

pomegranate

fast and flexible probabilistic modelling in python

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Acknowledgements



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ADVANCING DATA-INTENSIVE DISCOVERY IN ALL FIELDS

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Overview

pomegranate is **more flexible** than other packages, **faster**, is **intuitive to use**, and can do it all **in parallel**



Overview: supported models

Six Main Models:

1. Probability Distributions
2. General Mixture Models
3. Markov Chains
4. Hidden Markov Models
5. Bayes Classifiers / Naive Bayes
6. Bayesian Networks

Two Helper Models:

1. k-means++/kmeans||
2. Factor Graphs



Overview: model stacking in pomegranate

Distributions

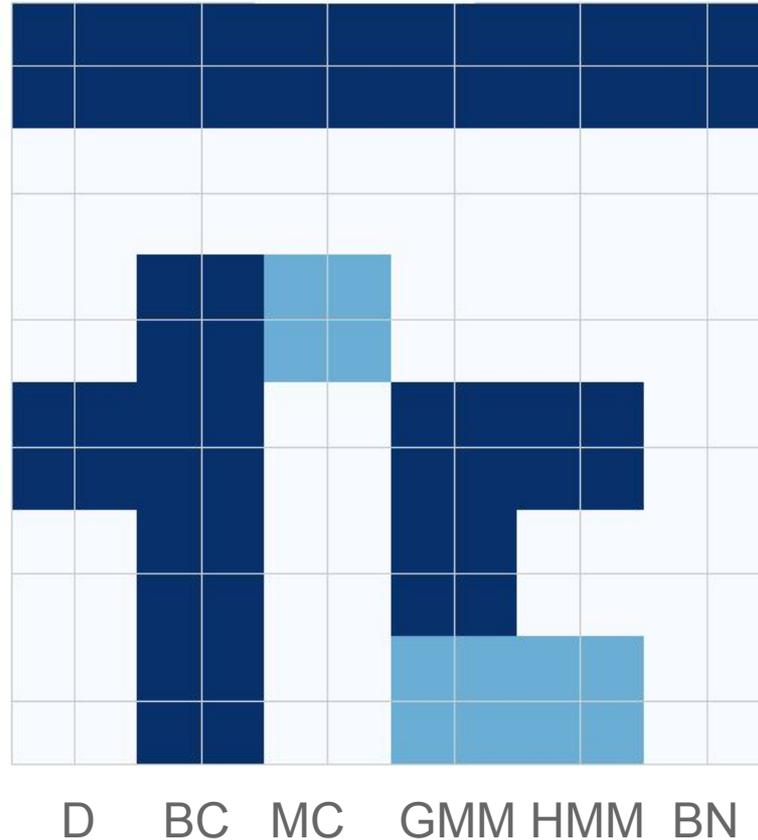
Bayes Classifiers

Markov Chains

General Mixture Models

Hidden Markov Models

Bayesian Networks





The API is common to all models

`model.log_probability(X) / model.probability(X)`

`model.sample()`

`model.fit(X, weights, inertia)`

All models have these methods!

`model.summarize(X, weights)`

`model.from_summaries(inertia)`

`Model.from_samples(X, weights)`

`model.predict(X)`

`model.predict_proba(X)`

`model.predict_log_proba(X)`

All models composed of distributions (like GMM, HMM...) have these methods too!



pomegranate supports many distributions

Univariate Distributions

1. UniformDistribution
2. BernoulliDistribution
3. NormalDistribution
4. LogNormalDistribution
5. ExponentialDistribution
6. BetaDistribution
7. GammaDistribution
8. DiscreteDistribution
9. PoissonDistribution

Kernel Densities

1. GaussianKernelDensity
2. UniformKernelDensity
3. TriangleKernelDensity

Multivariate Distributions

1. IndependentComponentsDistribution
2. MultivariateGaussianDistribution
3. DirichletDistribution
4. ConditionalProbabilityTable
5. JointProbabilityTable



pomegranate can be faster than numpy

Fitting Multivariate Gaussian to 10,000,000 samples of 10 dimensions

```
data = numpy.random.randn(10000000, 10)

print "numpy time:"
%timeit -n 10 data.mean(), numpy.cov(data.T)
print
print "pomegranate time:"
%timeit -n 10 MultivariateGaussianDistribution.from_samples(data)
```

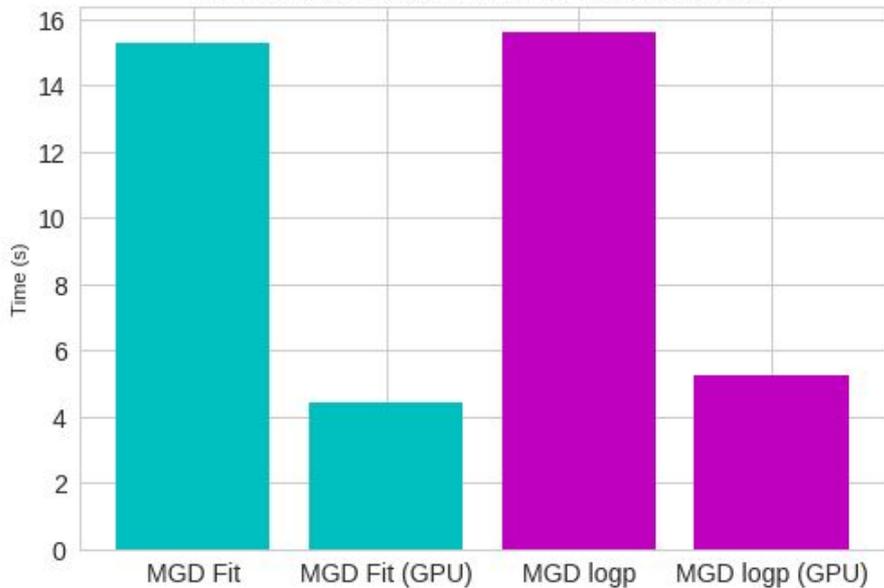
numpy time:
10 loops, best of 3: 1.02 s per loop

pomegranate time:
10 loops, best of 3: 799 ms per loop

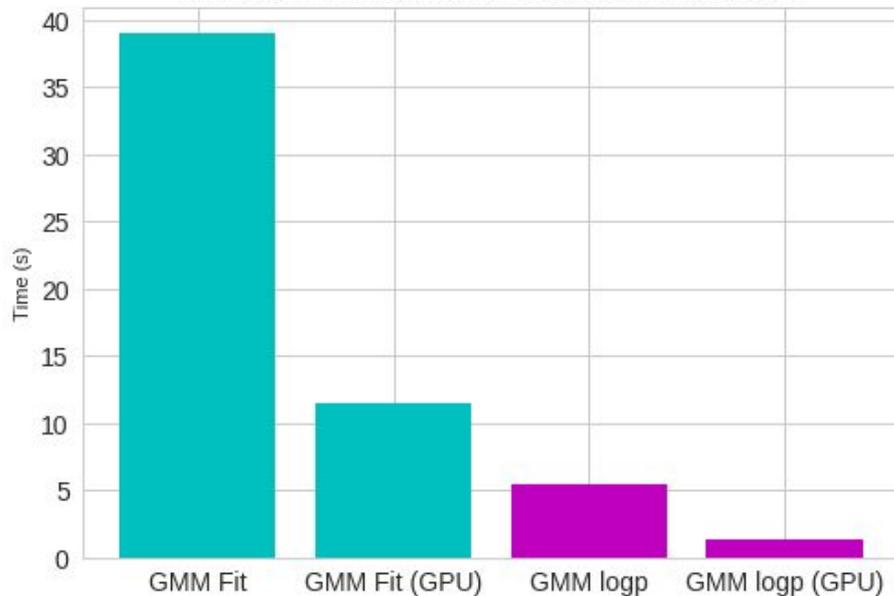


pomegranate just merged GPU support

Multivariate Gaussian with GPU Acceleration



Gaussian Mixture Model with GPU Acceleration





pomegranate uses additive summarization

pomegranate reduces data to sufficient statistics for updates and so only has to go datasets once (for all models).

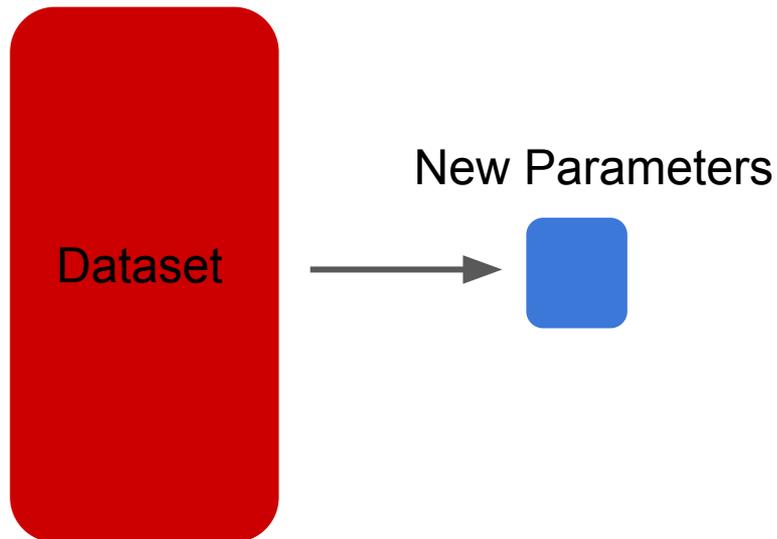
Here is an example of the Normal Distribution sufficient statistics

$$\sum_{i=1}^n w_i \quad \sum_{i=1}^n w_i x_i \quad \sum_{i=1}^n w_i x_i^2 \quad \longrightarrow \quad \begin{aligned} \mu &= \frac{\sum_{i=1}^n w_i x_i}{\sum_{i=1}^n w_i} \\ \sigma^2 &= \frac{\sum_{i=1}^n w_i x_i^2}{\sum_{i=1}^n w_i} - \frac{\left(\sum_{i=1}^n w_i x_i\right)^2}{\left(\sum_{i=1}^n w_i\right)^2} \end{aligned}$$



pomegranate supports out-of-core learning

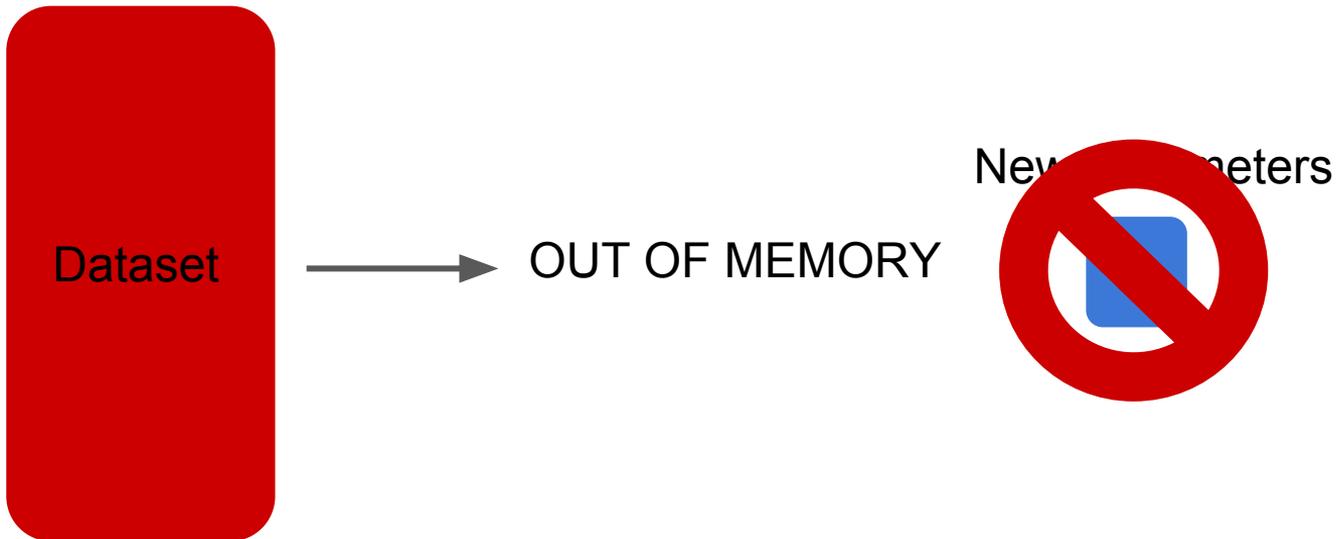
Typically, one wants to get new, better, parameters from data





