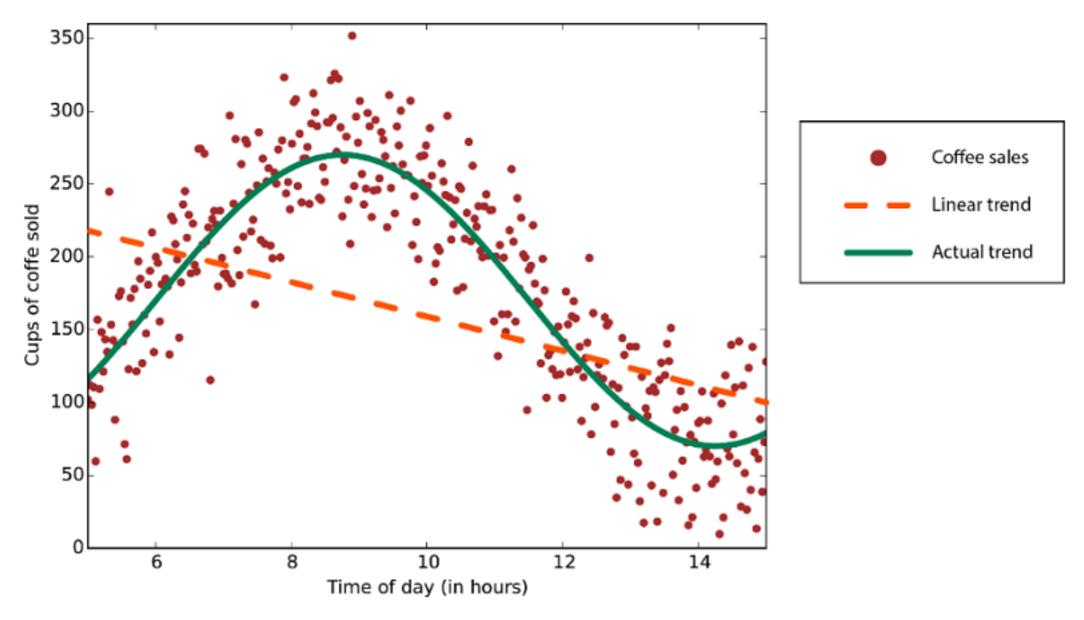
Linearity (or approach)

- Is the problem likely to have linear solution?
- Was your problem just finding the key variable?





Did you just need more flexibility?

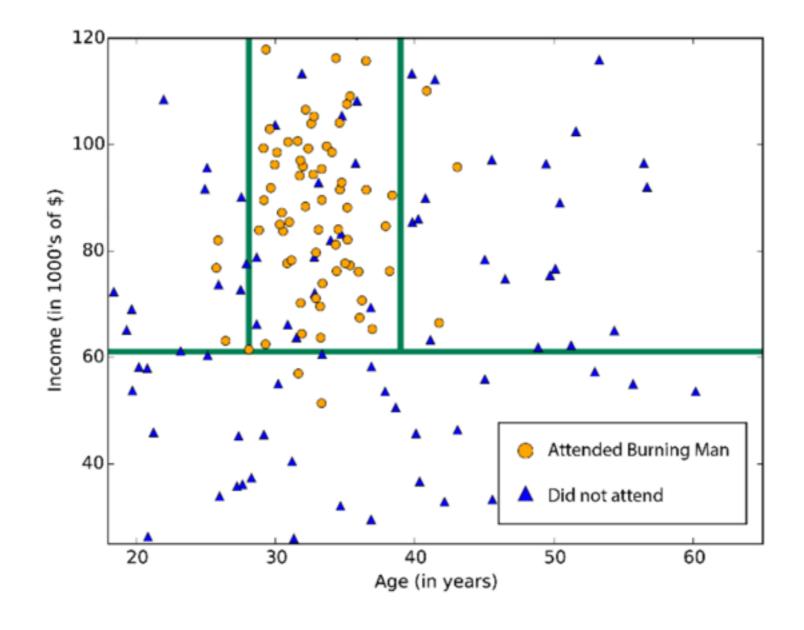


Data with a nonlinear trend - using a linear regression method would generate much larger errors than necessary

Despite their dangers, linear algorithms are very popular as a first line of attack. They tend to be algorithmically simple and fast to train.

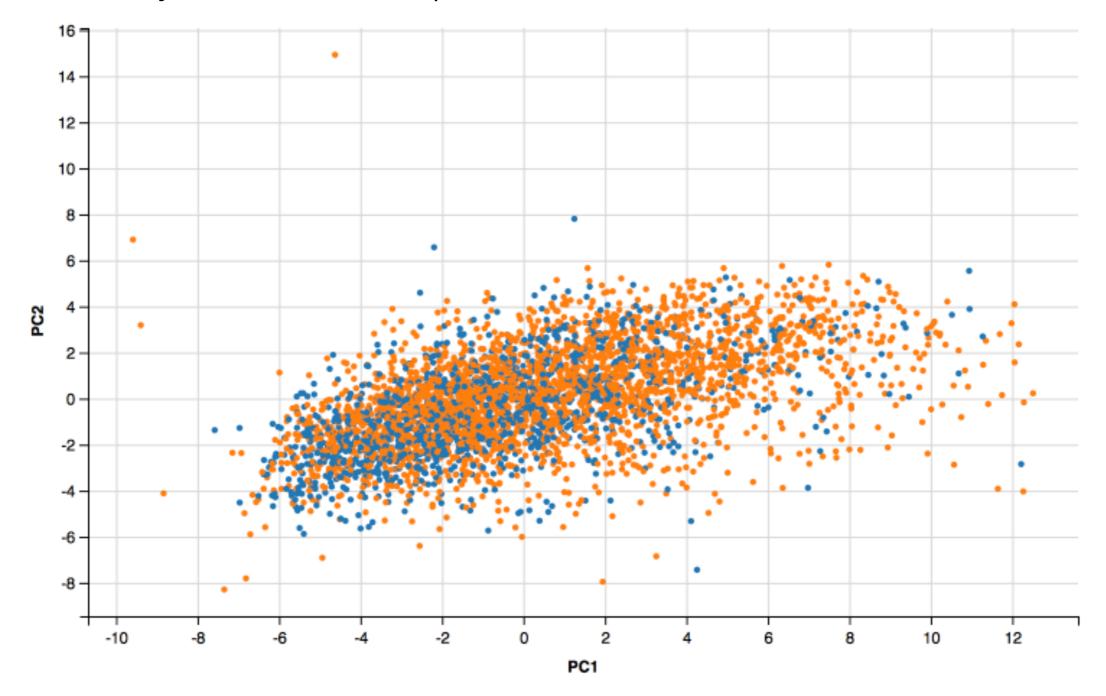
Linearity (or approach)

How about a non-parametric approach?



