

KNN Imputation

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Why KNN imputation?

Mol Breeding (2016)36:69 DOI 10.1007/s11032-016-0490-y



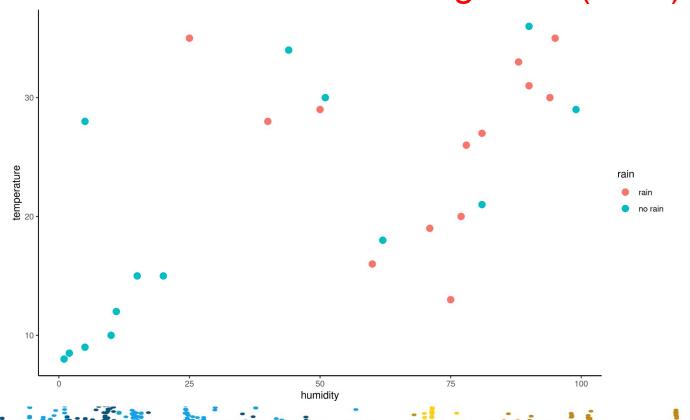
Marker imputation efficiency for genotyping-by-sequencing data in rice (*Oryza sativa*) and alfalfa (*Medicago sativa*)

Nelson Nazzicari · Filippo Biscarini · Paolo Cozzi · E. Charles Brummer · Paolo Annicchiarico





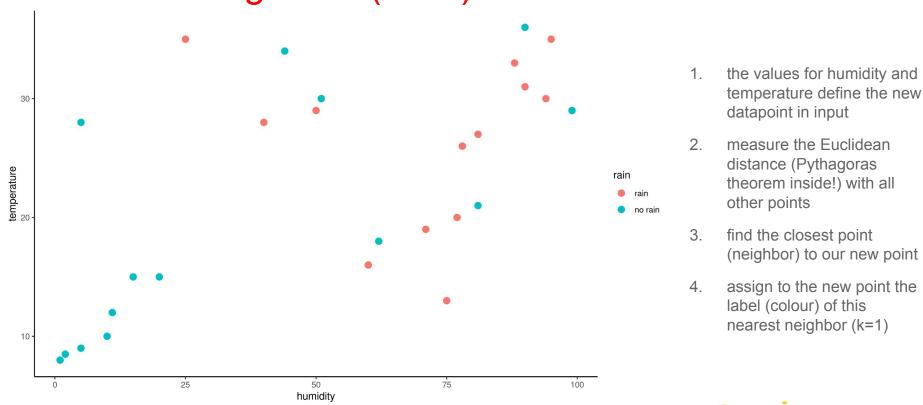
One-minute k-nearest neighbors (KNN)!



- 1. We collect data on temperature, humidity and rain, for a number of days
- 2. New day: will it rain?



k-nearest neighbors (KNN): a bit of math





One-minute k-nearest neighbors (KNN)!

If it looks like a duck, swims like a duck, and quacks like a duck, then it probably is a duck.

$$Pr(Y=j|X=x_0)$$



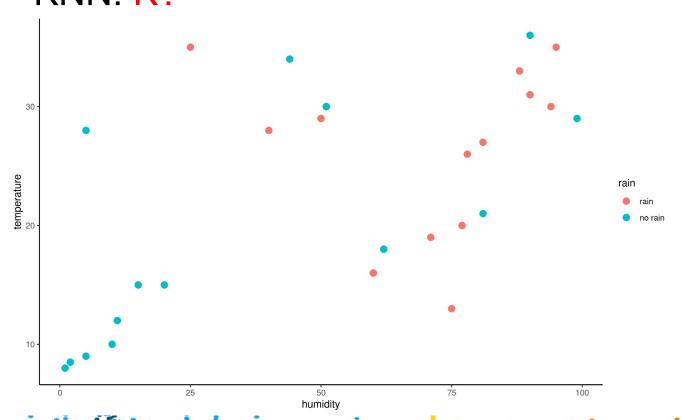


KNN: more than 3-D

- We saw a simple example with few datapoints and only 2 dimensions
- When we have (typically) many dimensions (e.g. many SNPs) and many datapoints (e.g. many samples) we need the machine to do it! (and we can no longer visualize it)