pomegranate supports out-of-core learning

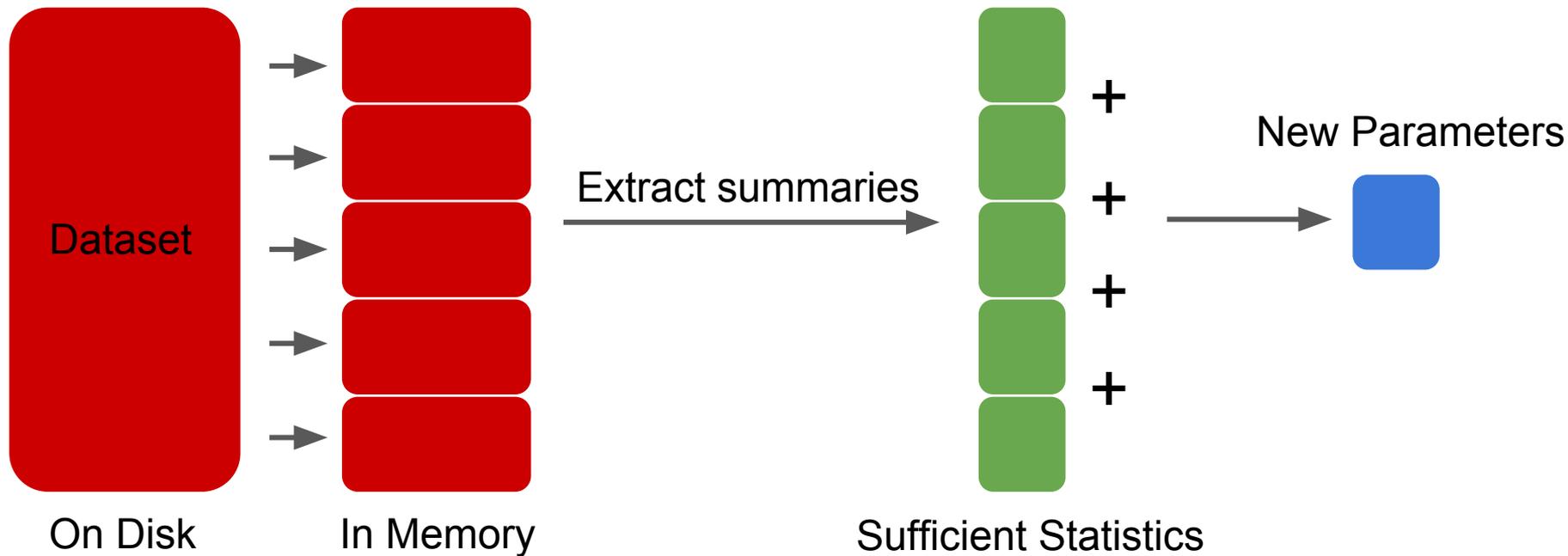
If the dataset is too big, sometimes what you get instead is an out of memory error.





pomegranate supports out-of-core learning

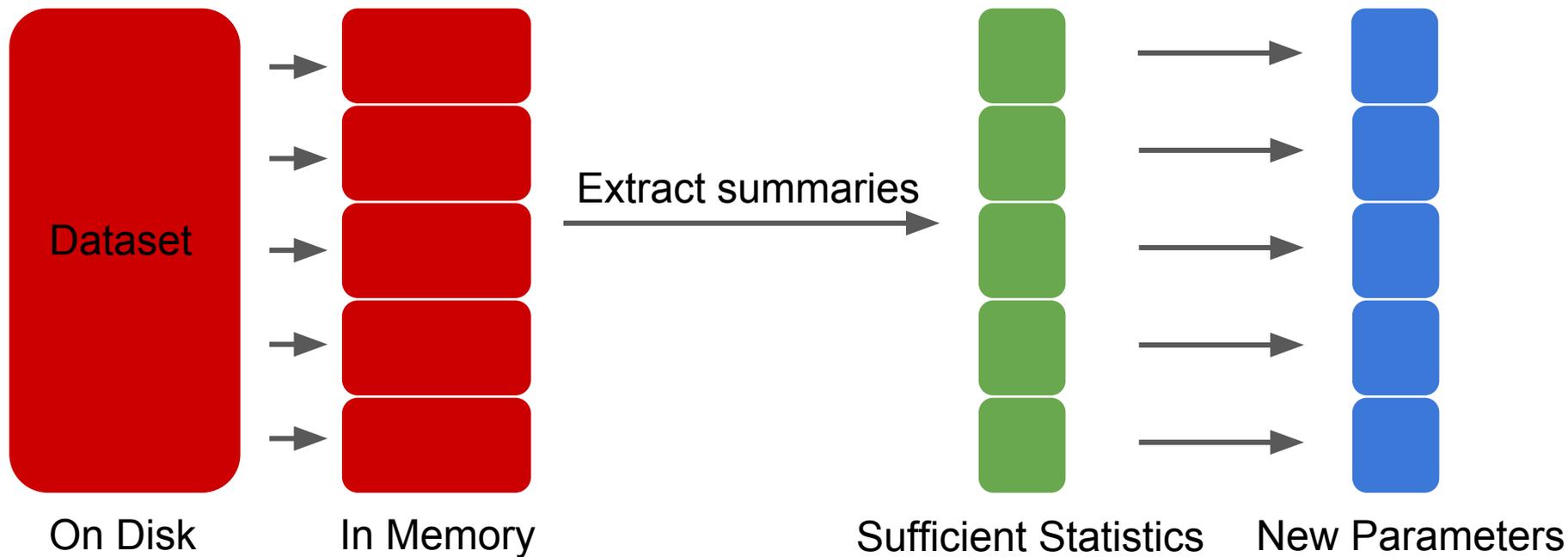
Batches from a dataset can be reduced to additive summary statistics, enabling exact updates from data that can't fit in memory.





pomegranate supports mini-batching

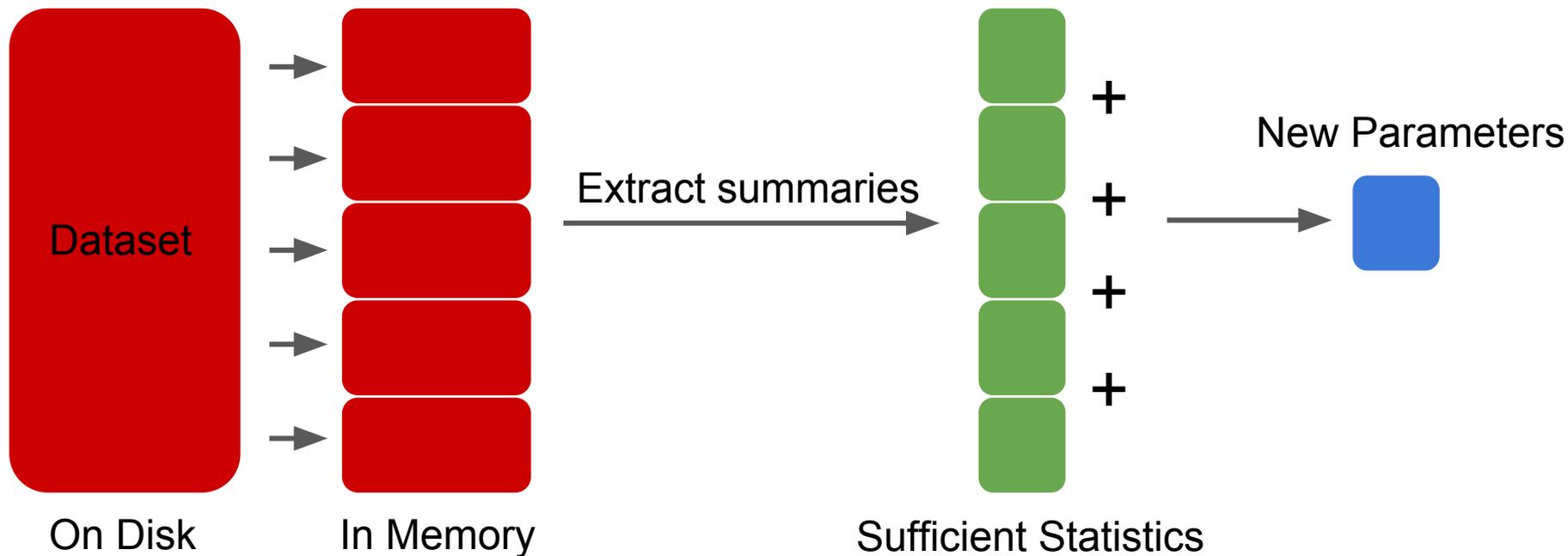
Instead of going through the full dataset before updating parameters, one could update parameters at each step.





pomegranate supports parallelization

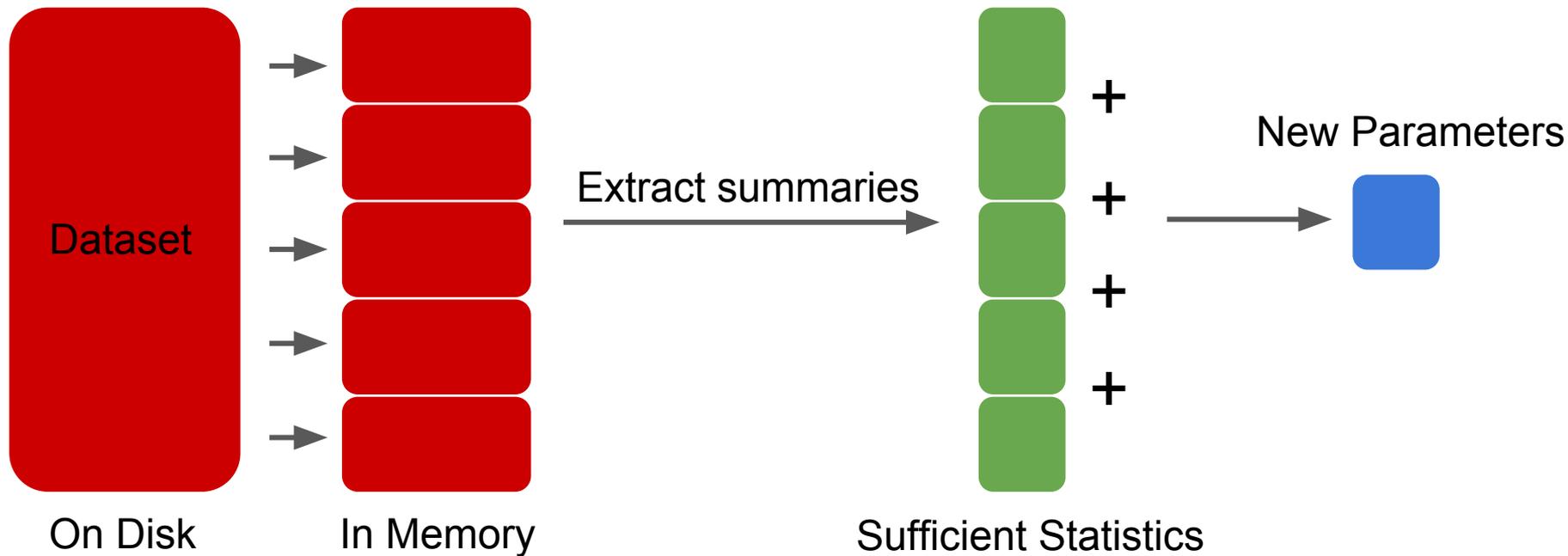
Batches from a dataset can be reduced to additive summary statistics, enabling exact updates from data that can't fit in memory.





