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will wait for the function to complete on the remote process, and produce the function's result. The following script creates three new processes,⁴ and instructs process #4 to factor a large number:⁵

```
addprocs(3)
using Primes.factor
@assert nprocs() == 4
future = remotecall(factor, 4, 21883298135690819)
@assert isa(future, Future)
factors = fetch(future)
@assert factors == Dict(234711901=>1,93234719=>1)
```

The **@spawn** macro is often used in place of **remotecall**; it not only takes advantage of a clearer syntax (since macros take unevaluated expressions as arguments), but it chooses a process for you:

```
addprocs(3)
using Primes.factor
ref = @spawn factor(21883298135690819)
factors = fetch(ref)
@assert factors == Dict(234711901=>1,93234719=>1)
```

Between the remote call and fetching the result, the calling process is free to do other work.

A common example of parallel programming across multiple cores computes an approximation of π by generating random points in the square (0, 0)...(1, 1). Counting the number of random points within the inscribed arc (the points (x, y) for which $x^2 + y^2 \leq 1$ in Figure 6.3) divided by the total number of points within the square approximates the value $\frac{\pi}{4}$. Let's generate a million random points in parallel on three cores:



Figure 6.3 Approximation of $\frac{\pi}{4}$

⁴Calling addprocs with an integer argument creates processors on the local machine to take advantage of multiple cores. There are other variants of addprocs to add processors on remote machines either via user-host-port strings or cluster managers; we will not cover these alternatives in this text.

⁵The factor function resides in the package Primes. You must first install this package by executing Pkg.add("Primes") in an interactive shell, or julia -e 'Pkg.add("Primes")' from the command line.