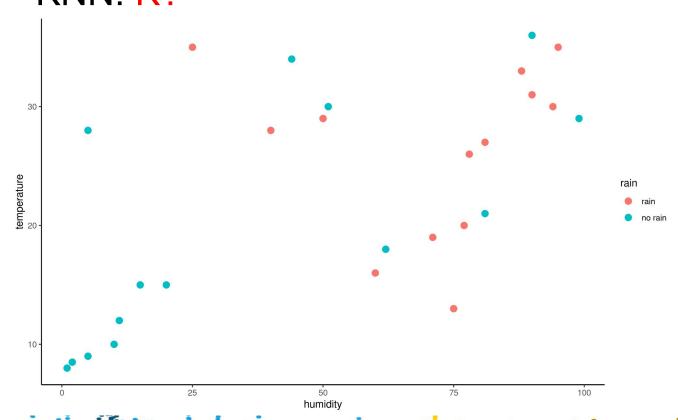




 K-neighborhood → majority!



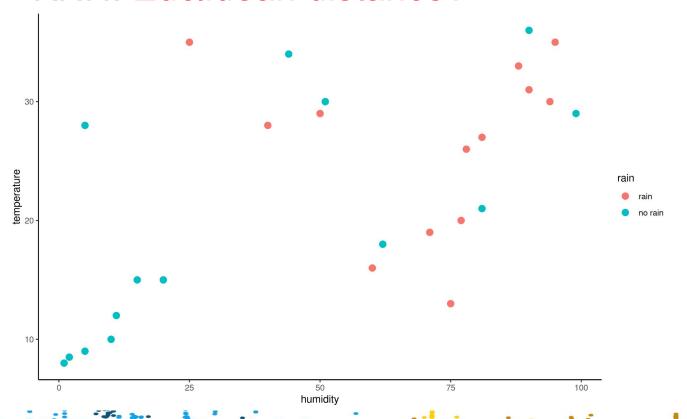
KNN: K?



- K-neighborhood \rightarrow majority!
- majority?



KNN: Euclidean distance?



- similarity↔dissimilarity
- [0, +∞]
- many possible
 distance metrics
 (Hamming,
 Chebyshev, Jaccard etc.) → see here

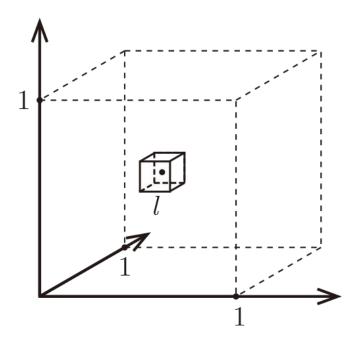


KNN: a few things to tweak

- size of neighborhood (K)
- type of distance
- type of assignment metric (majority, average, weighted metrics, etc.)



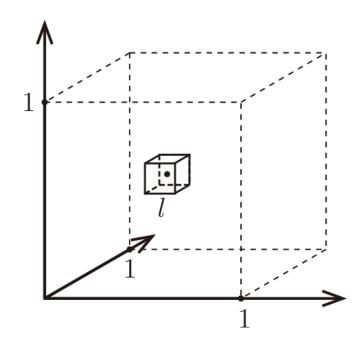
- KNN assumes that similar (close) points share similar labels/target
- unfortunately, in high dimensional spaces points tend to never be close together







 increasing the number of dimensions (parameters) of the problem increases and complicates the identification of k neighbors which are close enough to the data point to be classified/predicted



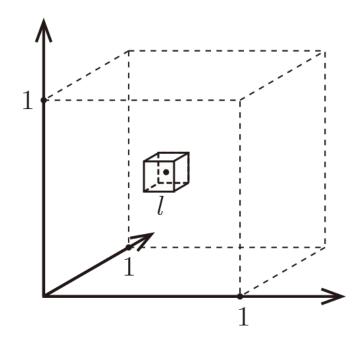




$$l=\left(rac{k}{n}
ight)^{1/d}$$

- I: side of the hypercube that include the k neighbours
- **n**: sample size
- d: n. of dimensions