pomegranate supports parallelization

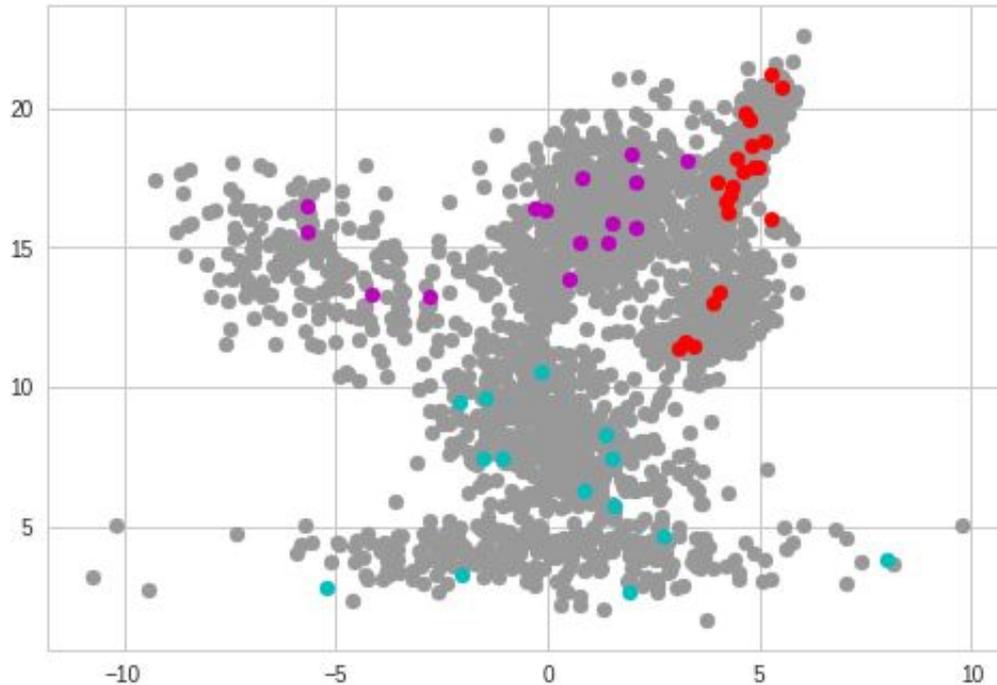
Batches from a dataset can be reduced to additive summary statistics, enabling exact updates from data that can't fit in memory.





pomegranate supports semisupervised learning

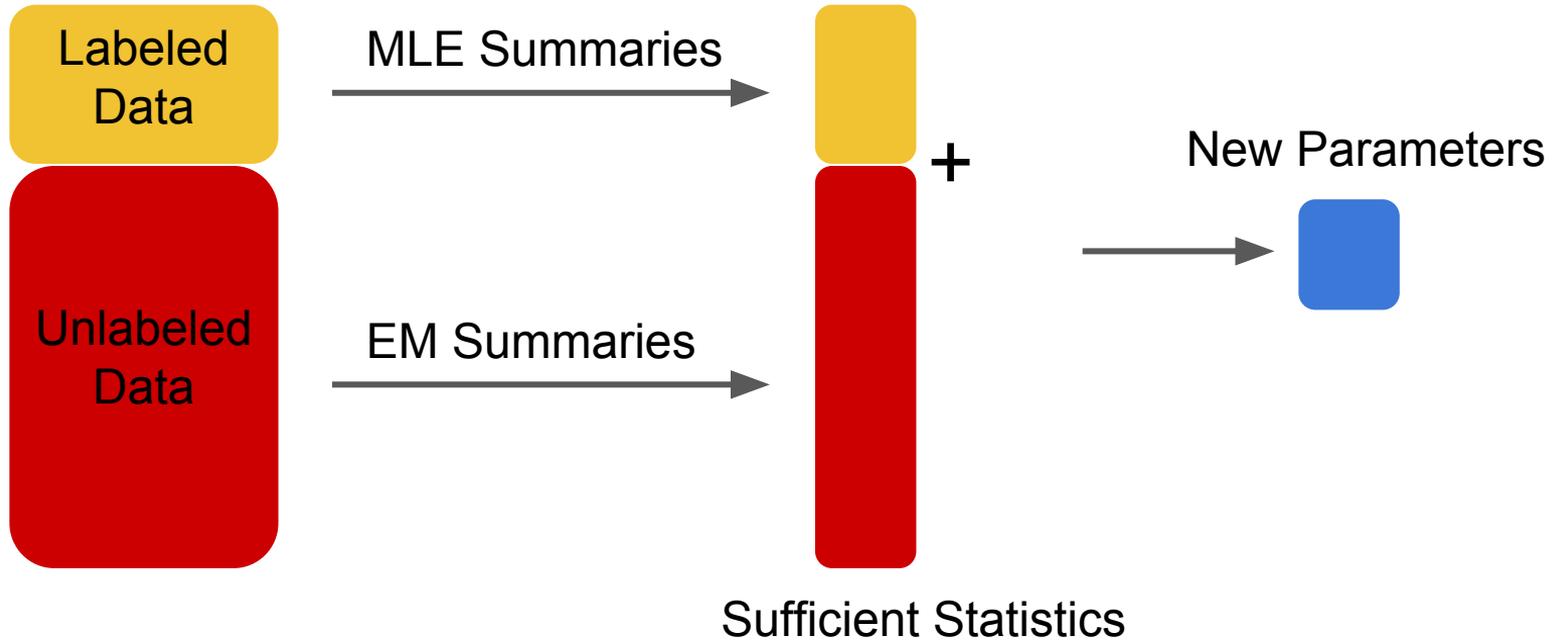
For many tasks, there is limited labeled data but a deluge of unlabeled data, and one wants to utilize both.





pomegranate supports semisupervised learning

Summaries from MLE on the labeled data can be added to summaries from EM on the unlabeled data

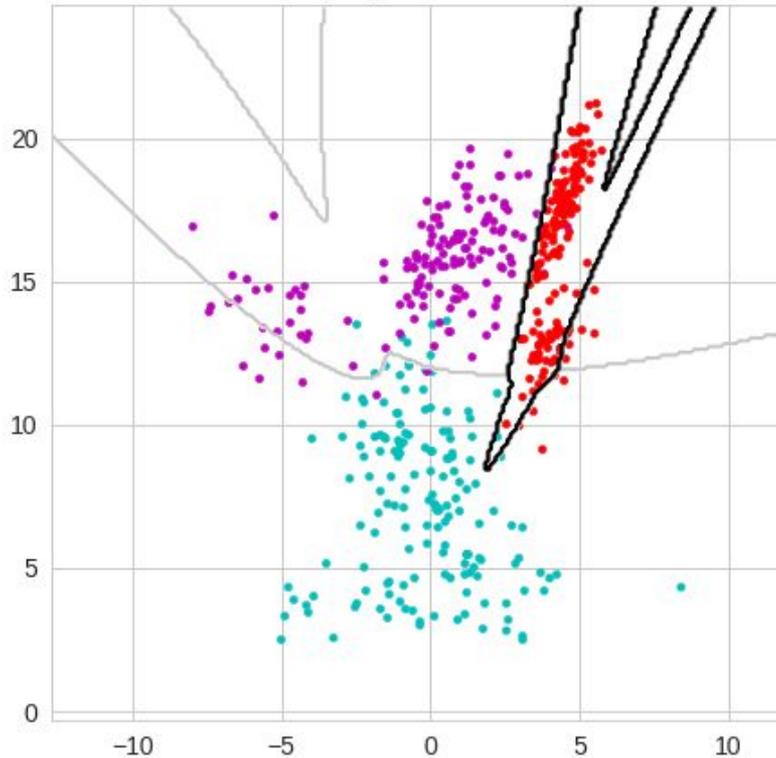




pomegranate supports semisupervised learning

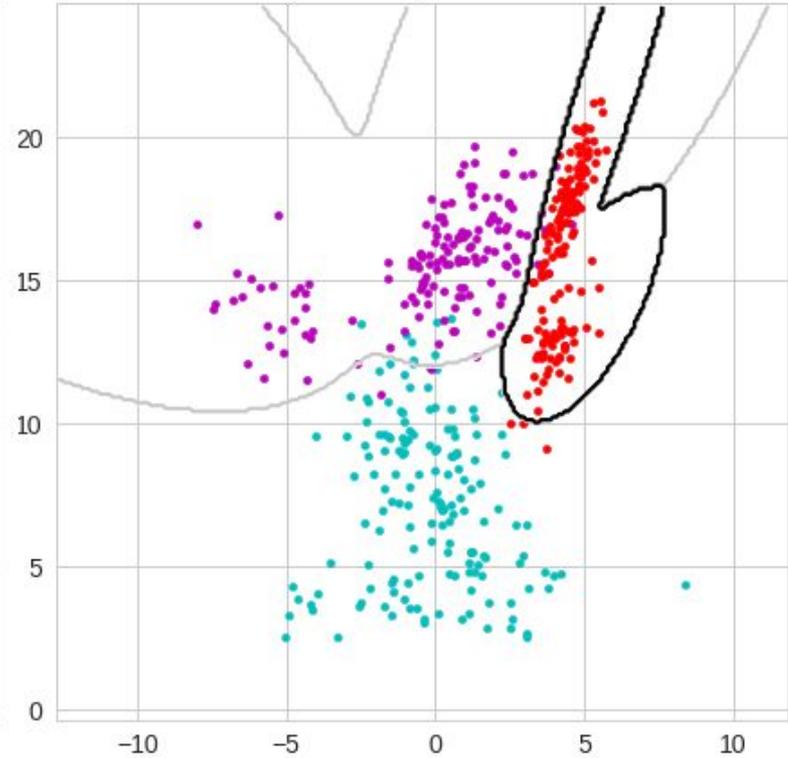
Supervised Accuracy: 0.93

Test Data, Supervised Boundaries



Semisupervised Accuracy: 0.96

Test Data, Semi-supervised Boundaries





pomegranate can be faster than scipy

```
mu, cov = numpy.random.randn(2000), numpy.eye(2000)
d = MultivariateGaussianDistribution(mu, cov)
X = numpy.random.randn(2000, 2000)
print "scipy time: ",
%timeit multivariate_normal.logpdf(X, mu, cov)
print "pomegranate time: ",
%timeit MultivariateGaussianDistribution(mu, cov).log_probability(X)
print "pomegranate time (w/ precreated object): ",
%timeit d.log_probability(X)
```

```
scipy time: 1 loop, best of 3: 1.67 s per loop
pomegranate time: 1 loop, best of 3: 801 ms per loop
pomegranate time (w/ precreated object): 1 loop, best of 3: 216 ms per loop
```



pomegranate uses aggressive caching

$$P(X|\mu, \sigma) = \frac{1}{\sqrt{2\pi}\sigma} \exp\left(-\frac{(x - \mu)^2}{2\sigma^2}\right)$$

$$\log P(X|\mu, \sigma) = -\log(\sqrt{2\pi}\sigma) - \frac{(x - \mu)^2}{2\sigma^2}$$

$$\log P(X|\mu, \sigma) = \alpha - \frac{(x - \mu)^2}{\beta}$$



GOSSIP GIRL



Example 'blast' from Gossip Girl

Spotted: Lonely Boy. Can't believe the love of his life has returned. If only she knew who he was. But everyone knows Serena. And everyone is talking. Wonder what Blair Waldorf thinks. Sure, they're BFF's, but we always thought Blair's boyfriend Nate had a thing for Serena.



Example 'blast' from Gossip Girl

Why'd she leave? Why'd she return? Send me all the deets.
And who am I? That's the secret I'll never tell. The only one.
—XOXO. Gossip Girl.



How do we encode these 'blasts'?

Better lock it down with Nate, B. Clock's ticking.

+1 Nate

-1 Blair



How do we encode these 'blasts'?

This just in: S and B committing a crime of fashion. Who doesn't love a five-finger discount. Especially if it's the middle one.