Linearity (or approach)

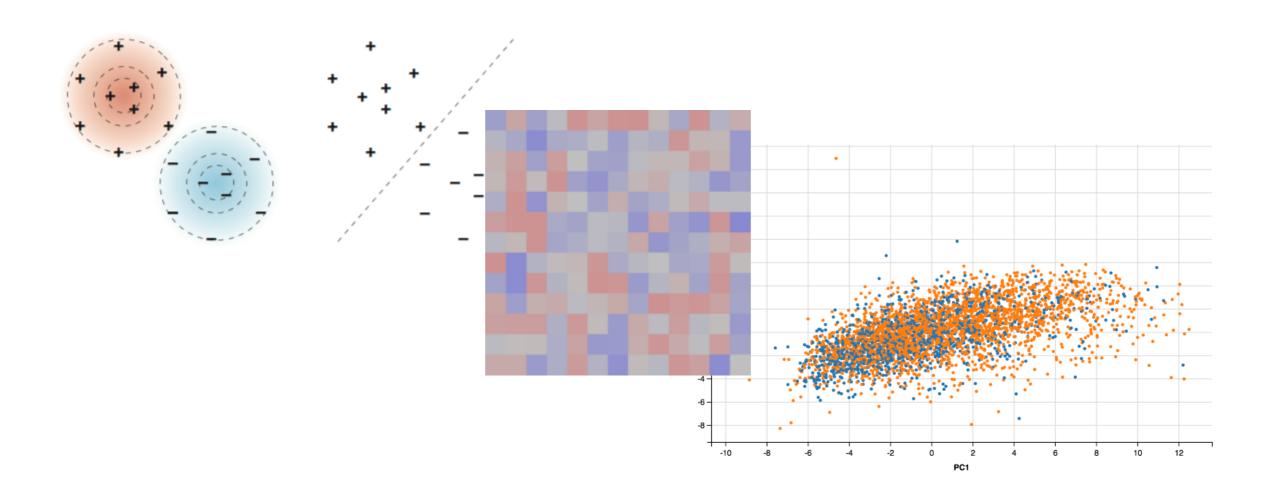
• If only it was that simple...



Inductive Bias: Algorithm selection necessarily introduces a bias

- We can only approximate a correct output
- Not necessarily a bad thing
 - That is how we learn!

...but we want to be as less wrong as possible.



Inductive Bias (reading)

Algorithm	Inductive Bias
Candidate-Elimination	The target concept c is contained in the hypothesis space H.
Linear Regression	The relationship between the attributes x and the output y is linear. The goal is to minimize the sum of squared errors.
Decision Trees	Shorter trees are preferred over longer trees. Trees that place high information gain attributes close to the root are preferred over those that do not.
Neural Networks with Backpropagation	Smooth interpolation between data points.
K-Nearest Neighbors	The classification of an instance x will be most similar to the classification of other instances that are nearby in Euclidean distance.
Support Vector Machines	Distinct classes tend to be separated by wide margins.
Naive Bayes	Each input depends only on the output class or label; the inputs are independent from each other.

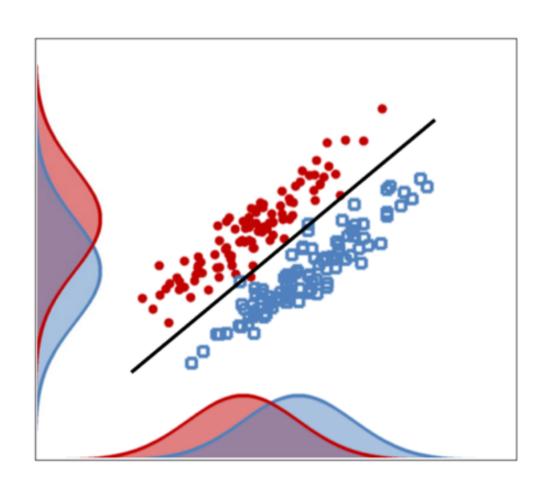
Interim summary

Working definition: using computer algorithms to identify patterns in data that help us **predict** things we care about

• Pragmatic approach, different to classic inference

Try them all, but think about:

- Accuracy
- Feasibility
- Inductive bias



What are the options?

ML framework called *caret*

Collection of functions that streamline ML processes

Contains tools for

- Data splitting
- Pre-processing
- Feature selection
- Model tuning
- Variable importance estimation
- Model validation
- Many models already integrated into this framework:

Boosted Classification Trees Bagged AdaBoost AdaBoost.M1 Adaptive Mixture Discriminant Analysis Adaptive-Network-Based Fuzzy Inference System Model Averaged Neural Network Naive Bayes Classifier with Attribute Weighting Tree Augmented Naive Bayes Classifier with Attribute Weighting Bagged Model Bagged MARS Bagged Flexible Discriminant Analysis Bagged Flexible Discriminant Analysis Bagged FDA using gCV Pruning Bayesian Additive Regression Trees Bayesian Generalized Linear Model Self-Organizing Map Binary Discriminant Analysis Boosted Tree The Bayesian lasso Bayesian Ridge Regression (Model Averaged) Random Forest with Additional Feature Selection Bayesian Regularized Neural Networks Boosted Linear Model Boosted Smoothing Spline Boosted Tree C5.0 Cost-Sensitive C5.0 Single C5.0 Ruleset Single C5.0 Tree Conditional Inference Random Forest CHi-squared Automated Interaction Detection SIMCA Conditional Inference Tree Cubist	Model		
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http://topepo.github.io/caret/modelList.html

