Factoring with conditionals

Earlier we saw that we can write a conditional distribution as the ratio of a joint and a marginal:

$$P(X \mid Y) = rac{P(X,Y)}{P(Y)}$$

A useful rearrangement of the above is

$$P(X,Y) = P(X \mid Y)P(Y)$$

That is, we can write a joint distribution as a product of a marginal and a conditional. (By symmetry it works the other way too: $P(X, Y) = P(Y \mid X)P(X)$.)

Since a conditional distribution is just like any other distribution, the same identity works inside a conditional:

$$P(X, Y \mid Z) = P(X \mid Y, Z)P(Y \mid Z)$$

So, we can use this identity repeatedly to break up a large joint distribution into a product of factors: e.g.,

$$P(X, Y, Z) = P(X, Y \mid Z)P(Z) = P(X \mid Y, Z)P(Y \mid Z)P(Z)$$

Bayes' rule

From the identity above, we know

$$P(X \mid Y)P(Y) = P(X, Y) = P(Y \mid X)P(X)$$

Rearranging, we have

$$P(Y \mid X) = \frac{P(X \mid Y)P(Y)}{P(X)}$$

This is Bayes' rule.

Bayes' rule is really useful since it tells us how to incorporate new evidence about an uncertain variable. For example, suppose there are two coins in a bag: one fair one, and one that's biased so that the probability of heads is 0.7. We draw one coin at random from the bag and start flipping it.

Initially we are equally likely to have drawn the fair coin or the biased one: P(fair) = 0.5. Now suppose we flip our coin and see heads. Intuitively, this outcome is more likely if we're holding the biased coin, so P(fair) should decrease.

Bayes' rule tells us that

$$P(\text{fair} \mid \text{flip} = H) = P(\text{flip} = H \mid \text{fair})P(\text{fair})/P(\text{flip} = H)$$

In words, we start from our *prior* probability table P(fair):

fair	0.5
¬fair	0.5

We multiply it by the evidence $P(\text{flip} = H \mid \text{fair})$

$H \mid fair$	0.5
$H \mid \neg$ fair	0.7

This gets us

$H \wedge fair$	$0.5\cdot 0.5=0.25$
$H \wedge \neg fair$	$0.5 \cdot 0.7 = 0.35$

We then divide by the marginal probability of our observation P(H). We can either calculate P(H) directly, or use a shortcut: $P(H) = P(H \land fair) + P(H \land \neg fair)$, so P(H) is the sum of the entries in the table above. Either way we get P(H) = 0.6, so our final answer becomes:

fair $\mid H$	0.25/0.6 = 5/12
$ egthinspace{-1mm}{\neg} fair \mid H$	0.35/0.6 = 7/12

This is called the *posterior* probability of fair (given the evidence of seeing H).

If we flip the coin again, we can repeat the exercise: our posterior after the first flip becomes our prior before the second flip, and we use Bayes' rule again to get our posterior after the second flip. This process — repeatedly updating a distribution over some variable using new evidence — is often called *tracking* or *filtering*.

As this example shows, the words "prior" and "posterior" don't refer to a single distribution. Instead they mean the distribution before and after we incorporate some piece of evidence. If we incorporate several pieces of evidence, we have to know which one we're talking about in order to know which distribution is the prior and which is the posterior. This ambiguity is a common source of confusion, so it's best to specify precisely if there's any chance of misinterpretation.

Exercise: suppose we see heads again. What is our posterior P(fair | H, H)?

Factoring a probability distribution

When we use probability spaces in practice, one of the most important tasks is to describe complicated relationships among many different possible events and random variables. A good way to organize these relationships is to *factor* our probability distribution: we write

$$P(X) = F_1(X)F_2(X)\dots F_n(X)$$

where X stands for a list of all the random variables or events that we might want to reason about, and each factor $F_k(X)$ encodes some understandable part of our overall model. This factorization means that the probability of X taking a given value x is

$$P(X=x)=F_1(x)F_2(x)\ldots F_n(x)$$

There are lots of possible kinds of factors we might want to include. We can't cover them all, but the rest of this set of notes will look at a few useful kinds, and how to work with them.

In all of our examples, the factors $F_k(X)$ will be simpler probability distributions. As we'll see, this is nice for interpretability. But there do exist factored distributions where the factors can't be interpreted this way.

Independence

The simplest kind of relationship is none at all: suppose we can split our random variables into two or more subsets that don't influence one another. Then these subsets are called *independent* of one another.

We can represent independence by writing our overall probability as a product of two or more factors, where the factors have *disjoint* sets of arguments. Like this:

$$P(X_1, X_2) = P(X_1)P(X_2)$$

For example, if we flip a coin twice, it makes sense to assume that the two flips are independent: getting heads on one flip doesn't influence our chance of getting heads on the other. The above formula could represent our joint distribution, if we take X_1 to represent the first flip and X_2 to represent the second.

If we write x_1 for a value that X_1 might take, and x_2 for a value that X_2 might take, the above factorization stands for a table in which

$$P(X_1 = x_1, X_2 = x_2) = P(X_1 = x_1)P(X_2 = x_2)$$

For example, if our first coin has probability 0.7 of showing heads, while our second coin has probability 0.6, we get the following table:

X_1	X_2	p
Н	Н	$0.7\cdot 0.6=0.42$
Н	Т	$0.7 \cdot 0.4 = 0.28$
Т	Н	$0.3 \cdot 0.6 = 0.18$
Т	Т	$0.3 \cdot 0.4 = 0.12$

We can see that this table encodes independence by calculating: suppose we observe that $X_1 = T$. Then we can use the rule for conditional probabilities to find $P(X_2 | X_1 = T)$. To do so, we cross out the rows in the table that are inconsistent with our observation; in this case, we cross out the first and second rows, and keep the third and fourth. Then we renormalize the remaining rows: the two remaining entries are 0.18, 0.12 and their sum is 0.3, so the result is

X_2	p
Н	0.18/0.3 = 0.6
Т	0.12/0.3 = 0.4

The distribution of the second coin is still a 60% probability of heads, unchanged from before we made our observation of the first coin. This property — the distribution of one random variable is unchanged when we make observations about the other — is what defines independence.

