

EVALUATION

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CS 158 – Fall 2023

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Admin

Assignment 3

- Autograder

Reading

Mentor hours this week: Thursday and Sunday, 7-9pm

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So far...

- Throw out outlier examples
- Remove noisy features
- Pick "good" features
- Normalize feature values
 - center data
 - scale data (either variance or absolute)
- Normalize example length
- Finally, train your model!

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What about testing?

training data (labeled examples)

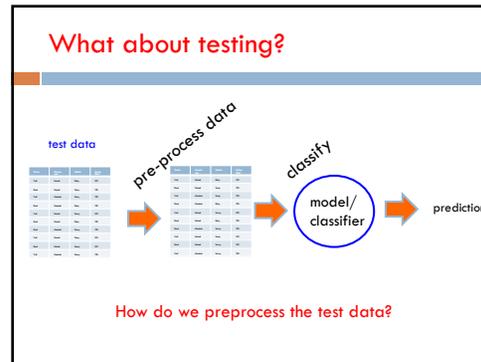
Pre-process data

better training data

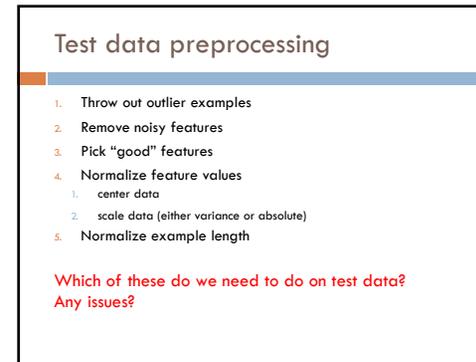
learn

model/classifier

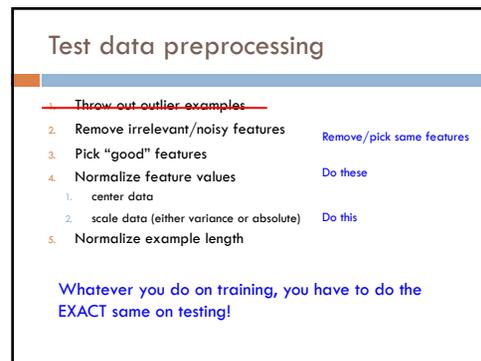
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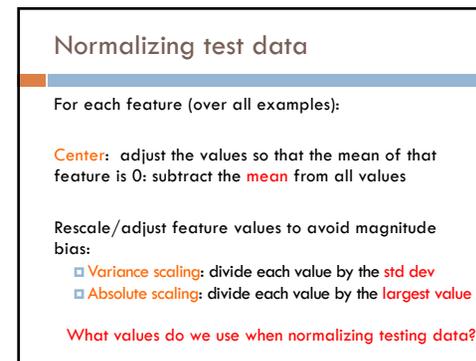
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Normalizing test data

For each feature (over all examples):

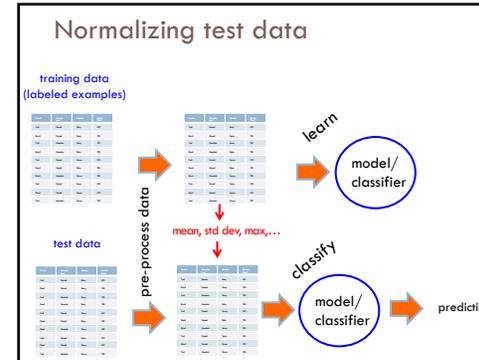
Center: adjust the values so that the mean of that feature is 0: subtract the **mean** from all values

Rescale/adjust feature values to avoid magnitude bias:

- ▢ **Variance scaling:** divide each value by the **std dev**
- ▢ **Absolute scaling:** divide each value by the **largest value**

Save these from training normalization!

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Features pre-processing summary

Many techniques for preprocessing data

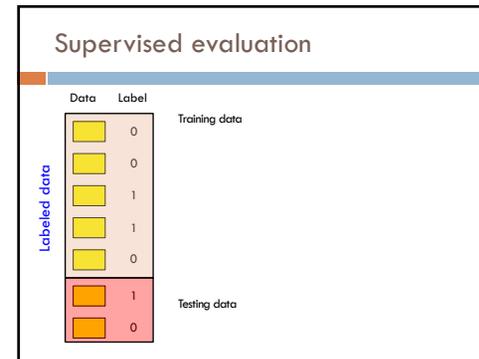
Which ones will work well will depend on the data and the classifier

