Larson and Marx Section 3.1

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Given a sample space S, we have defined probabilities in terms of a probability function that maps the power set $\mathcal{P}(S)$ into the real numbers

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It is often the case that the power set $\mathcal{P}(S)$ has far more granularity than applications require, and in fact things can be greatly simplified by "redefining" the sample space to elimate some of the detail.

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Six dice are thrown. By the multiplication rule, there are

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However, chances are we are only interested in the **sum** of the six faces that come up.

Consequently, we really only care which of the 31 possible totals

 $\{6, 7, 8, \dots, 34, 35, 36\}$

has occurred.

One way we can "simplify" a sample space with 46,656 elements down to one with 31 mutually exclusive outcomes is to define a *function* from the original sample space into the numbers from 6 to 36.

Every outcome from the original sample space is mapped into exactly one number in the set $\{6, 7, \ldots, 35, 36\}$.

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In fact, this is standard practice in the study of probability, leading to the following

Definition:

A real-valued function defined on a sample space is called a **random variable**

When we say that the function is defined on a sample space, what we mean is that the independent variable is a point in a sample space, i.e., the outcome of an experiment.

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At least in the beginning, it is somewhat helpful to think "random function"

when you see the term "random variable".

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In the context of our tossing six dice experiment, if X was a random variable defined to be the sum of the six faces, this is how we would denote the outcome that the sum was 17.

If we think of X as a function defined on S (which it is),

 $X: S \mapsto \{6, 7, \dots, 35, 36\}$

we could write

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This notation would indicate more clearly the nature of the situation, namely that an event $s \in S$ occurred and the function defining the random variable X maps this event to the number 17.

But this notation is never used, so we have to live with X = 17.

Suppose *X* is a random variable.

Let

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Random variables which assume only a finite or countable number of values are called **discrete random variables**

For any element of R, say

$$x_i \in R$$
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the collection of all sample points on which X assumes the value x_i constitutes the **event** that

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The probability of this event is denoted by

$$P(X = x_i)$$

Definition:

The function

$$f(x_i) = P(X = x_i), \quad i = 1, 2, \dots$$

which is defined on the set of values R assumed by the random variable X is called the

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Consequently, a probability distribution function can assume only nonnegative values:

 $f(x_i) \geq 0 \quad \forall i$

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Furthermore, the sum of the values of the probability distribution function over all values assumed by the random variable must be one:

$$\sum_{i} f(x_i) = 1$$

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We define a random variable *X* representing the number of heads. Actually, *X* is a function from the sample space *S* into the set $A = \{0, 1, 2, 3\}$, defined by the following table:

$$X: S \mapsto A \quad \text{such that} \begin{cases} \{t, t, t\} & \to & 0\\ \{t, t, h\}, \{t, h, t\}, \{h, t, t\} & \to & 1\\ \{t, h, h\}, \{h, t, h\}, \{h, h, t\} & \to & 2\\ \{h, h, h\} & \to & 3 \end{cases}$$

The event

X = 2

will occur if the experiment produces one of the outcomes

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We can now define a **probability distribution** f on A, the set of possible values of the random variable X.

f will be defined as:

$$f(x_i) = P(X = x_i) \qquad x_i \in \{0, 1, 2, 3\}$$

We will define f by the following table:

$$f: X \mapsto [0, 1] \quad \text{such that} \begin{cases} 0 \quad \to \quad P(X = 0) \\ 1 \quad \to \quad P(X = 1) \\ 2 \quad \to \quad P(x = 2) \\ 3 \quad \to \quad P(x = 3) \end{cases}$$

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Defined in this fashion, f is called a **probability distribution** for the random variable X.

In this situation, we also say that the random variable X has distribution f.

If the coin is fair, each of the eight outcomes in the sample space S has probability 1/8, so we expect the following probabilities for the random variable X:

$$f: X \mapsto [0,1] \quad \text{such that} \begin{cases} 0 \quad \to \quad P(X=0) = 1/8 \\ 1 \quad \to \quad P(X=1) = 3/8 \\ 2 \quad \to \quad P(x=2) = 3/8 \\ 3 \quad \to \quad P(x=3) = 1/8 \end{cases}$$

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If we add up the values of f(X) over the domain of f, which is $\{0, 1, 2, 3\}$, the set of values that the random variable X can assume, we have

$$\sum f(x_i) = \sum P(X = x_i), \quad x_i \in \{0, 1, 2, 3\} = 1$$

as required.

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Since these are the only possible outcomes of the experiment, our probability theorems require that

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Since these are the only possible outcomes of the experiment, our probability theorems require that

$$P(success) = 1 - P(failure)$$

We will denote the probability of success by p.

Our experiment now had two outcomes, "success" and "failure".

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We define a random variable *X* representing the number of successes. Actually, *X* is a function from the sample space *S* into the set $A = \{0, 1, 2, 3\}$, defined by the following table:

$$X: S \mapsto A \quad \text{such that} \begin{cases} \{f, f, f\} & \to & 0\\ \{f, f, s\}, \{f, s, f\}, \{s, f, f\} & \to & 1\\ \{f, s, s\}, \{s, s\}, \{s, s, f\} & \to & 2\\ \{h, h, h\} & \to & 3 \end{cases}$$