Exercise: give a joint distribution for two real variables X, Y such that their correlation is zero, but they are not independent.

Conditional independence

Sometimes our random variables aren't independent to start out, but they *become* independent after we observe something. This is called *conditional independence*. Under conditional independence, our distribution isn't a product of independent factors to start, but it becomes a product of independent factors after we make an observation.

A good example is a clustered distribution:



In this distribution, the X and Y coordinates are not independent: if X is big, then we're more likely to be in the second cluster, meaning that Y is more likely to be big as well. But once we know which cluster we're in, X and Y are independent.

If the conditional distribution is independent, that means that it factors into the product of the probability of X and the probability of Y:

$$P(X, Y \mid Z) = P(X \mid Z)P(Y \mid Z)$$

where X, Y are the coordinates of a sample and Z is the indicator of which cluster the sample is in. So, the joint is

$$P(X, Y, Z) = P(Z)P(X, Y \mid Z) = P(Z)P(X \mid Z)P(Y \mid Z)$$

We can test that this formula gives us conditional independence by conditioning on an observation — say Z = 1, i.e., the point is in the first cluster. The rule for conditional probabilities gives us

$$P(X, Y \mid Z = 1) = P(X, Y, Z = 1) / P(Z = 1)$$

= $P(Z = 1)P(X \mid Z = 1)P(Y \mid Z = 1) / P(Z = 1)$
= $P(X \mid Z = 1)P(Y \mid Z = 1)$

Since we've fixed a value for Z, P(X | Z = 1) depends only on X, and P(Y | Z = 1) depends only on Y. So, our conditional distribution is a product of two independent factors, as claimed.

Samples

One of the most common uses of independence or conditional independence is when we repeat an experiment many times to collect a sample. In this situation it makes sense to assume that each run of the experiment is independent from all of the other runs. If our sample is X_1, X_2, \ldots, X_T , that means our overall distribution factors as

$$P(X_1, X_2, \ldots, X_T) = P(X_1)P(X_2) \ldots P(X_T)$$

In a sample like this, we might have an unknown parameter vector $\theta \in \mathbb{R}^d$ that influences the distribution of our samples — e.g., θ might contain the mean and variance of a sample X_t . If we think of θ as fixed but unknown, we could emphasize the dependence on θ by writing

$$P_{ heta}(X_1, X_2, \dots, X_T) = P_{ heta}(X_1) P_{ heta}(X_2) \dots P_{ heta}(X_T)$$

On the other hand, we might want to think of θ itself as a random variable. In this case we would say that the samples X_t are *conditionally* independent given θ :

$$P(X_1, X_2, \dots, X_T, \theta) = P(\theta)P(X_1 \mid \theta)P(X_2 \mid \theta) \dots P(X_T \mid \theta)$$

These are two alternate views of the world: either the parameters take some fixed but unknown value, or the parameters are themselves random. Both views are reasonable; they often lead to similar conclusions about θ , but they can also be subtly different.

Group at a time

In general, we might have a complicated pattern of dependence, where each of our random variables depends on some of the others. A good way to organize our factorization in this case is to sort our random variables in some order, and imagine that we pick their values one or a few at a time according to our order. Each group that we pick at the same time leads to a factor in our probability distribution.

For example, we might have three random variables: whether it's raining today, whether our sprinkler is turned on, and whether the grass is wet. We could pick them in the order rain, sprinkler, wet, and write

 $P(\text{rain, sprinkler, wet}) = P(\text{rain}) P(\text{sprinkler} \mid \text{rain}) P(\text{wet} \mid \text{rain, sprinkler})$

Each factor is the conditional probability distribution of one variable or group of variables, given all of the preceding ones.

Sometimes we don't need every single one of the preceding variables to predict our current one. For example, if our sprinkler is attached to a timer, then the probability that it's turned on when we look at it doesn't depend on whether it's raining. In that case we'd omit rain as a conditioning variable for sprinkler, and write

 $P(\text{rain, sprinkler, wet}) = P(\text{rain}) P(\text{sprinkler}) P(\text{wet} \mid \text{rain, sprinkler})$

The variables that we include when predicting X_i are called its *parents*; so here, rain and sprinkler have no parents, while the parents of wet are rain and sprinkler.

Often there's one distinguished order for our variables that makes the most sense. For example, in a medical study we pick a treatment first, and observe its outcome later; or in our example above, the weather happens first, and helps determine whether the grass becomes wet later. Other times there's no one order that's better than the others; in this case we can pick any convenient order.

In either case, the order is our own modeling decision: we're free to use the "wrong" order, and the only cost will be to the interpretability and accuracy of our model. The decision of what parents to use for any given variable is ours as well: if we pick a good set of parents we might get a more accurate model, but the more parents we pick, the more complicated our factorization will become.

The only restriction is that there always has to be an ordering where each variable's parents come before it — if not, we would have a loop, where A determines B and B determines A. If we tried to write a factorization this way, the result would typically not be a valid probability distribution.