Try them out and evaluate how they affect performance on dev data

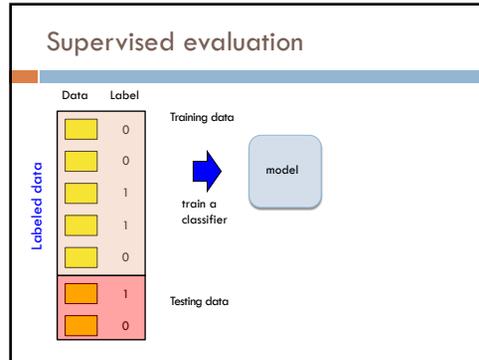
Make sure to do the **exact same** pre-processing on train and test

1. Throw out outlier examples
2. Remove noisy features
3. Pick "good" features
4. Normalize feature values
 1. center data
 2. scale data (either variance or absolute)
5. Normalize example length

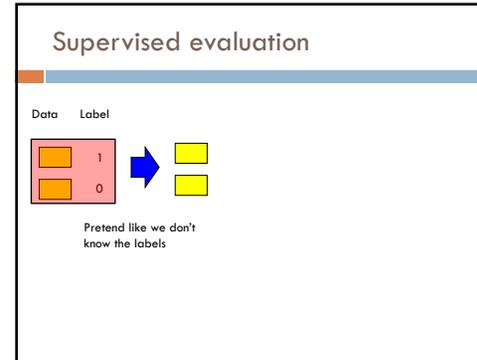
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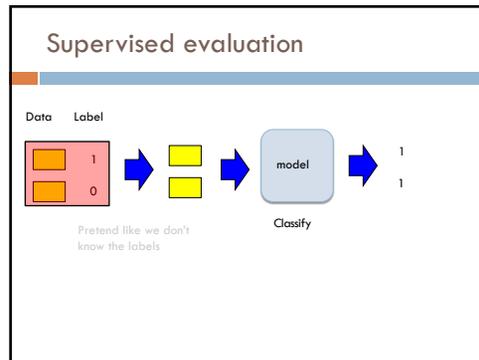
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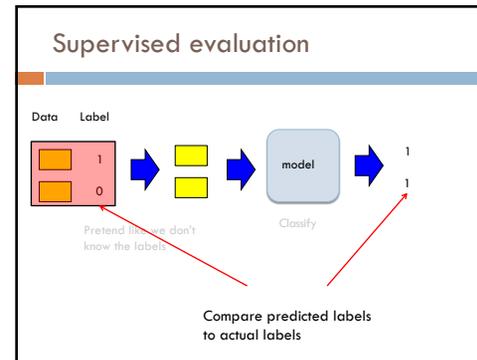
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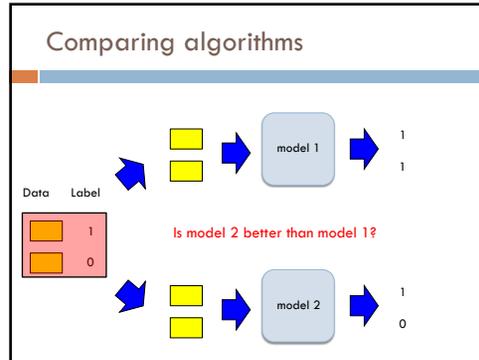
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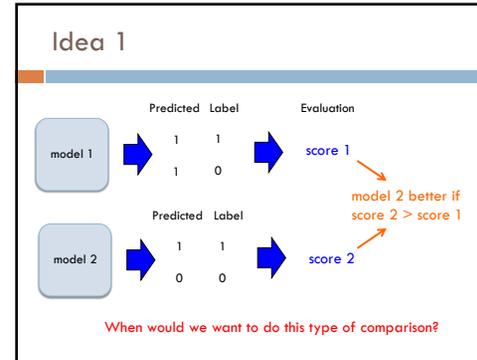
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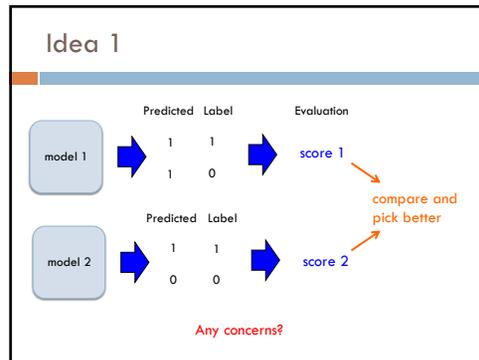
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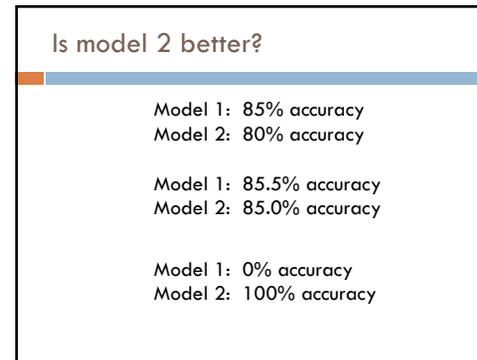
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Comparing scores: significance

Just comparing scores on one data set isn't enough!