-1 Blair

-1 Serena

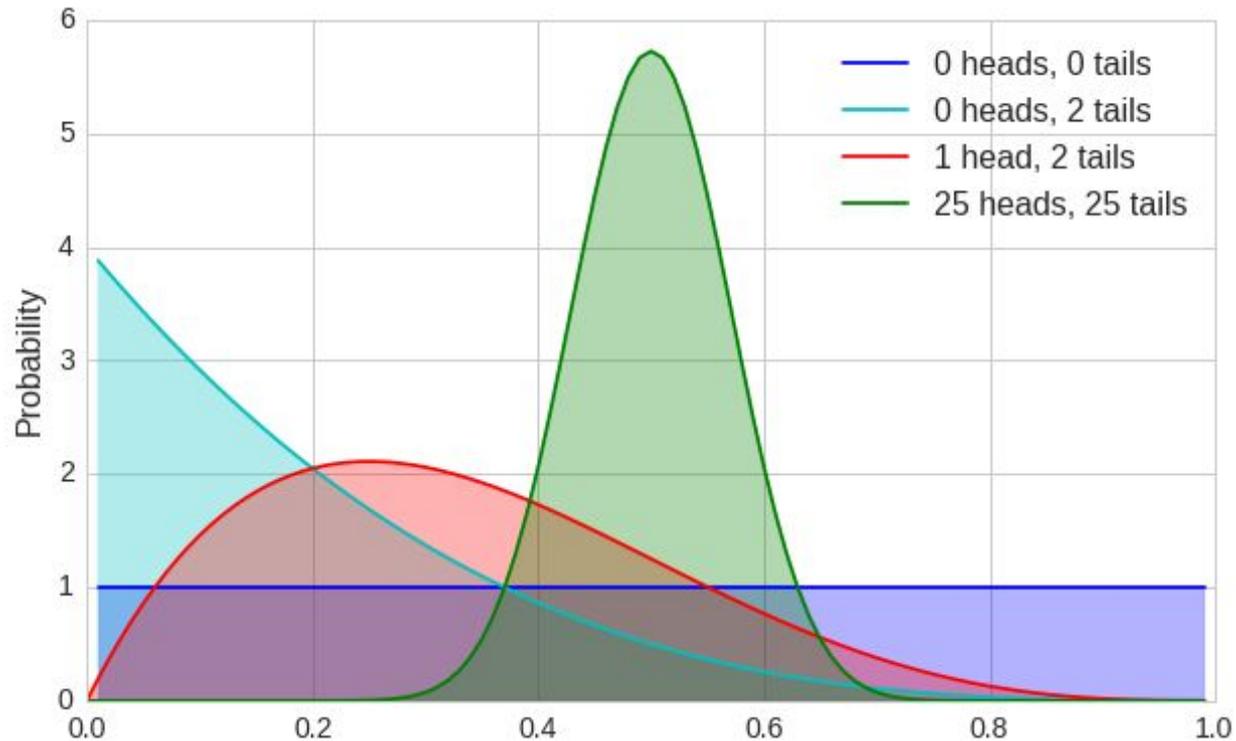


Simple summations don't work well



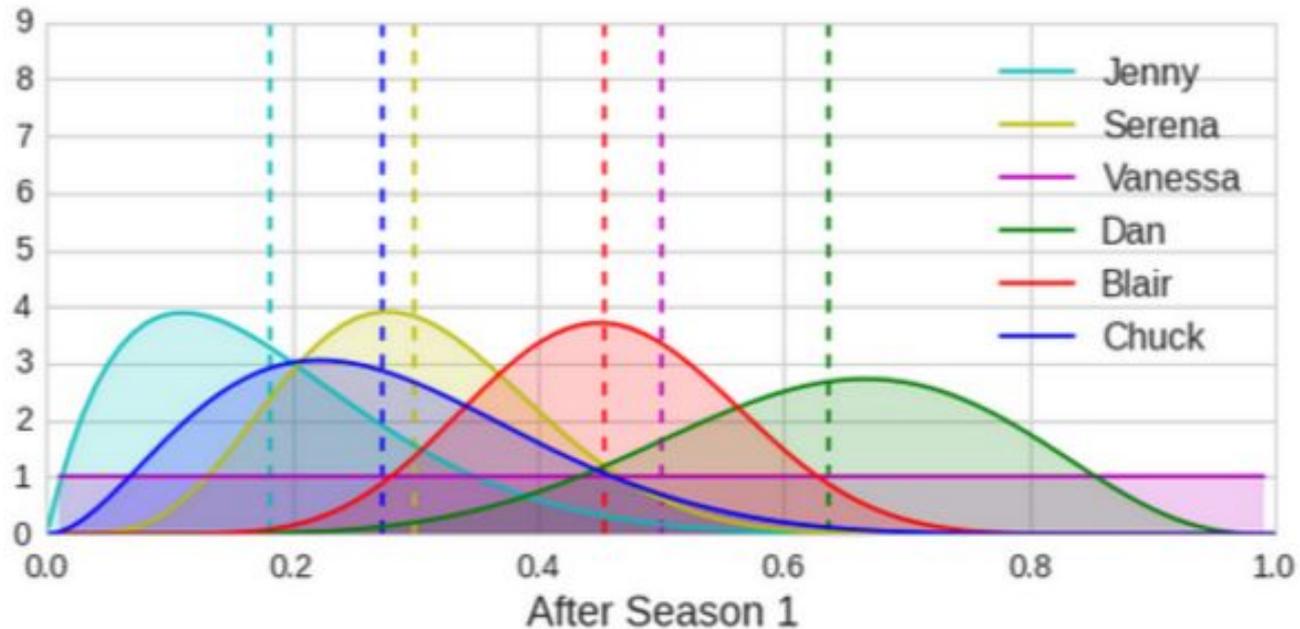


Beta distributions can model uncertainty



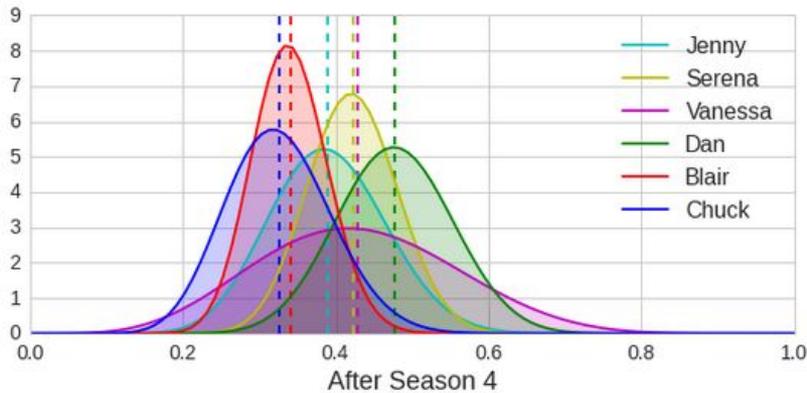
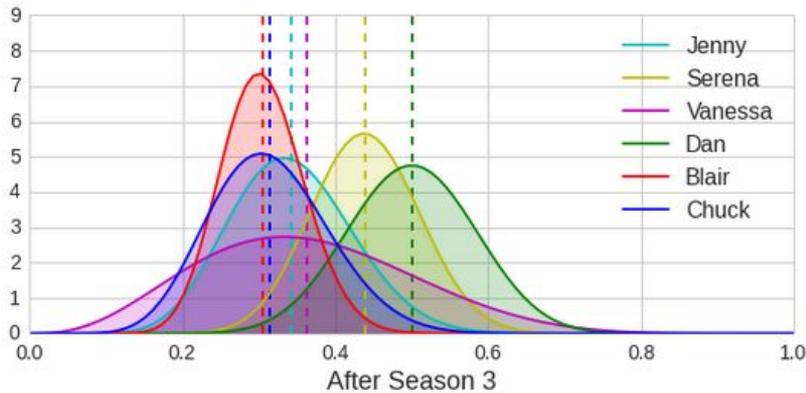
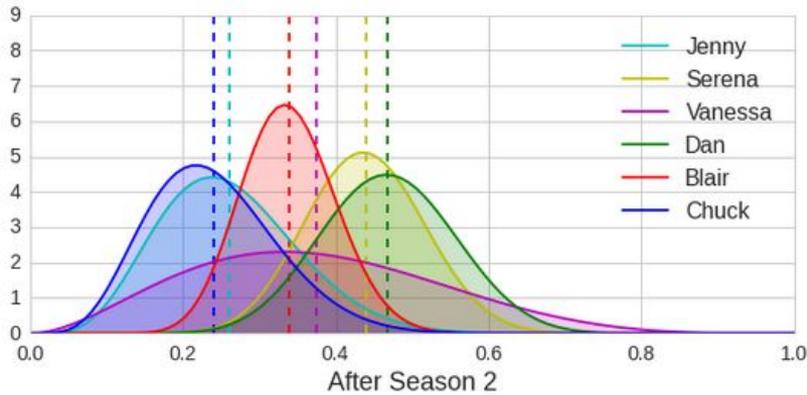
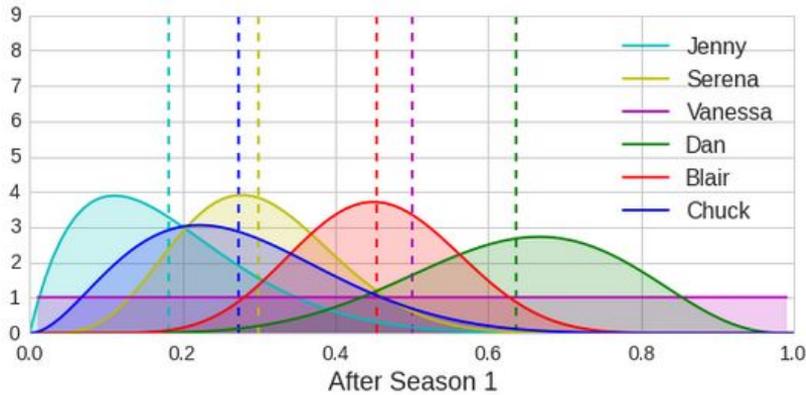


Beta distributions can model uncertainty



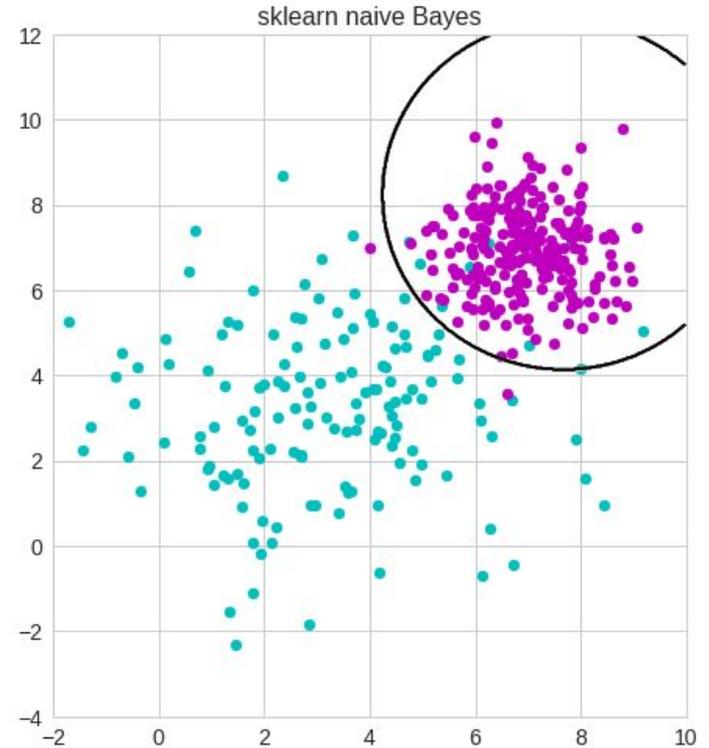
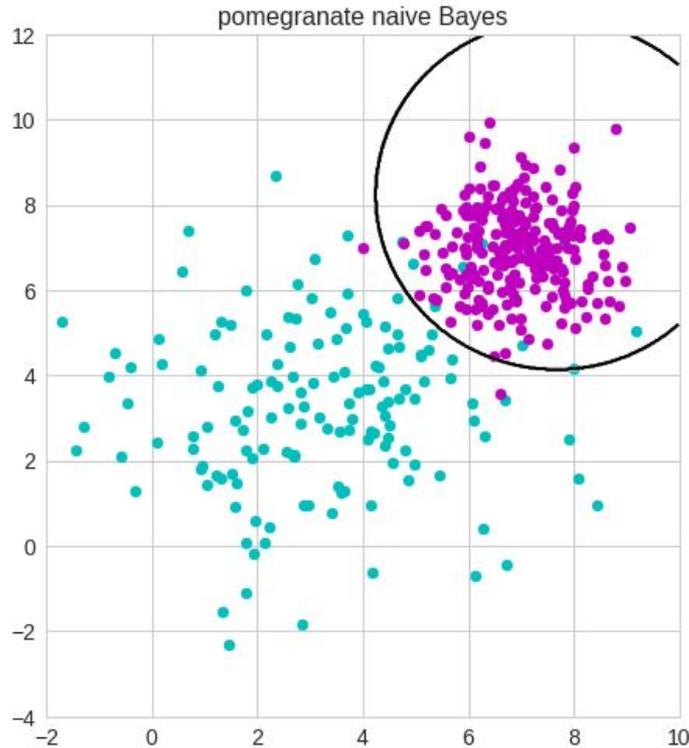