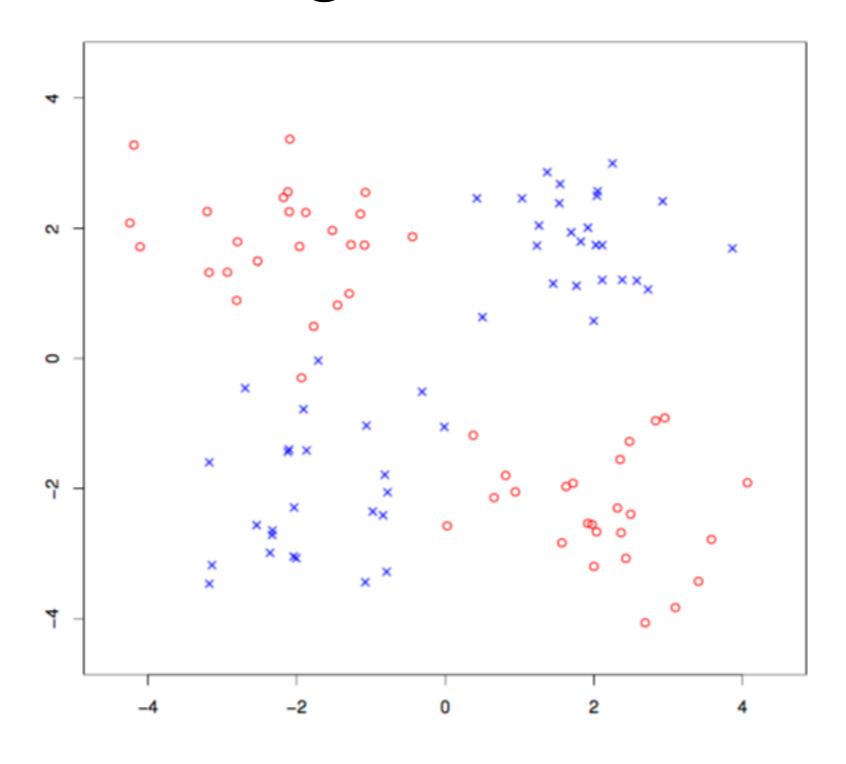
What are the options?

High-level overview

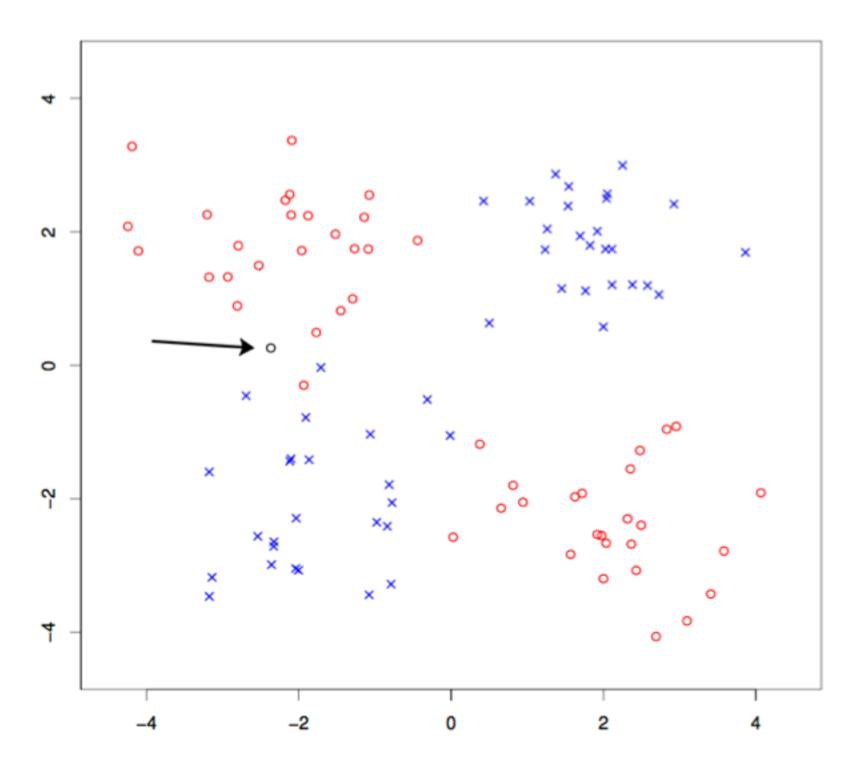
- No stats
- No equations
- No jargon**
- All "off the shelf"

Cover some popular algorithms

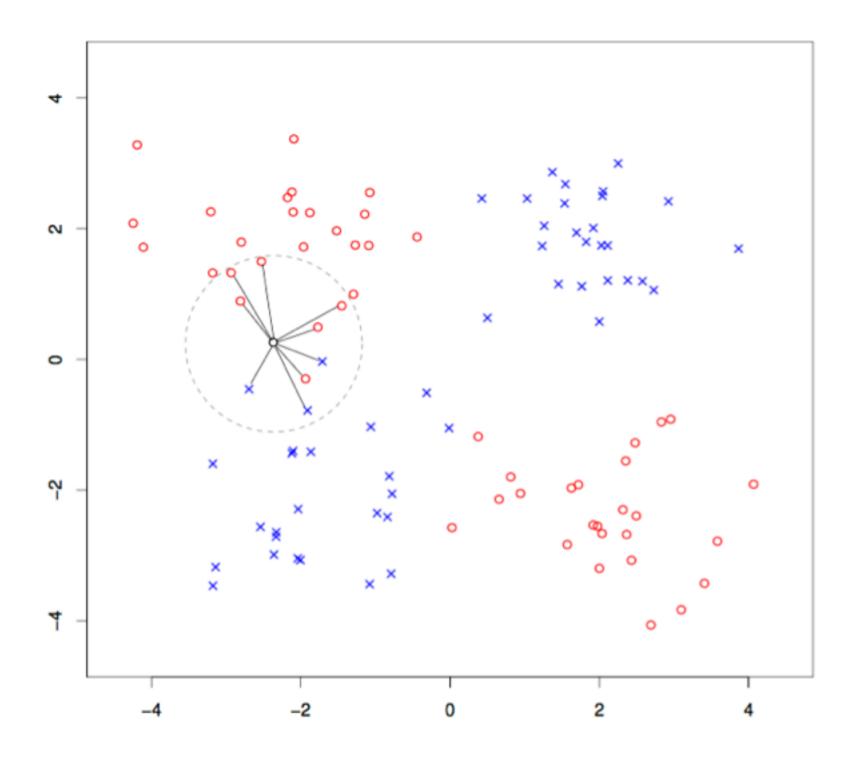
- k-Nearest Neighbours
- Support vector machine
 - linear
 - non-linear (radial)
- Decision tree
- Penalized regression