We don't just want to know which system is better on *this particular data*, we want to know if model 1 is better than model 2 *in general*

Put another way, we want to be confident that the difference is real and not just due to chance or eccentricities about the particular dataset

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Idea 2

	Predicted	Label	Evaluation
model 1	1	1	score 1
	1	0	
model 2	1	1	score 2
	0	0	

model 2 better if $score\ 2 + c > score\ 1$

Is this any better?

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Idea 2

	Predicted	Label	Evaluation
model 1	1	1	score 1
	1	0	
model 2	1	1	score 2
	0	0	

model 2 better if $score\ 2 + c > score\ 1$

NO!
Key: we don't know the variance of the output

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Variance

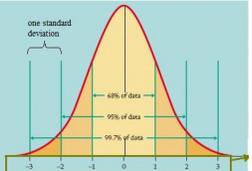
Recall that variance (or standard deviation) helped us predict how likely certain events are:

How do we know how variable a model's accuracy is?

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Variance

Recall that variance (or standard deviation) helped us predict how likely certain events are:



We need multiple accuracy scores! Ideas?

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Repeated experimentation

	Data	Label	
Labeled data	Yellow	0	Training data
	Yellow	0	
	Yellow	1	
	Yellow	1	
	Yellow	0	
	Orange	1	Testing data
	Orange	0	

Rather than just splitting once, split multiple times

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Repeated experimentation

	Data	Label	Data	Label	Data	Label	
Training data	Yellow	0	Yellow	0	Orange	0	...
	Yellow	0	Orange	0	Yellow	0	
	Yellow	1	Yellow	1	Yellow	1	
	Yellow	1	Orange	1	Orange	1	
	Orange	0	Yellow	0	Yellow	0	
	Orange	1	Yellow	1	Yellow	1	
			Yellow	0	Yellow	0	
			Yellow	1	Yellow	1	

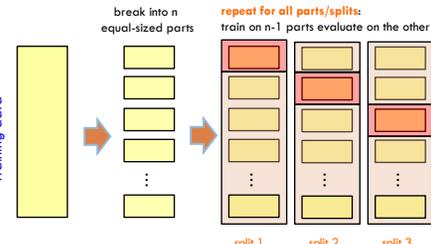
= train
 = development

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n-fold cross validation

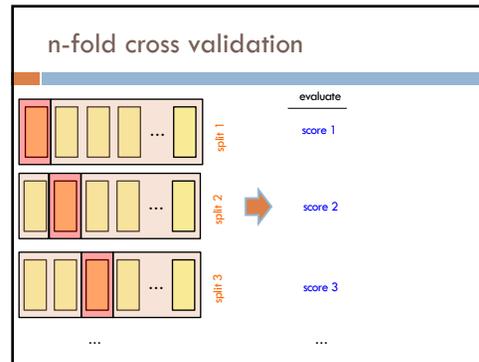
break into n equal-sized parts

repeat for all parts/splits
train on n-1 parts evaluate on the other



split 1 split 2 split 3

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n-fold cross validation

- better utilization of labeled data
- more robust: doesn't just rely on one test/development set to evaluate the approach (or for optimizing parameters)
- multiplies the computational overhead by n (have to train n models instead of just one)
- 10 is the most common choice of n

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Leave-one-out cross validation

n -fold cross validation where n = number of examples

aka "jackknifing"

pros/cons?

when would we use this?

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Leave-one-out cross validation

- Can be very expensive if training is slow and/or if there are a large number of examples
- Useful in domains with limited training data: maximizes the data we can use for training
- Some classifiers are very amenable to this approach (e.g.?)

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Random splits

n -fold cross-validation explicitly divides the data into n partitions and each data point gets used exactly once for testing

Another approach is do some number of random X/Y splits (like we did on Assignment 2)

Pros/cons?

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Random splits

Lose the guarantee that all examples are used once for testing

Can be slower (if we do more splits)

Can allow for more samples

Both are fine approaches!

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Comparing systems: sample 1

split	model 1	model 2
1	87	88
2	85	84
3	83	84
4	80	79
5	88	89
6	85	85
7	83	81
8	87	86
9	88	89
10	84	85
average:	85	85

Is model 2 better than model 1?