Beta distributions can model uncertainty





Naive Bayes produces ellipsoid boundaries



```
model = NaiveBayes.from_samples(NormalDistribution, X, y)
```



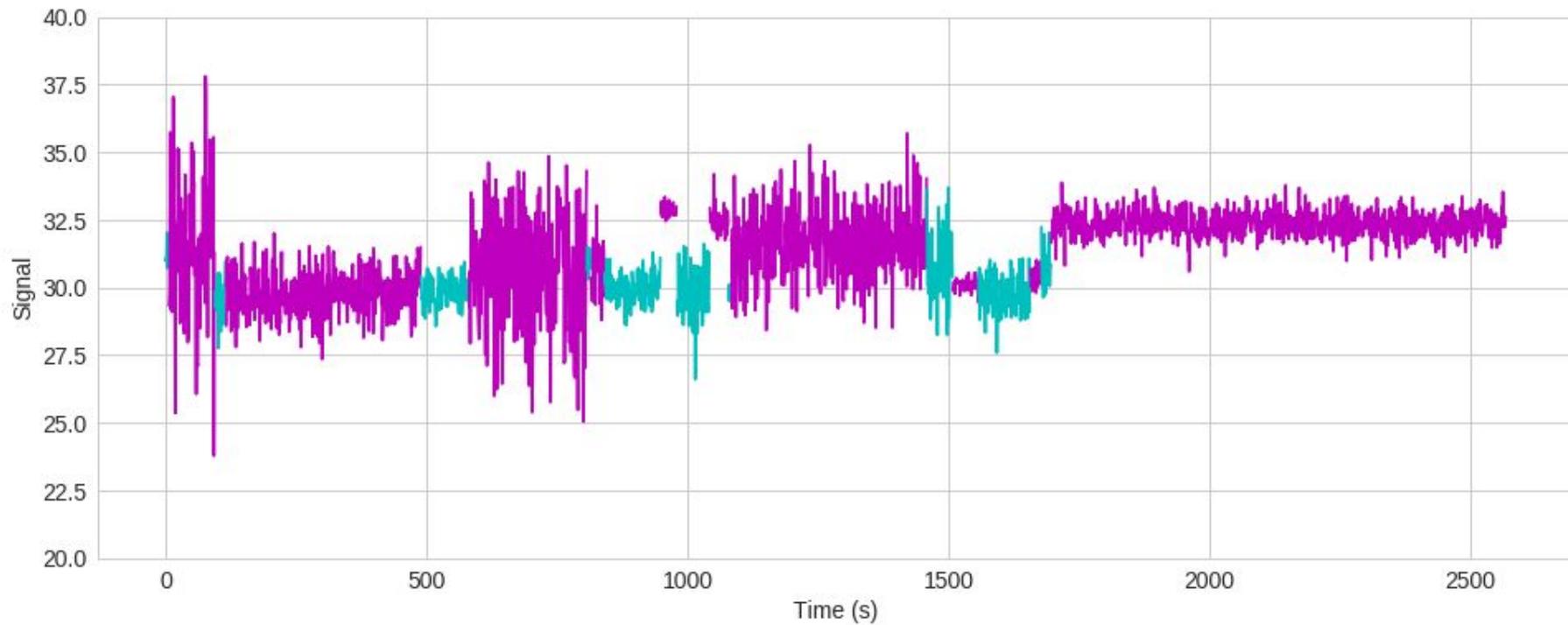
Naive Bayes assumes independent features

$$\textit{Posterior} = \frac{\textit{Likelihood} * \textit{Prior}}{\textit{Normalization}}$$

$$P(M|D) = \frac{\prod_{i=1}^d P(D_i|M)P(M)}{\sum_M \prod_{i=1}^d P(D_i|M)P(M)}$$



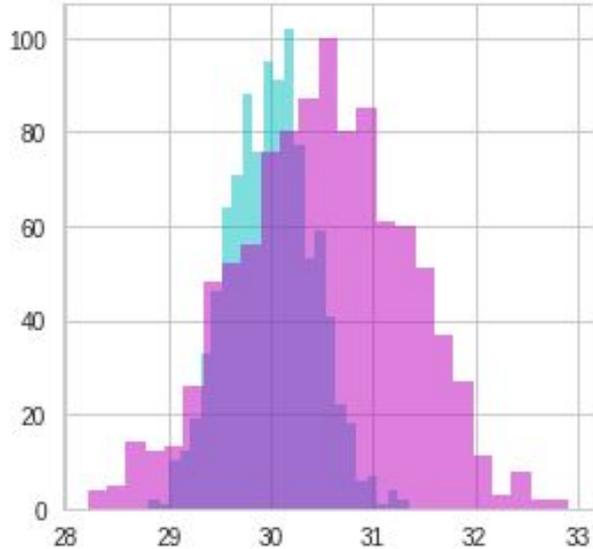
Naive Bayes can be heterogenous



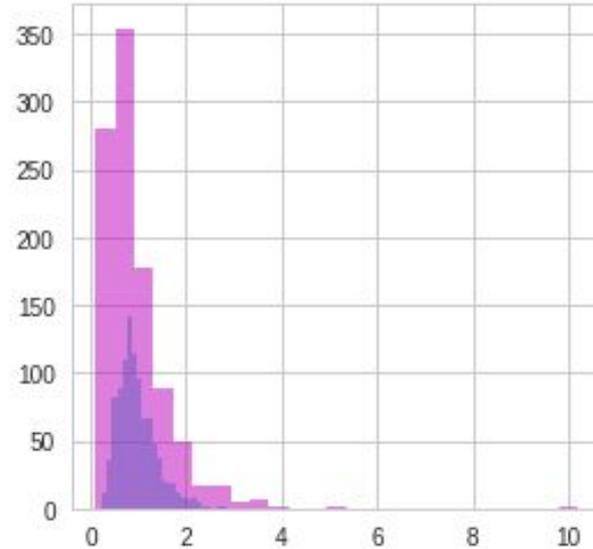


Data can fall under different distributions

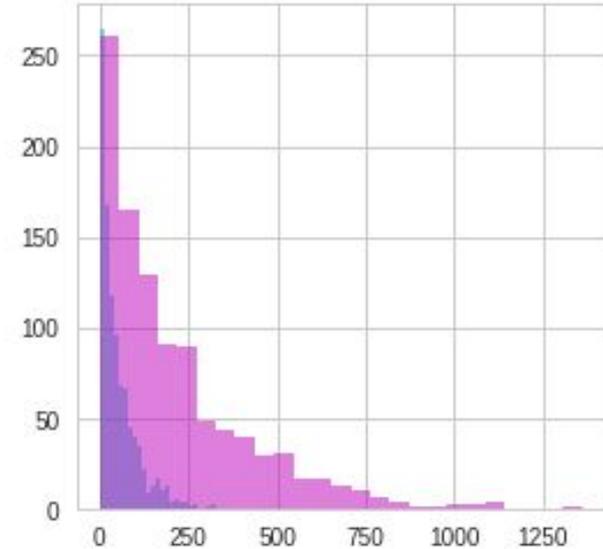
Mean



Standard Deviation



Duration





Using appropriate distributions is better

```
model = NaiveBayes.from_samples(NormalDistribution, X_train, y_train)
print "Gaussian Naive Bayes: ", (model.predict(X_test) == y_test).mean()
```

```
clf = GaussianNB().fit(X_train, y_train)
print "sklearn Gaussian Naive Bayes: ", (clf.predict(X_test) == y_test).mean()
```

```
model = NaiveBayes.from_samples([NormalDistribution, LogNormalDistribution,
ExponentialDistribution], X_train, y_train)
print "Heterogeneous Naive Bayes: ", (model.predict(X_test) == y_test).mean()
```