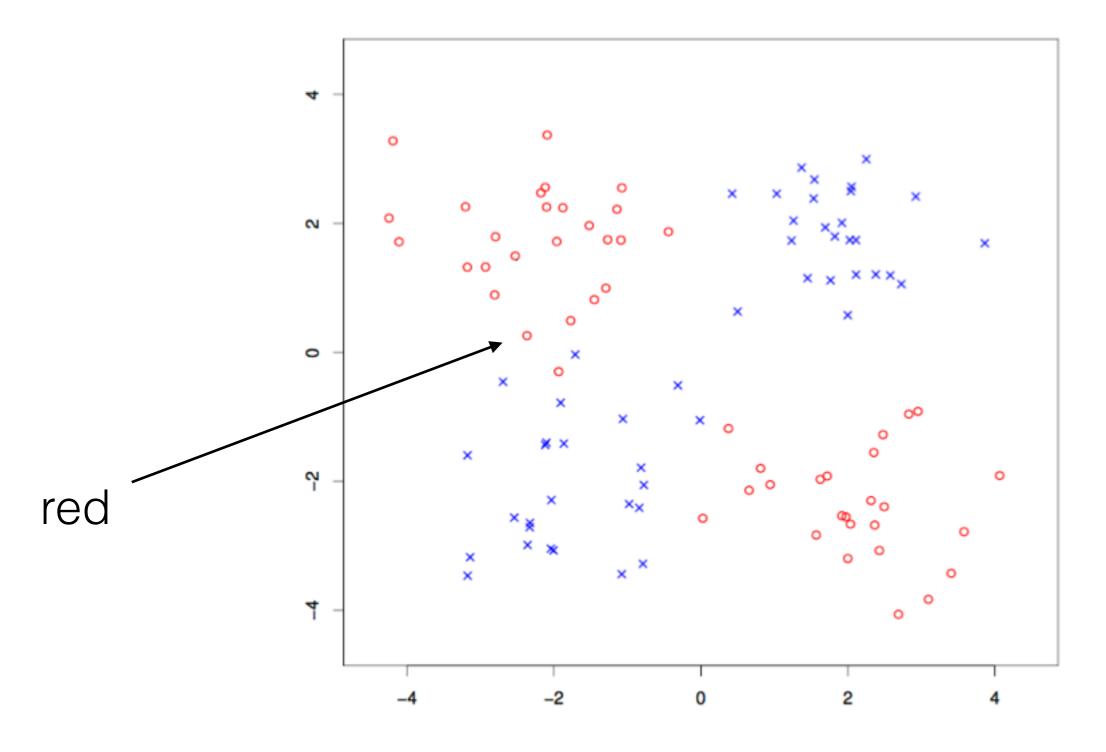
1 - Get data



2 - bring a new observation



3 - find the k most similar observations to that one



4 - prediction is the most common outcome in the circle

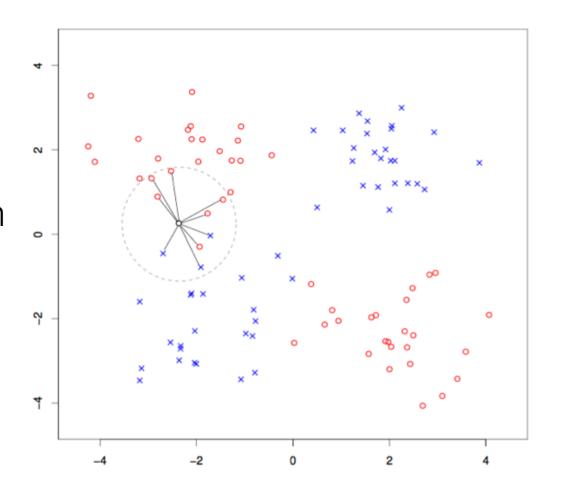
k-NN - finer details (optional)

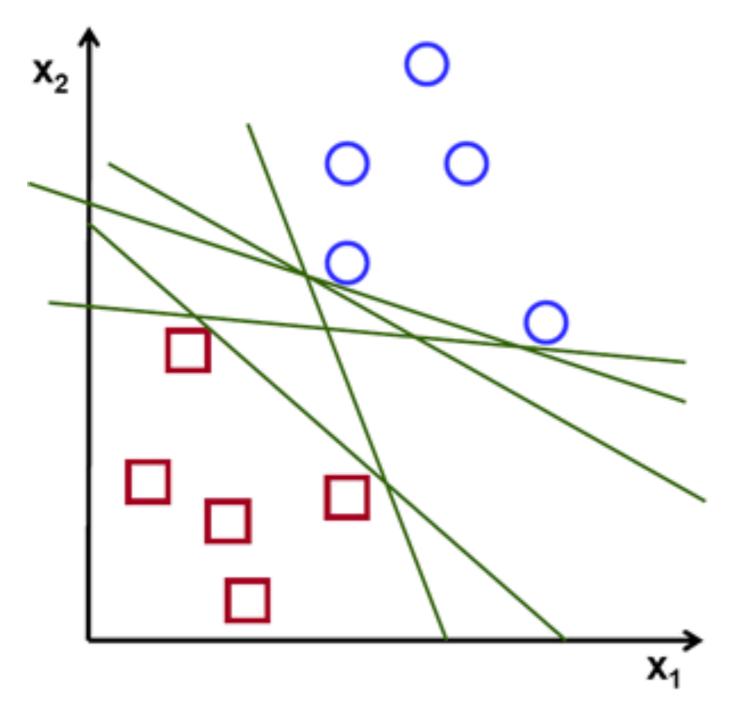
Defining neighbours

- Which "distance measure"?
 - Pearson/spearman correlation
- How many neighbours?
 - Choose k through cross-validation
- Are all neighbours equal?
 - Weighted circle of neighbours

Dimensionality?

