## Q

## Bernoulli Naive Bayes

If *X* is a Bernoulli-distributed random variable, it can have only two possible outcomes (for simplicity, let's call them 0 and 1) and their probability is this:

$$p(X) = \left\{egin{array}{ll} p & if & X=1 \ q & if & X=0 \end{array}
ight. egin{array}{ll} where & q=1-p & and & 0$$

In general, the input vectors  $x_i$  are assumed to be multivariate Bernoulli distributed and each feature is binary and independent. The parameters of the model are learned according to a frequency count. Hence, if there are n samples with m features, the probability for the i<sup>th</sup> feature is this  $(N_x^n)$  counts the number of times the i<sup>th</sup> = 1):

$$p_i = rac{N_{ar{x}^{(i)}} = 1}{n}$$

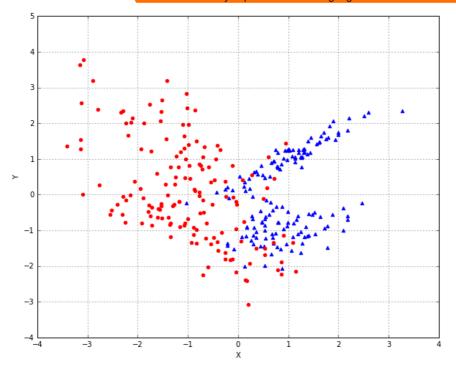
To test this algorithm with scikit-learn, we're going to generate a dummy dataset. Bernoulli Naive Bayes expects binary feature vectors; however, the BernoulliNB class has a binarize parameter, which allows us to specify a threshold that will be used internally to transform the features:

from sklearn.datasets import make\_classification

nb samples = 300

X, Y = make\_classification(n\_samples=nb\_samples, n\_features=2, n\_informative=2, n\_re

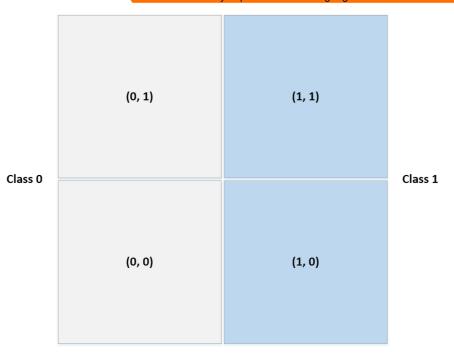
We have generated the bidimensional dataset shown in the following graph:



## Dataset for Bernoulli Naive Bayes test

We have decided to use 0.0 as a binary threshold, so each point can be characterized by the quadrant where it's located. Of course, this is a rational choice for our dataset, but Bernoulli Naive Bayes is envisaged for binary feature vectors or continuous values, which can be precisely split with a predefined threshold:

The score is rather good, but if we want to understand how the binary classifier worked, it's useful to see how the data has been internally binarized:



Structure of the binarized dataset

Now, checking the Naive Bayes predictions, we obtain the following:

```
data = np.array([[0, 0], [0, 1], [1, 0], [1, 1]])
print(bnb.predict(data))
[0, 0, 1, 1]
```