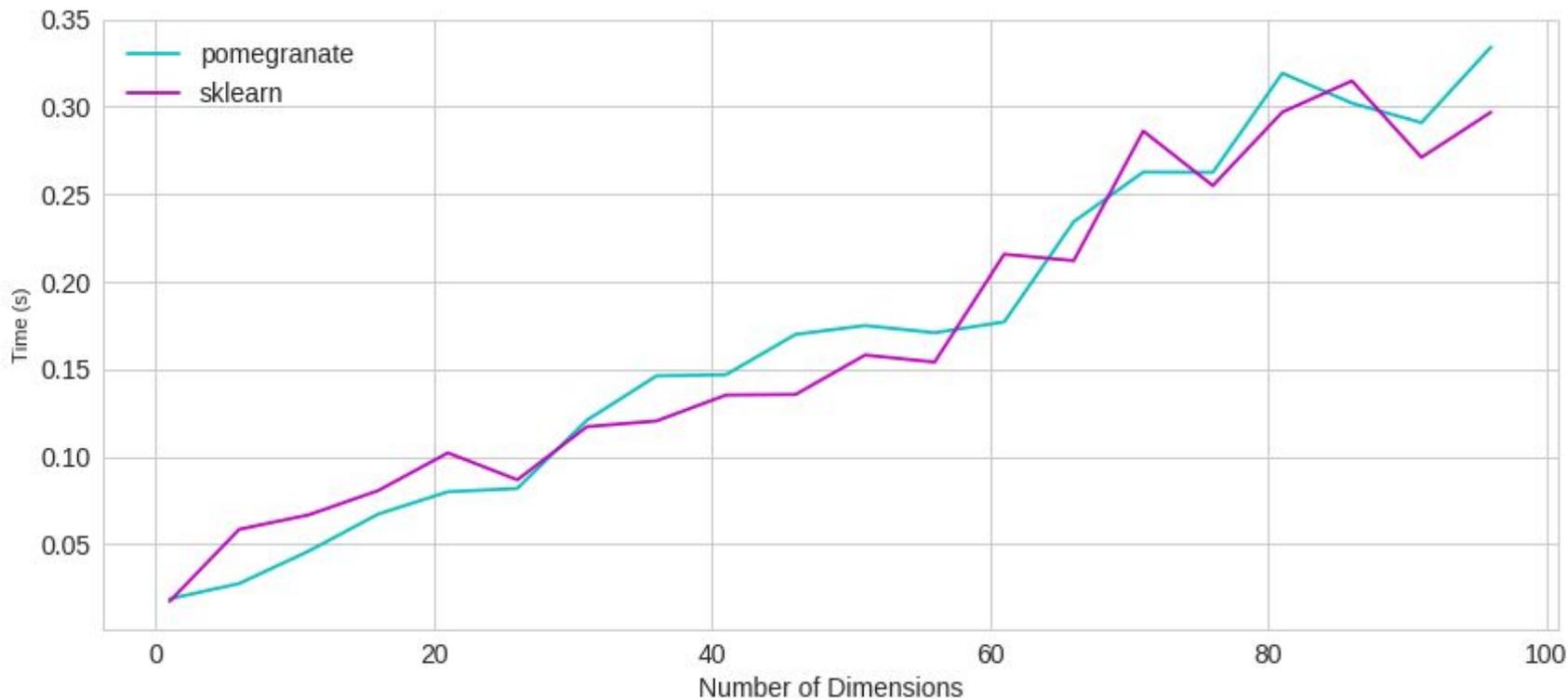
Gaussian Naive Bayes: 0.798

sklearn Gaussian Naive Bayes: 0.798

Heterogeneous Naive Bayes: 0.844



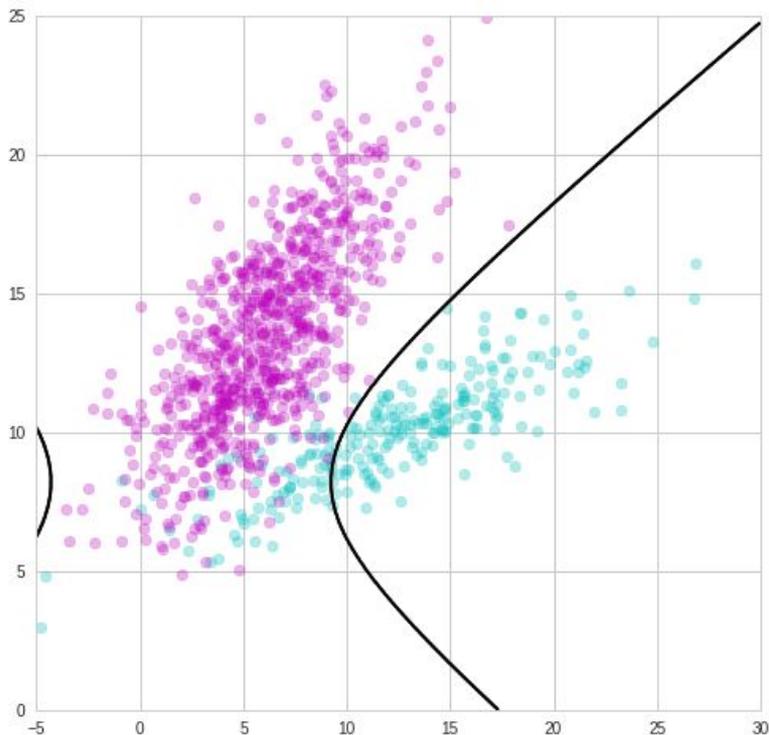
This additional flexibility is just as fast



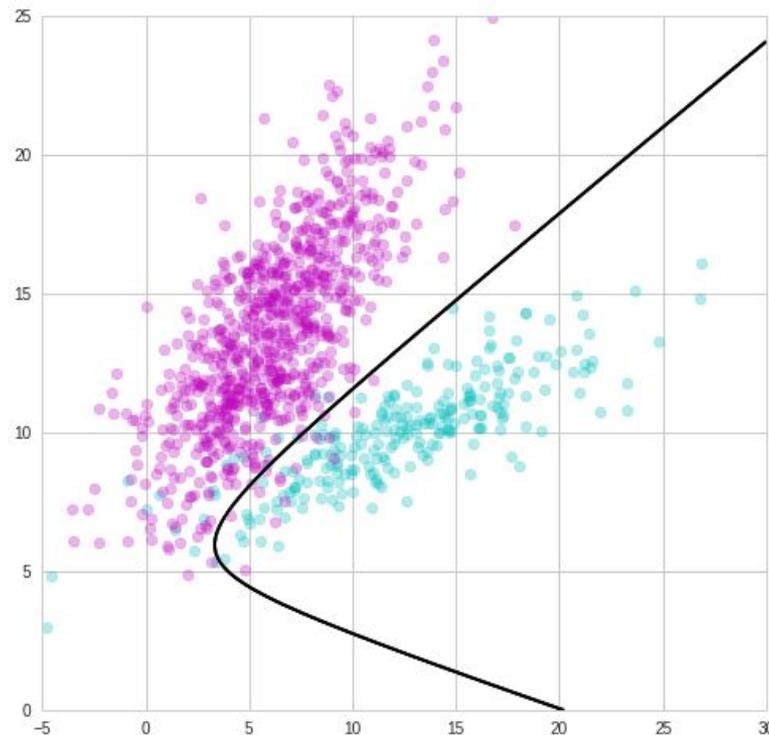


Bayes classifiers don't require independence

naive accuracy: 0.929

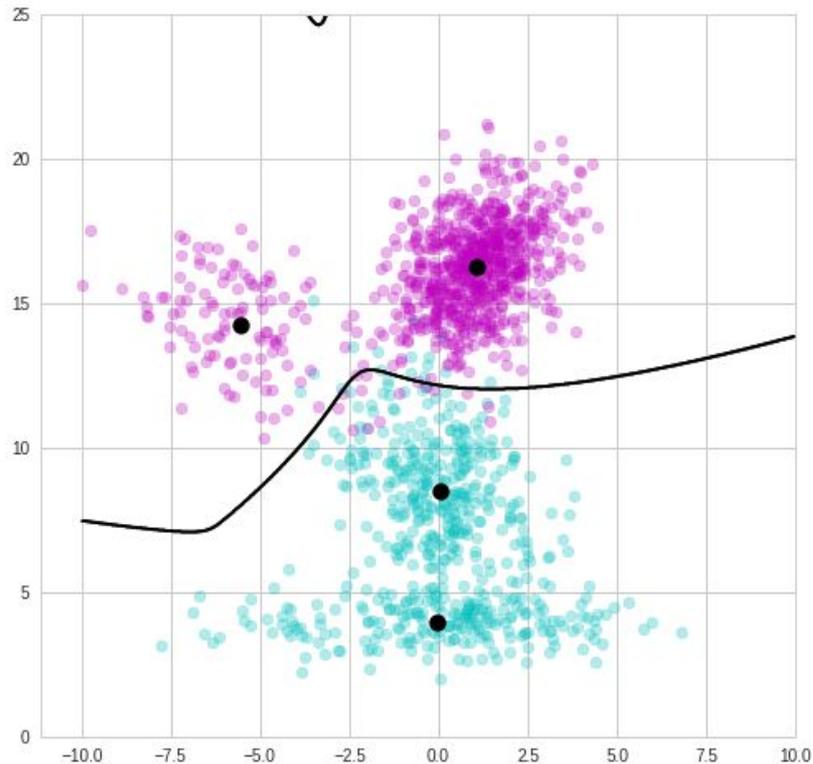
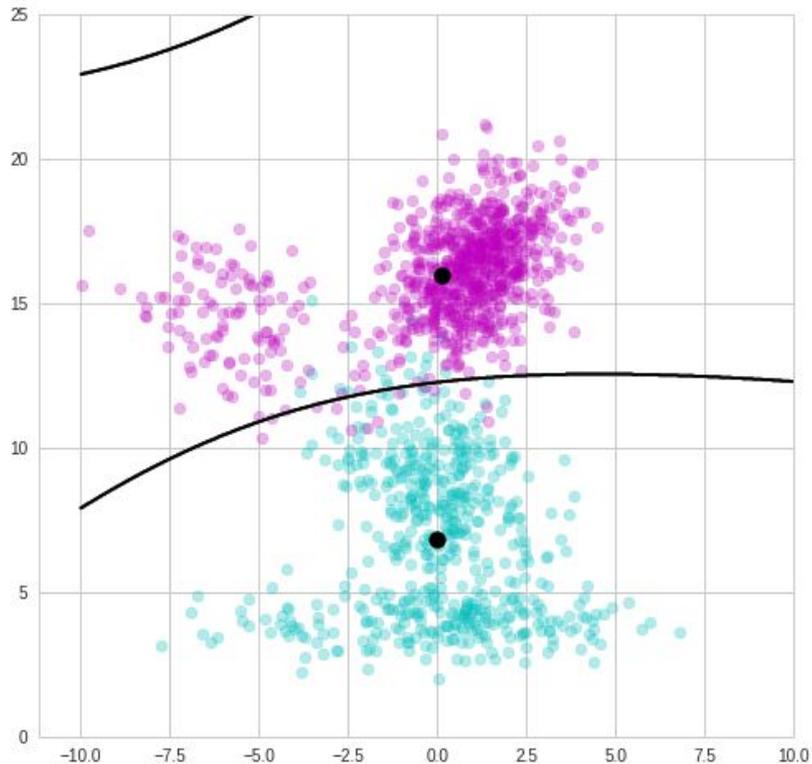


bayes classifier accuracy: 0.966





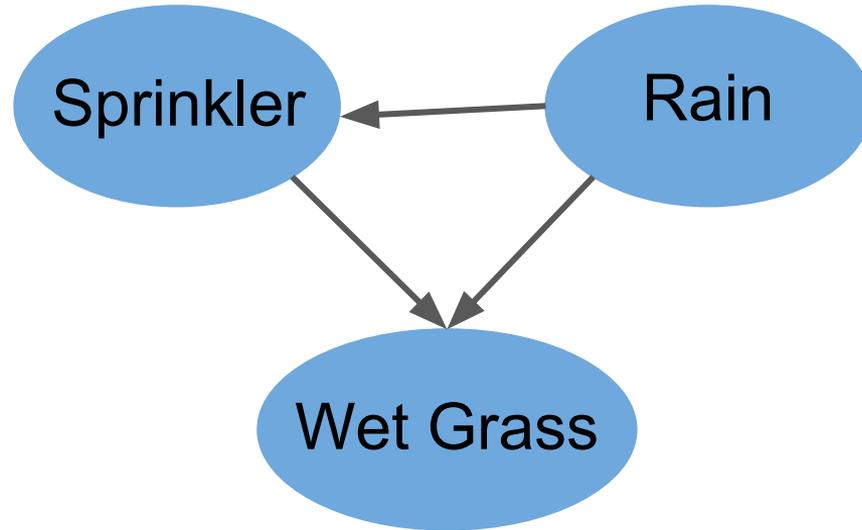
Gaussian mixture model Bayes classifier





Bayesian networks

Bayesian networks are powerful inference tools which define a dependency structure between variables.

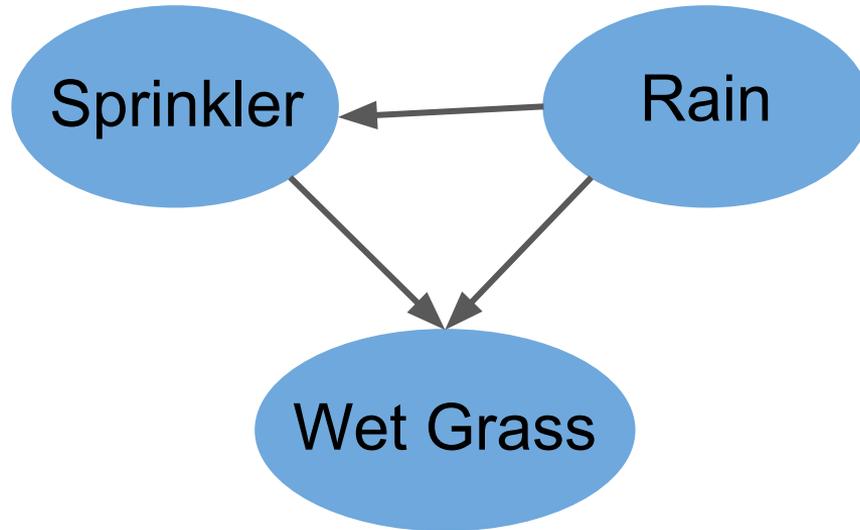




Bayesian networks

Two main difficult tasks:

- (1) Inference given incomplete information
- (2) Learning the dependency structure from data





Bayesian network structure learning

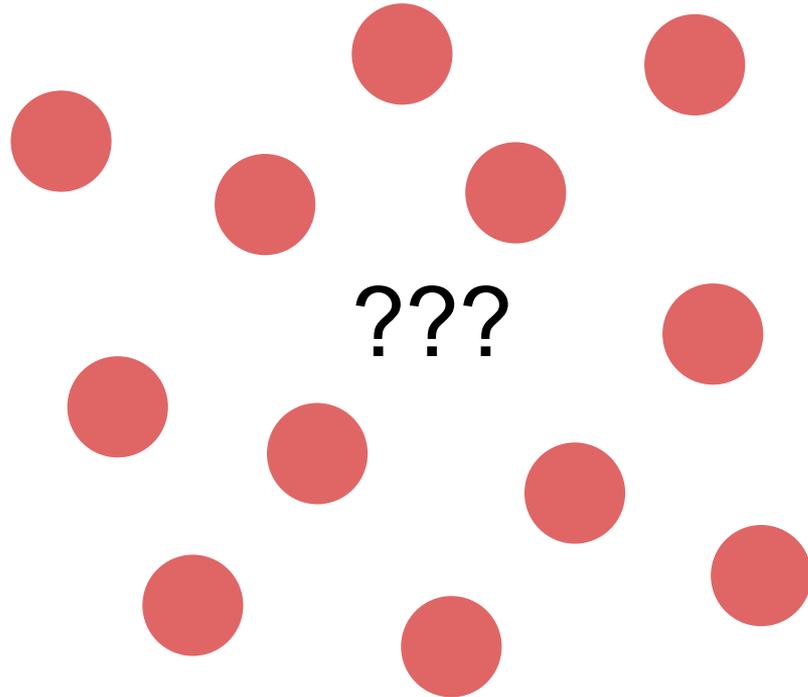
???