PCA first, then k-NN



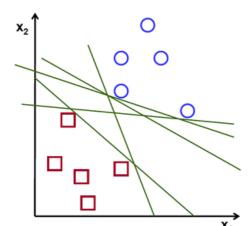


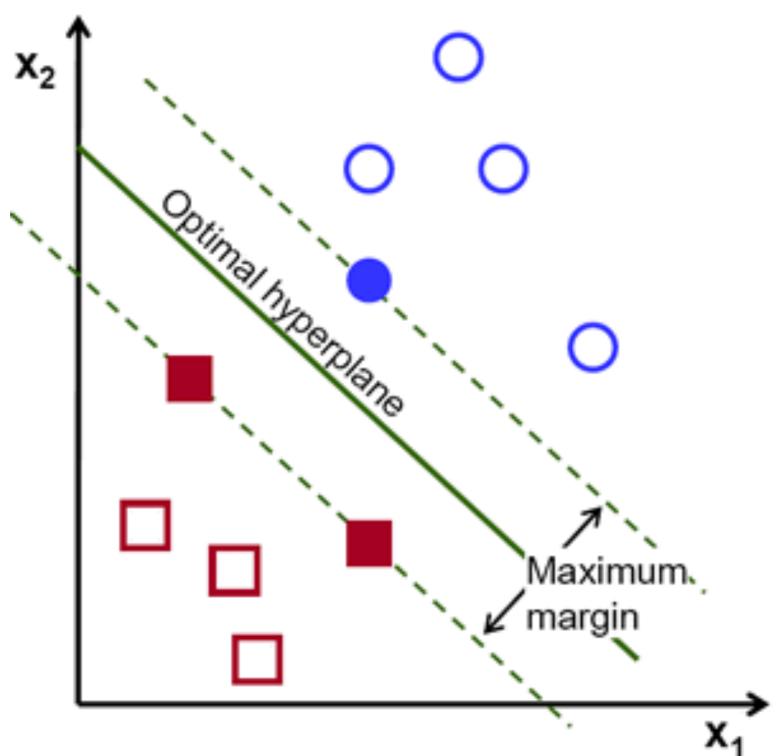
Aim: find a boundary that (optimally) separates our classes

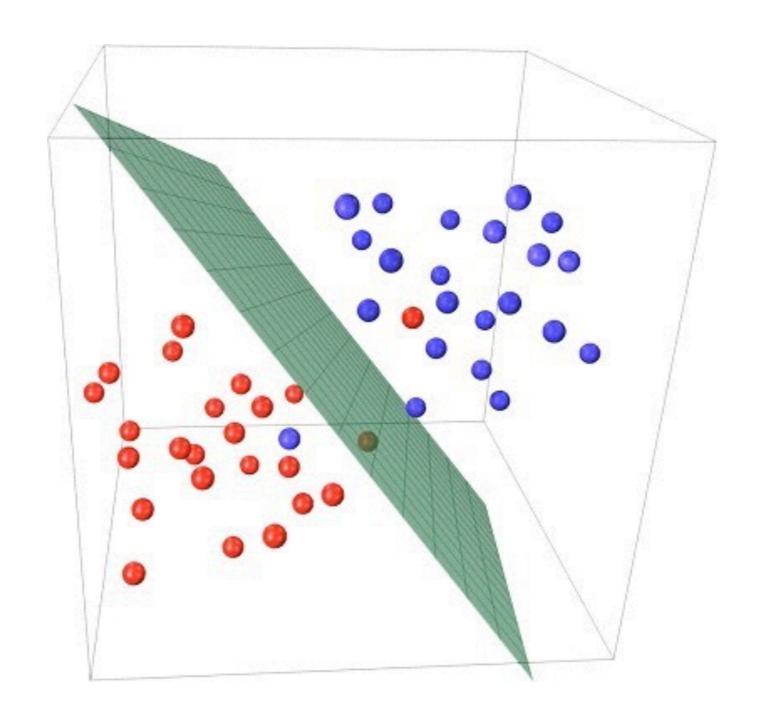
Support Vector Machine Class 2 Class 2 Class 1 Class 1

 Max-margin theory: boundary should be as far away from the data as possible

(Bad boundaries)