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Comparing systems: sample 2

split	model 1	model 2
1	87	87
2	92	88
3	74	79
4	75	86
5	82	84
6	79	87
7	83	81
8	83	92
9	88	81
10	77	85
average:	82	85

Is model 2 better than model 1?

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Comparing systems: sample 3

split	model 1	model 2
1	84	87
2	83	86
3	78	82
4	80	86
5	82	84
6	79	87
7	83	84
8	83	86
9	85	83
10	83	85
average:	82	85

Is model 2 better than model 1?

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Comparing systems

split	model 1	model 2
1	84	87
2	83	86
3	78	82
4	80	86
5	82	84
6	79	87
7	83	84
8	83	86
9	85	83
10	83	85
average:	82	85

split	model 1	model 2
1	87	87
2	92	88
3	74	79
4	75	86
5	82	84
6	79	87
7	83	81
8	83	92
9	88	81
10	77	85
average:	82	85

What's the difference?

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Comparing systems

split	model 1	model 2
1	84	87
2	83	86
3	78	82
4	80	86
5	82	84
6	79	87
7	83	84
8	83	86
9	85	83
10	83	85
average:	82	85
std dev	2.3	1.7

split	model 1	model 2
1	87	87
2	92	88
3	74	79
4	75	86
5	82	84
6	79	87
7	83	81
8	83	92
9	88	81
10	77	85
average:	82	85
std dev	5.9	3.9

Even though the averages are same, the variance is different!

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Comparing systems: sample 4

split	model 1	model 2
1	80	82
2	84	87
3	89	90
4	78	82
5	90	91
6	81	83
7	80	80
8	88	89
9	76	77
10	86	88
average:	83	85
std dev	4.9	4.7

Is model 2 better than model 1?

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Comparing systems: sample 4

split	model 1	model 2	model 2 - model 1
1	80	82	2
2	84	87	3
3	89	90	1
4	78	82	4
5	90	91	1
6	81	83	2
7	80	80	0
8	88	89	1
9	76	77	1
10	86	88	2
average:	83	85	
std dev	4.9	4.7	

Is model 2 better than model 1?

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Comparing systems: sample 4

split	model 1	model 2	model 2 - model 1
1	80	82	2
2	84	87	3
3	89	90	1
4	78	82	4
5	90	91	1
6	81	83	2
7	80	80	0
8	88	89	1
9	76	77	1
10	86	88	2
average:	83	85	
std dev	4.9	4.7	

Model 2 is ALWAYS better

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Comparing systems: sample 4

split	model 1	model 2	model 2 - model 1
1	80	82	2
2	84	87	3
3	89	90	1
4	78	82	4
5	90	91	1
6	81	83	2
7	80	80	0
8	88	89	1
9	76	77	1
10	86	88	2
average:	83	85	
std dev	4.9	4.7	

How do we decide if model 2 is better than model 1?

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Statistical tests

Setup:

- Assume some default hypothesis about the data that you'd like to *disprove*, called the **null hypothesis**
- e.g. model 1 and model 2 are not statistically different in performance

Test:

- Calculate a test statistic from the data (often assuming something about the data)
- Based on this statistic, with *some probability* we can **reject the null hypothesis**, that is, show that it does not hold

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t-test

Determines whether two samples come from the same underlying distribution or not

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t-test

Null hypothesis: model 1 and model 2 accuracies are no different, i.e. come from **the same** distribution

Assumptions: there are a number that often aren't completely true, but we're often not too far off

Result: probability that the difference in accuracies is due to random chance (low values are better)

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Calculating t-test

For our setup, we'll do what's called a "paired t-test"

- The values can be thought of as pairs, where they were calculated under the same conditions
- In our case, the same train/test split
- Gives more power than the unpaired t-test (we have more information)

For almost all experiments, we'll do a "two-tailed" version of the t-test

Can calculate by hand or in code, but why reinvent the wheel: use excel or a statistical package

http://en.wikipedia.org/wiki/Student's_t-test

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p-value

The result of a statistical test is often a p-value

p-value: the probability that the null hypothesis holds. Specifically, if we re-ran this experiment multiple times (say on different data) what is the probability that we would reject the null hypothesis incorrectly (i.e. the probability we'd be wrong)

Common values to consider "significant": 0.05 (95% confident), 0.01 (99% confident) and 0.001 (99.9% confident)

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Comparing systems: sample 1

split	model 1	model 2
1	87	88
2	85	84
3	83	84
4	80	79
5	88	89
6	85	85
7	83	81
8	87	86
9	88	89
10	84	85
average:	85	85

Is model 2 better than model 1?