Three primary ways:

- “Search and score” / Exact
- “Constraint Learning” / PC
- Heuristics



Bayesian network structure learning



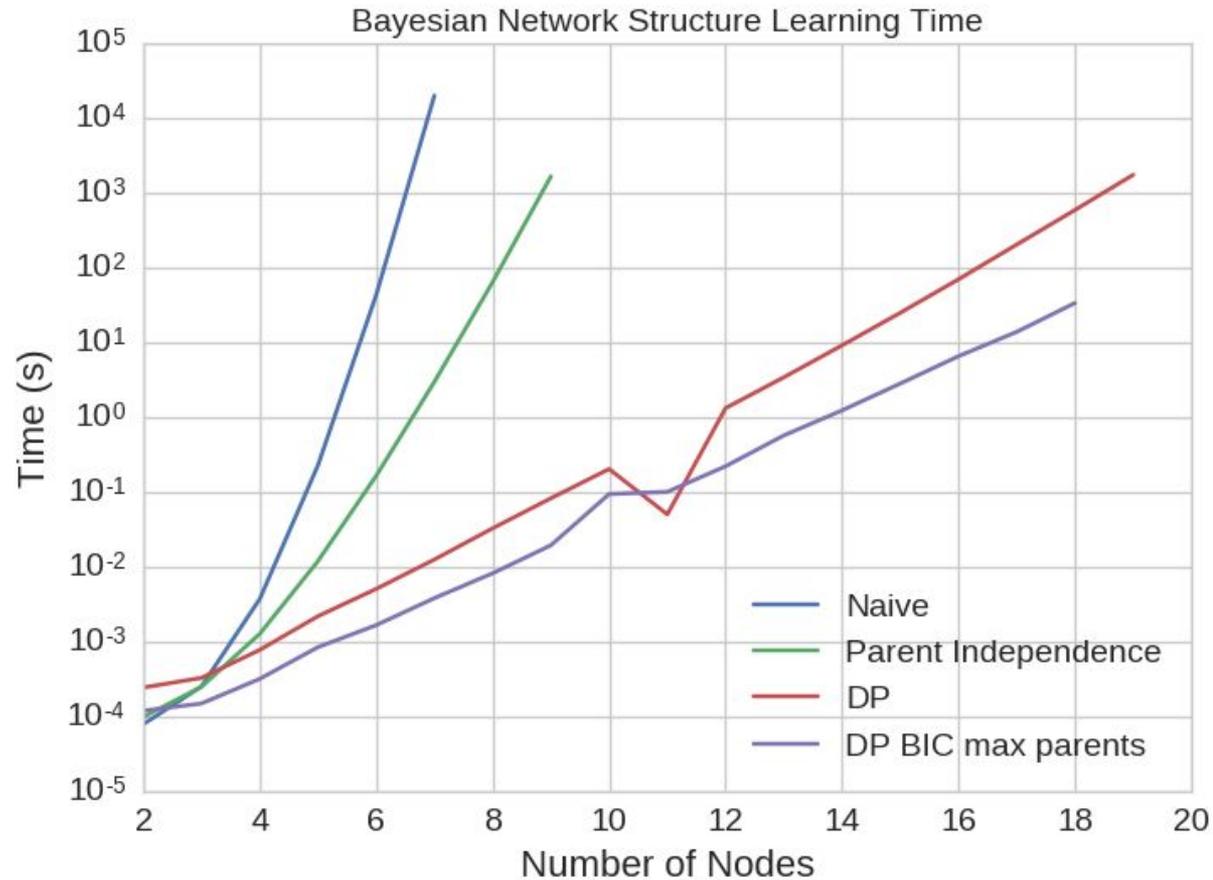
pomegranate supports:

- “Search and score” / Exact
- “Constraint Learning” / PC
- Heuristics



Exact structure learning is intractable

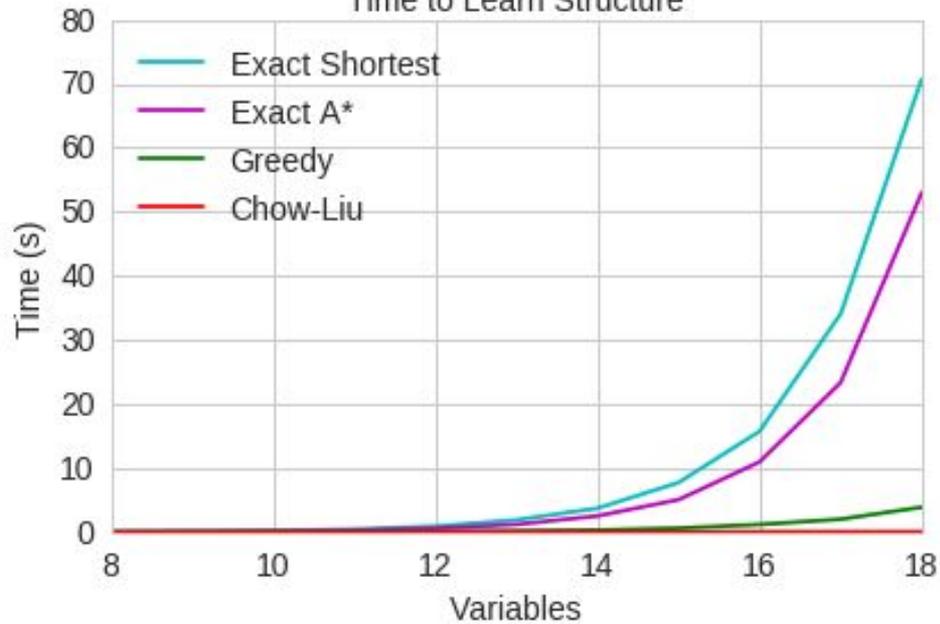
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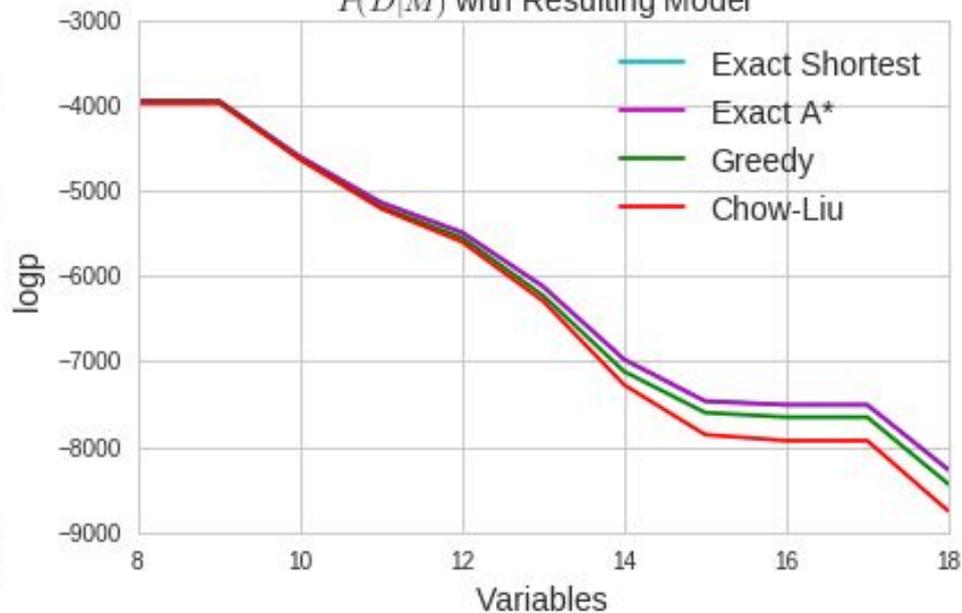


pomegranate supports four algorithms

Time to Learn Structure

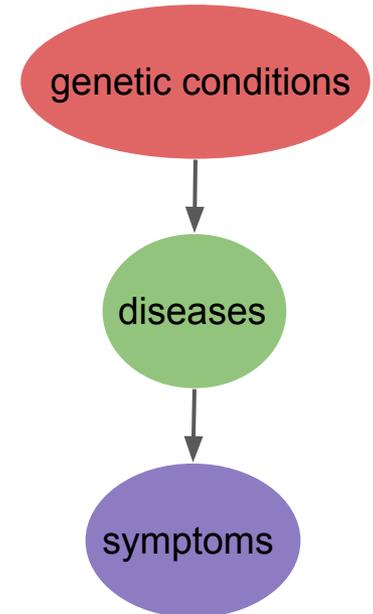
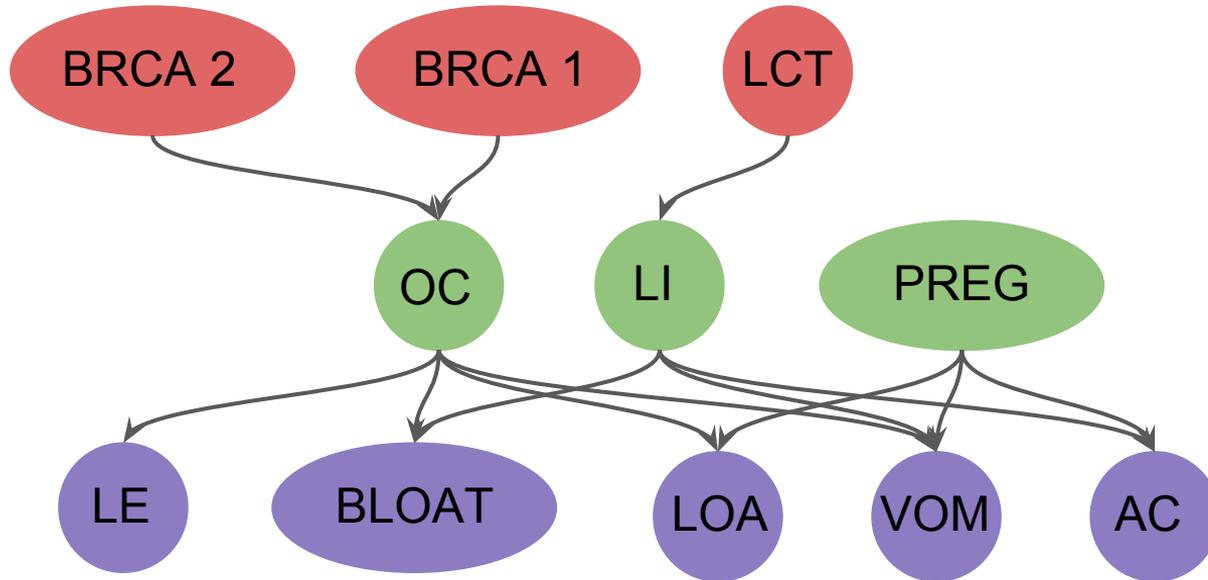