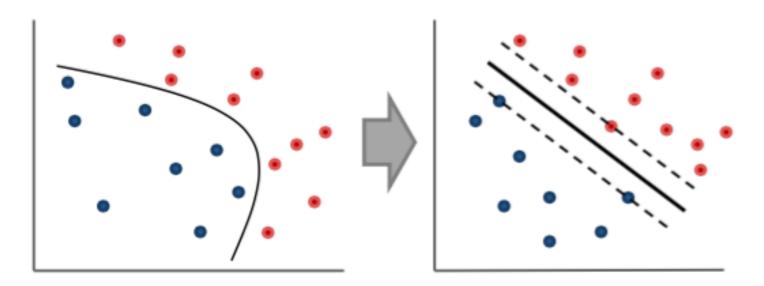


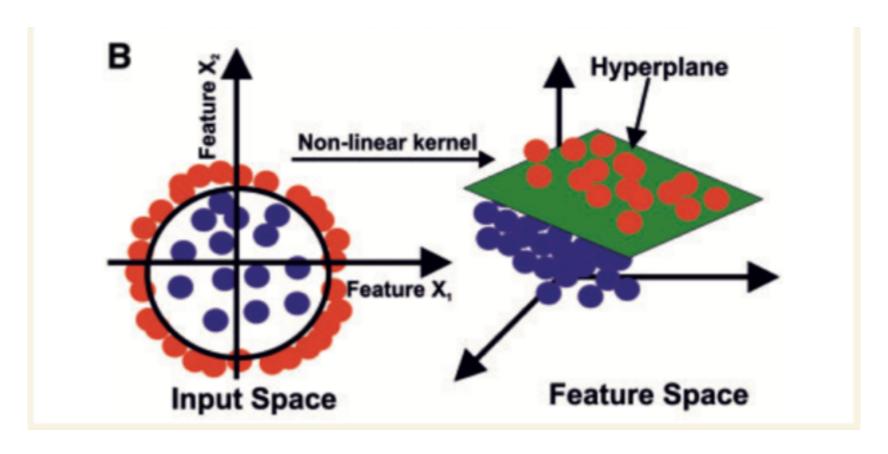


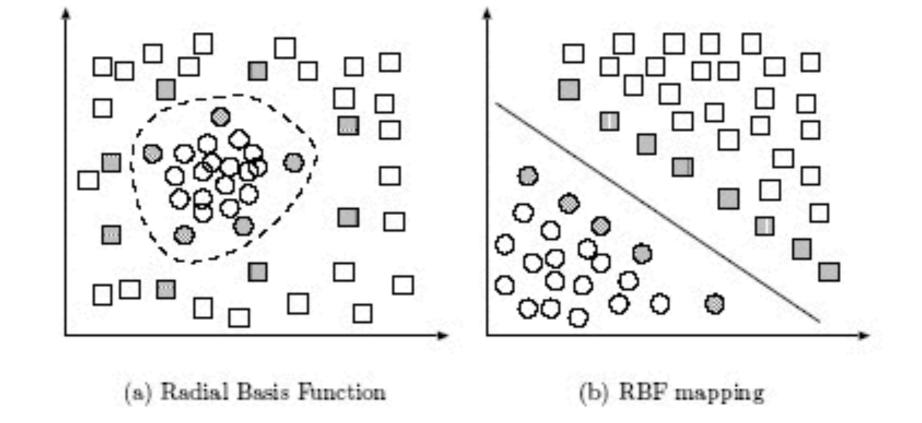


(hyperplane is not just in 2-3D)

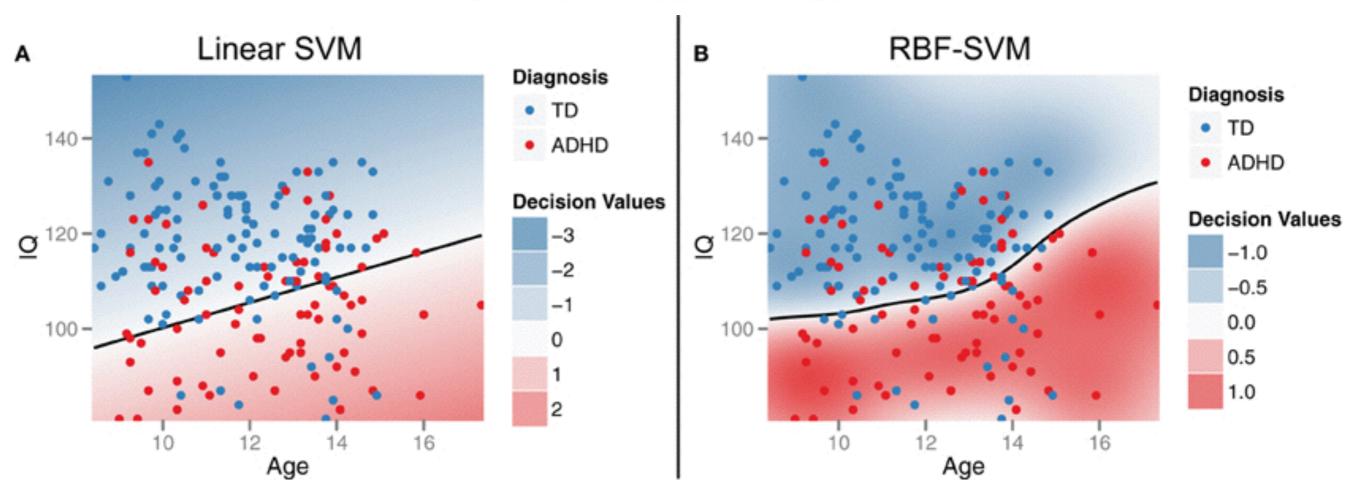
Nonlinear SVM



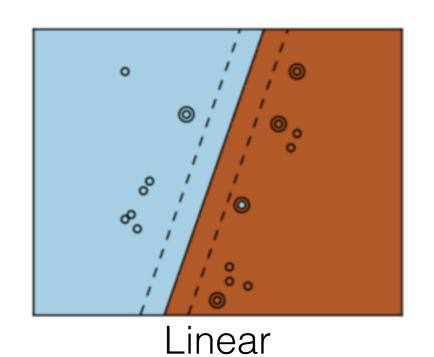


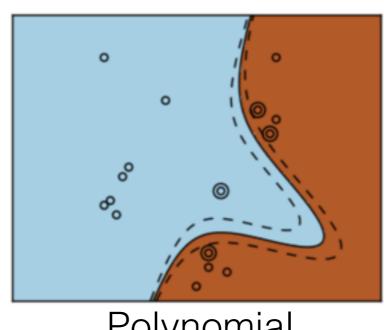


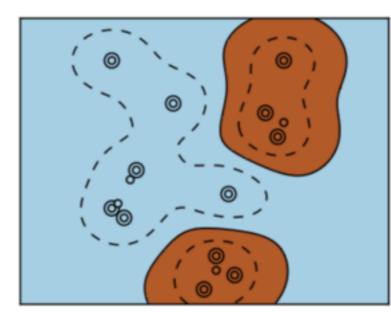
Separable classification with Radial Basis kernel functions in different space. Left: original space. Right: feature space.