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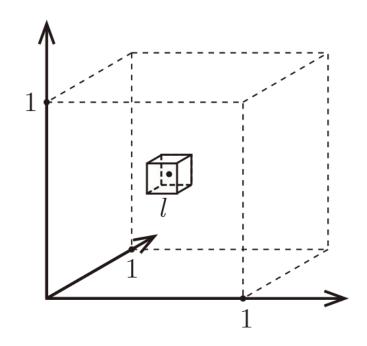




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- I: side of the hypercube that include the k neighbours
- **n**: sample size
- d: n. of dimensions

 with n constant (data size), the hypercube in which the k neighbors lie gets bigger as d increases



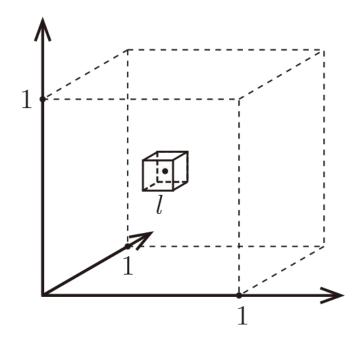




$$l=\left(rac{k}{n}
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- **n** = 1000
- **k** = 5

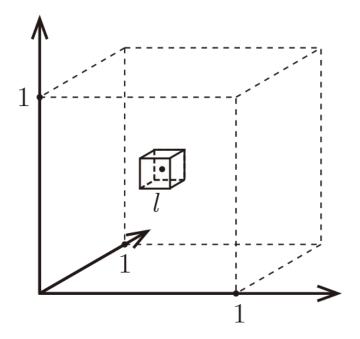
d	1
2	0.071
10	0.588
50	0.899
100	0.948
1000	0.994





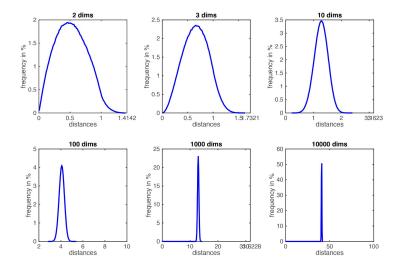
- when $d \gg 1$ almost the entire space is needed to find the k NN
- this breaks down the KNN assumptions, because the k NN are not particularly closer (and therefore more similar) than any other data points in the dataset
- why would the test point share the label with those k-nearest neighbors, if they are not actually similar to it?

d	1
2	0.071
10	0.588
50	0.899
100	0.948
1000	0.994





distributions of all pairwise distances between randomly drawn points within d-dimensional unit hypercubes: as the number of dimensions *d* grows, all distances concentrate within a very small range ("the night where all cows are black")





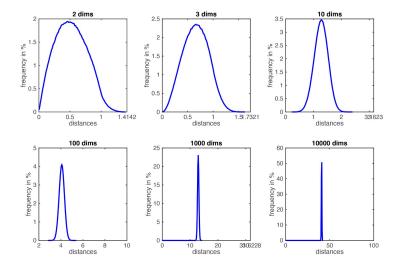


distributions of all pairwise distances between randomly drawn points within d-dimensional unit hypercubes: as the number of dimensions *d* grows, all distances concentrate within a very small range ("the night where all cows are black")

increase the number of samples *n* (data size)?

$$n=rac{k}{I^d}=k\cdot 10^d$$

- exponential growth
- for d > 100 we would need more data points than there are electrons in the universe!



From: https://www.cs.cornell.edu/courses/cs4780/2018fa/lectures/lecturenote02_kNN.html



KNN: as lazy as it gets

- KNN is a **lazy algorithm**: each time new datapoints are added (e.g. to be predicted) pairwise distances with all existing datapoints (over all dimensions) must be calculated
- calculations are slow
- However:
 - when new data are available, there's no need to retrain the model (no parameters to estimate or fine-tune) → excellent for applications where data are added incrementally (e.g. on-line learning, update predictions)





What about imputation?

- Ok, we learnt about KNN, but imputation?
- The imputation of missing SNP genotypes is a type of prediction (slightly "sui generis"), where non-missing values can be considered as training observations and missing values as the test observations



NEXT LECTURE

Genotype imputation with KNN: a demonstration (R code)