$P(D|M)$ with Resulting Model



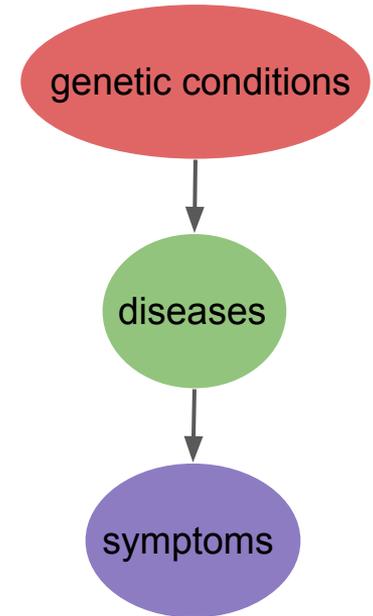
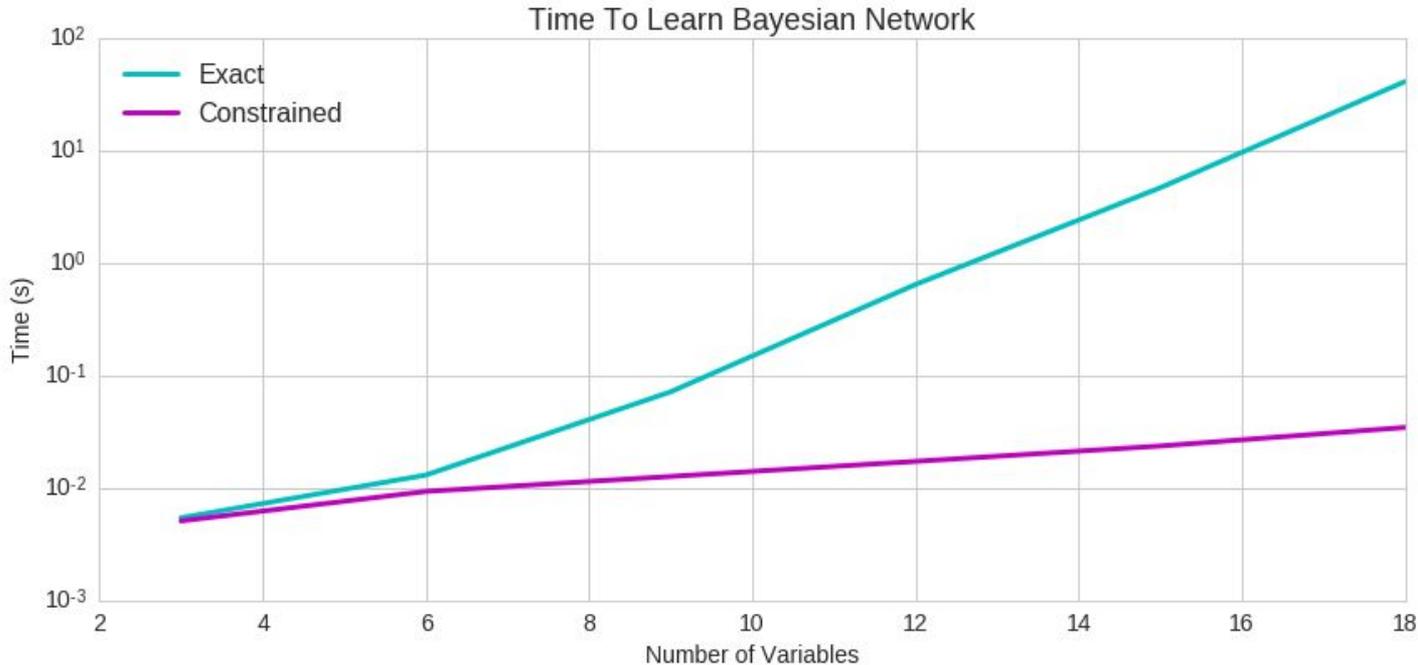


Constraint graphs merge data + knowledge



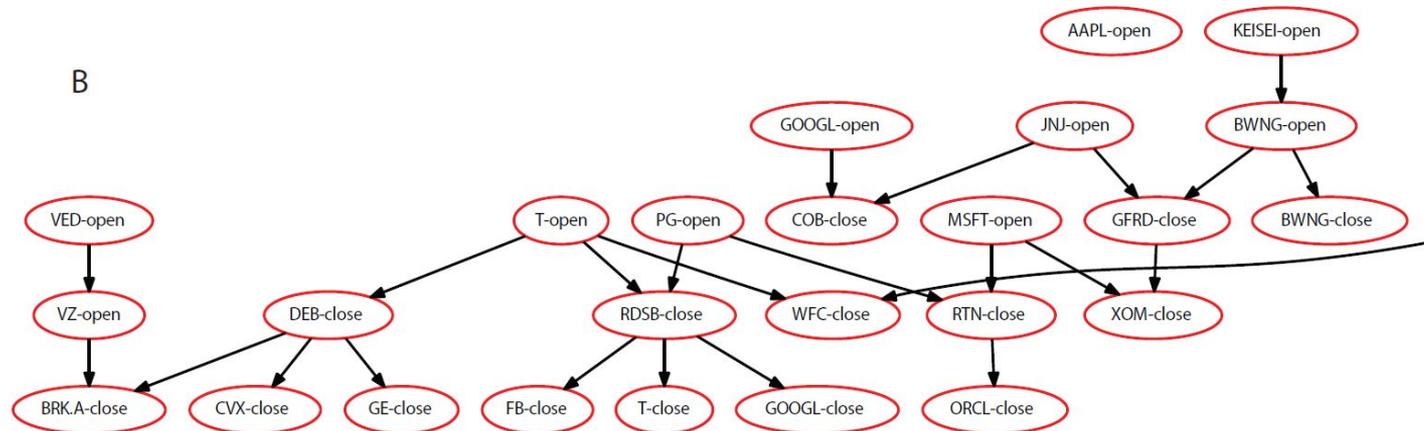
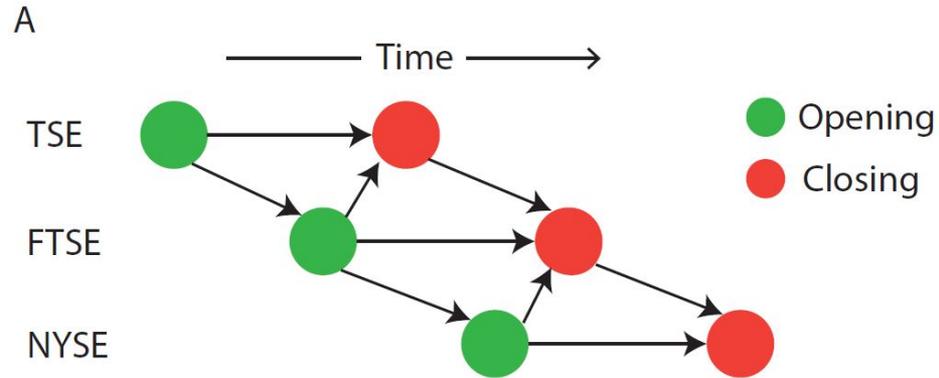


Constraint graphs merge data + knowledge





Modeling the global stock market





Finding the optimal Bayesian network given a constraint graph

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ABSTRACT

Despite recent algorithmic improvements, learning the optimal structure of a Bayesian network from data is typically infeasible past a few dozen variables. Fortunately, domain knowledge can frequently be exploited to achieve dramatic computational savings, and in many cases domain knowledge can even make structure learning tractable. Several methods have previously been described for representing this type of structural prior



Overview

pomegranate is **more flexible** than other packages, **faster**, is **intuitive to use**, and can do it all **in parallel**



Paper preprint available on arxiv!

pomegranate: fast and flexible probabilistic modeling in python

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Abstract

We present pomegranate, an open source machine learning package for probabilistic modeling in Python. Probabilistic modeling encompasses a wide range of methods that explicitly describe uncertainty using probability distributions. Three widely used probabilistic models implemented in pomegranate are general mixture models, hidden Markov models, and Bayesian networks. A primary focus of pomegranate is to abstract away the complexities of training models from their definition. This allows users to focus on specifying the correct model for their



pomegranate is now NumFOCUS affiliated



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pomegranate

pomegranate is a Python module for fast and flexible probabilistic modeling inspired by the design of scikit-learn. A primary focus of pomegranate is to abstract away the intricacies of a model from its definition, allowing users to easily prototype with complex models and training strategies. Its modular implementation allows for probability distributions to be swapped in or out for each other with ease and for models to be stacked within each other, yielding such delights as a mixture of Bayesian networks or a Gaussian mixture model Bayes classifier.

<https://www.numfocus.org/open-source-projects/affiliated-projects/>



Documentation available at Readthedocs

🏠 pomegranate
latest

GETTING STARTED

Home

- Installation
- FAQ
- Release History

FEATURES

- Out of Core Learning
- Semi-Supervised Learning
- Parallelism
- GPU Usage

MODELS

- Probability Distributions
- General Mixture Models
- Hidden Markov Models
- Bayes Classifiers and Naive Bayes
- Markov Chains

Docs » Home Edit on GitHub

pomegranate

build passing build passing docs passing

Home

pomegranate is a python package which implements fast, efficient, and extremely flexible probabilistic models ranging from probability distributions to Bayesian networks to mixtures of hidden Markov models. The most basic level of probabilistic modeling is the a simple probability distribution. If we're modeling language, this may be a simple distribution over the frequency of all possible words a person can say.

- [Probability Distributions](#)

The next level up are probabilistic models which use the simple distributions in more complex ways. A markov chain can extend a simple probability distribution to say that the probability of a certain word depends on the word(s) which have been said previously. A hidden Markov model may say that the probability of a certain words depends on the latent/hidden state of the previous word,

<https://pomegranate.readthedocs.io/en/latest/>



Tutorials available on github

Branch: master ▾ pomegranate / tutorials / Create new file Upload files Find file History

 jmschrei	ADD bayes backend	Latest commit 724510d 10 hours ago
..		
 GGBlasts.xlsx	PyData Chicago 2016	8 months ago
 PyData_2016_Chicago_Tutorial.ipynb	FIX markov chain notebooks	3 months ago
 README.md	Update README.md	2 years ago
 Tutorial_0_pomegranate_overview.ipynb	Minor typos	3 months ago
 Tutorial_1_Distributions.ipynb	ENH tutorials	2 years ago
 Tutorial_2_General_Mixture_Models.ipynb	FIX hmm dimensionality	11 months ago
 Tutorial_3_Hidden_Markov_Models.ipynb	edit tutorial 3 to remove deprecated bake	7 months ago
 Tutorial_4_Bayesian_Networks.ipynb	ENH pomegranate vs libpgm tutorial	7 months ago
 Tutorial_4b_Bayesian_Network_Structure_Learning.i...	ENH a* search	28 days ago
 Tutorial_5_Bayes_Classifiers.ipynb	ADD bayes backend	10 hours ago
 Tutorial_6_Markov_Chain.ipynb	FIX markov chain notebooks	3 months ago
 Tutorial_7_Parallelization.ipynb	ADD tutorial 7 parallelization	8 months ago

<https://github.com/jmschrei/pomegranate/tree/master/tutorials>

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PyMC3, Edward, PyStan?

Pomegranate implements probabilistic models that do not require samplers perform inference with, whereas these packages focus on the implementation of efficient samplers

Model hyperparameters in pomegranate are numbers, whereas they are typically distributions in these other packages. This allows uncertainty in model parameters to be explicitly captured.

Pomegranate focuses on discrete latent state (but discrete/continuous observed state) whereas these focus on continuous latent state