SVM - finer details (optional)

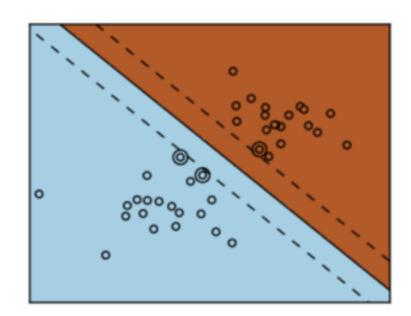


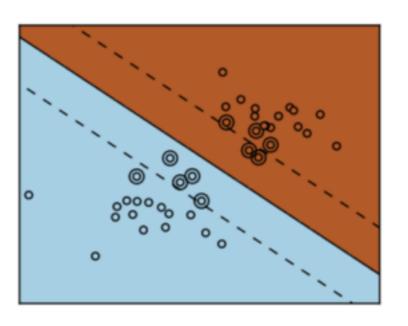




Polynomial

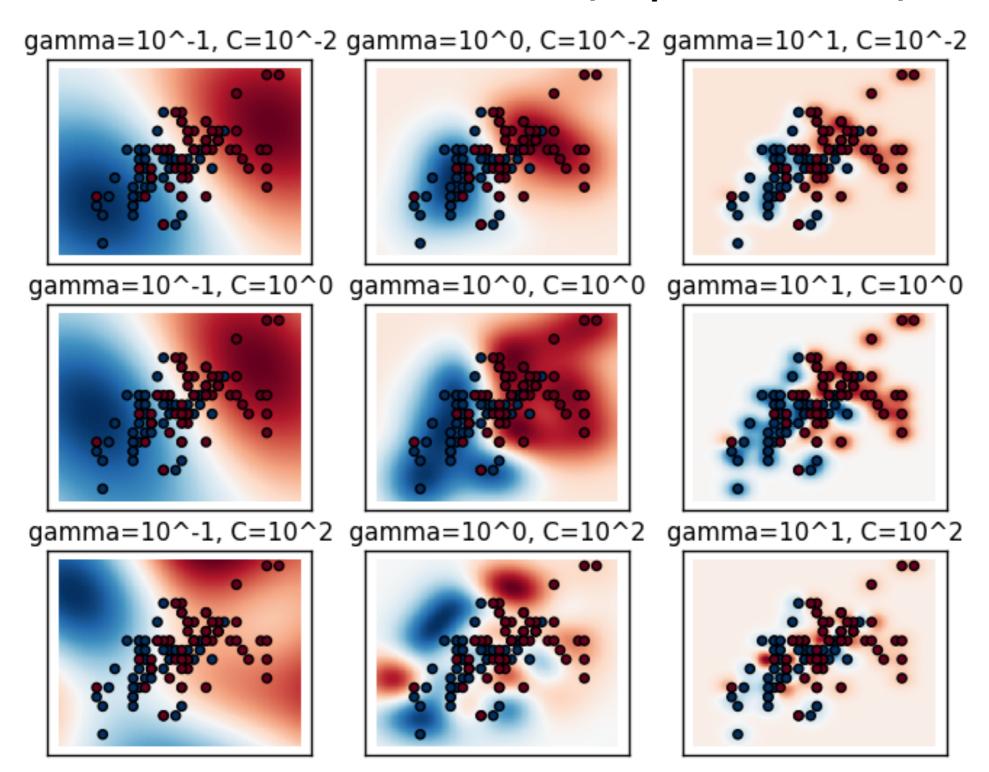
Gaussian/radial basis





Varying "cost", C

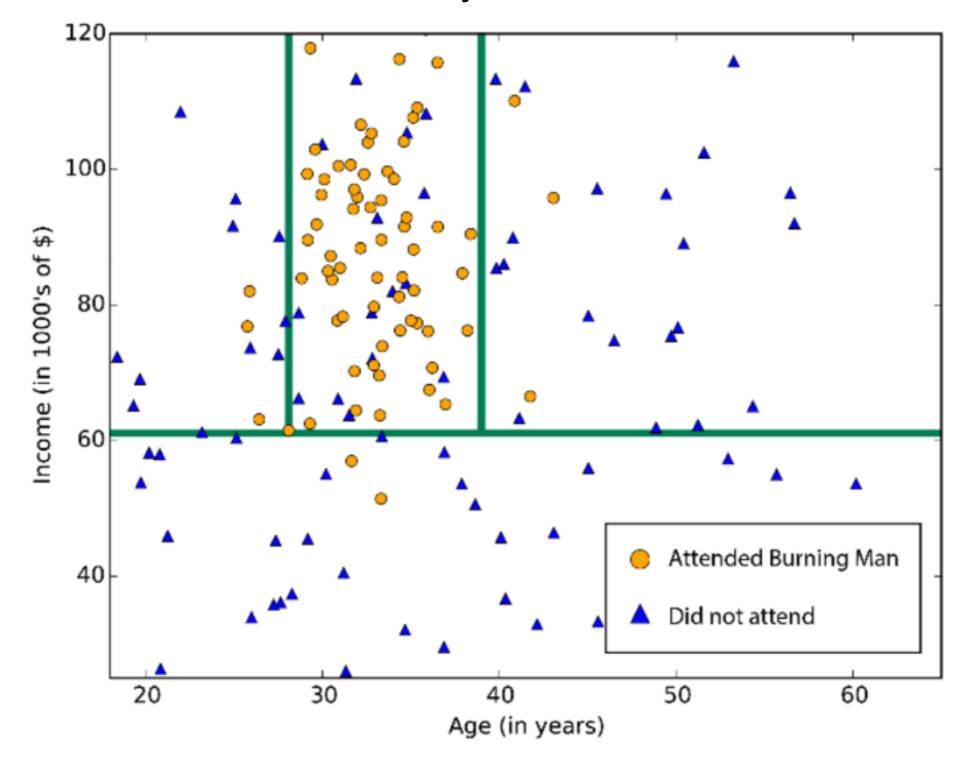
SVM - finer details (optional)