They are the same with:
 $p = 1$

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Comparing systems: sample 2

split	model 1	model 2
1	87	87
2	92	88
3	74	79
4	75	86
5	82	84
6	79	87
7	83	81
8	83	92
9	88	81
10	77	85
average:	82	85

Is model 2 better than model 1?

They are the same with:
 $p = 0.15$

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Comparing systems: sample 3

split	model 1	model 2
1	84	87
2	83	86
3	78	82
4	80	86
5	82	84
6	79	87
7	83	84
8	83	86
9	85	83
10	83	85
average:	82	85

Is model 2 better than model 1?

They are the same with:
 $p = 0.007$

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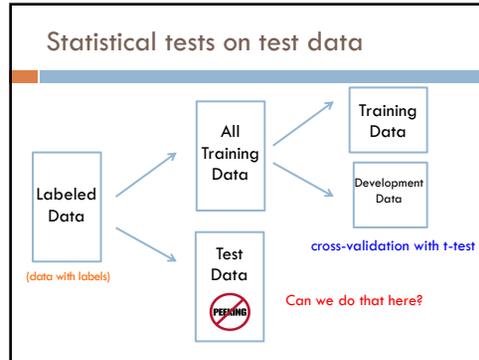
Comparing systems: sample 4

split	model 1	model 2
1	80	82
2	84	87
3	89	90
4	78	82
5	90	91
6	81	83
7	80	80
8	88	89
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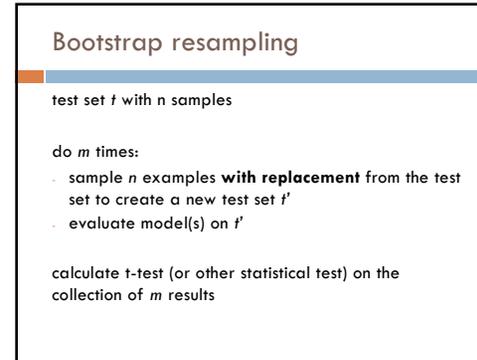
Is model 2 better than model 1?

They are the same with:
 $p = 0.001$

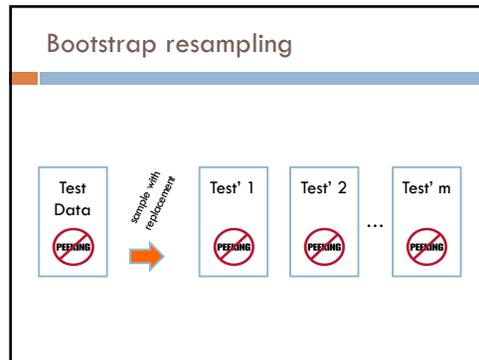
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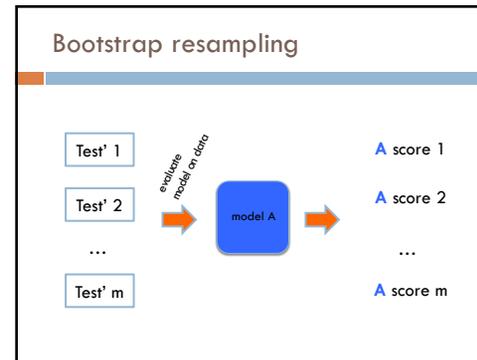
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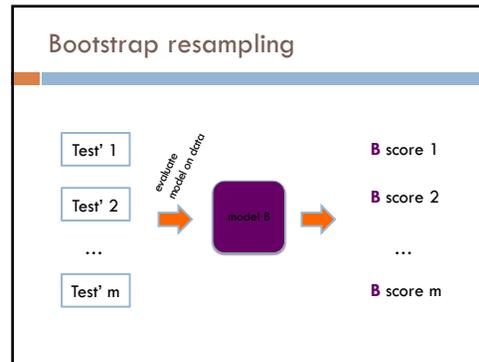
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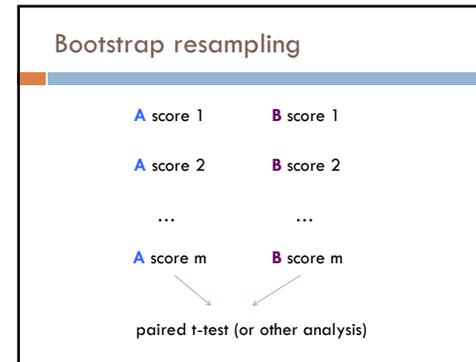
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Experimentation good practices

Never look at your test data!

During development

- ▣ Compare different models/hyperparameters on development data
- ▣ use cross-validation to get more consistent results
- ▣ If you want to be confident with results, use a t-test and look for $p = 0.05$ (or even lower)

For final evaluation, use bootstrap resampling combined with a t-test to compare final approaches

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