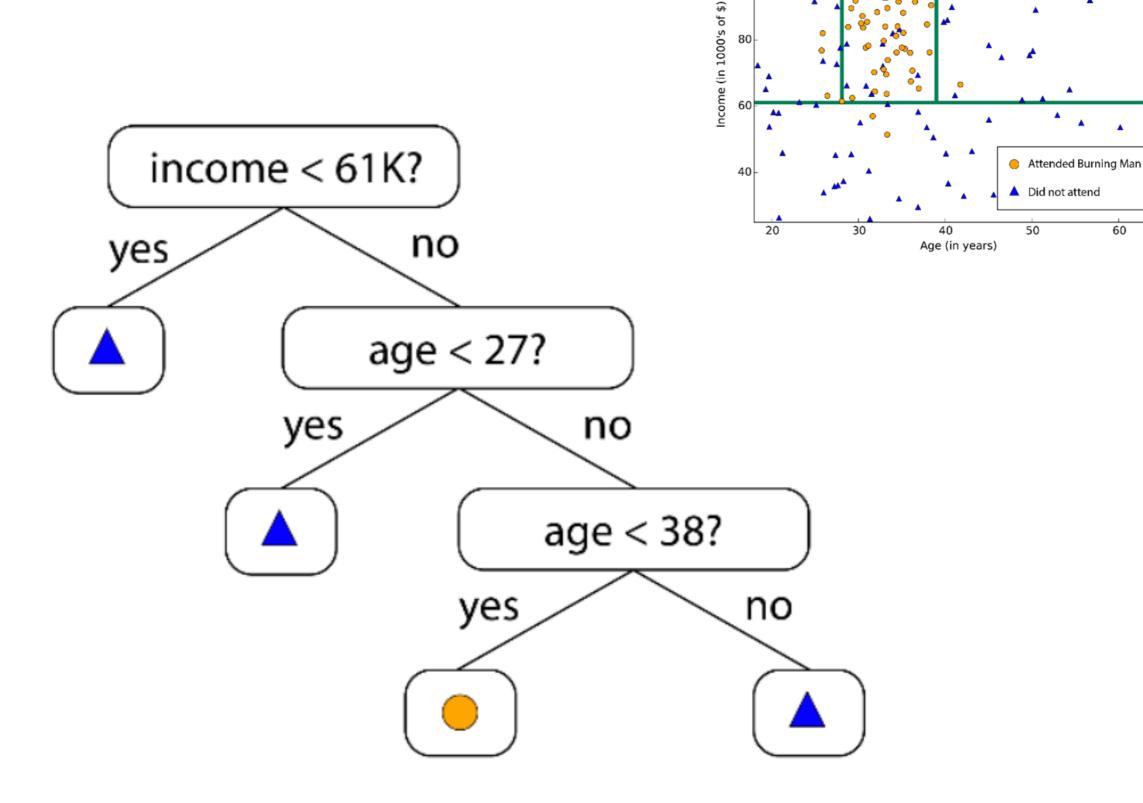
—-> See scikit learn SVM documentation for more information

Decision trees

Disclaimer: I think they are awesome



Decision trees



Decision trees - downsides

Not a one-size-fits-all

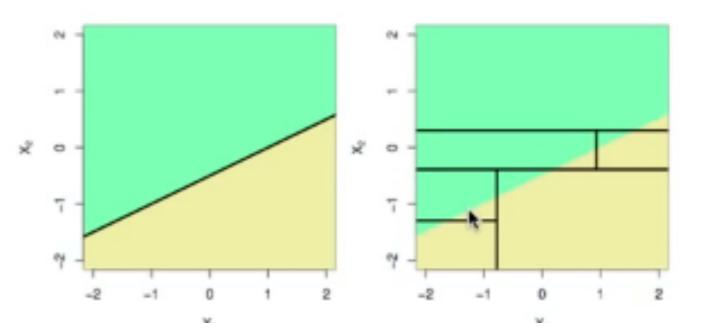
can be clumsy

Finer details

- how do we choose the variable to split on?
- when do we stop splitting?
- what if the tree gets too large?
- define accuracy?

Low bias, but high variance

Accurate, but can overfit



Trees vs linear models

Decision trees - advanced

What if we make loads of really small trees?

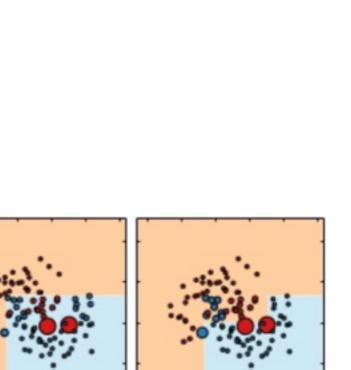
Bagging - "forests" of trees

- Take bootstrap samples (e.g. 1000)
- Fit a small tree (bad accuracy)
- Average predictions of all 1000 trees

Boosting

- As above, but every time we fit a new tree, upweight the misclassified observations
- "each tree is an expert on the errors of its

predecessor"



NB: when these trees only use

a few (randomly chosen)

features, they are called a

"random" forest

Penalized regression

Penalization, or regularization, is the imposing of constraints on a model to limit its complexity

Key concept:

- Just because our model is accurate, doesn't mean it will fit new data
- Maybe a simpler, less accurate model would be better in future?

(Statistically:

- Inclusion of more terms in the model causes high variance in coefficient estimates
- Will re-visit when we talk about bias-variance tradeoff)

Penalized regression

Only include predictors if they are "worth" it

- i.e. they reduce RSS enough to justify inclusion
- two typical methods: LASSO and Ridge regression

L1-regularization:

- aka. LASSO regression
- minimize total absolute beta weight

minimize
$$SSE + \lambda |\beta|$$
 coefficients

Penalized regression

L2-regularization:

- aka. Ridge regression
- minimize total squared beta weight

minimize
$$SSE + \lambda |\beta^2|$$
 coefficients

Elastic net regression: blend of both penalties

Conclusions

- ML algorithms help us identify patterns in the data that predict outcomes we care about
- No one algorithm fits all, so try multiple. But consider:
 - Accuracy, feasibility, and the nature of your problem
- Starting points?
 - k-NN, SVM, Trees (and fancy trees, or penalized methods)
- This was just an intro!
- Try them out— best way to learn good parameters

BUT

- ML is not a replacement for *lots* of good, useful data.
- Fancy algorithms not a substitute for good predictors
 - Don't fall down the rabbit hole.....