

which is related to the gamma function by

$$B(z, w) = \frac{\Gamma(z)\Gamma(w)}{\Gamma(z+w)} \quad (6.1.9)$$

hence

```
#include <math.h>

float beta(float z, float w)
Returns the value of the beta function B(z, w).
{
    float gammln(float xx);

    return exp(gammln(z)+gammln(w)-gammln(z+w));
}
```

#### CITED REFERENCES AND FURTHER READING:

- Abramowitz, M., and Stegun, I.A. 1964, *Handbook of Mathematical Functions*, Applied Mathematics Series, vol. 55 (Washington: National Bureau of Standards; reprinted 1968 by Dover Publications, New York), Chapter 6.
- Lanczos, C. 1964, *SIAM Journal on Numerical Analysis*, ser. B, vol. 1, pp. 86–96. [1]

## 6.2 Incomplete Gamma Function, Error Function, Chi-Square Probability Function, Cumulative Poisson Function

The incomplete gamma function is defined by

$$P(a, x) \equiv \frac{\gamma(a, x)}{\Gamma(a)} \equiv \frac{1}{\Gamma(a)} \int_0^x e^{-t} t^{a-1} dt \quad (a > 0) \quad (6.2.1)$$

It has the limiting values

$$P(a, 0) = 0 \quad \text{and} \quad P(a, \infty) = 1 \quad (6.2.2)$$

The incomplete gamma function  $P(a, x)$  is monotonic and (for  $a$  greater than one or so) rises from “near-zero” to “near-unity” in a range of  $x$  centered on about  $a - 1$ , and of width about  $\sqrt{a}$  (see Figure 6.2.1).

The complement of  $P(a, x)$  is also confusingly called an incomplete gamma function,

$$Q(a, x) \equiv 1 - P(a, x) \equiv \frac{\Gamma(a, x)}{\Gamma(a)} \equiv \frac{1}{\Gamma(a)} \int_x^\infty e^{-t} t^{a-1} dt \quad (a > 0) \quad (6.2.3)$$

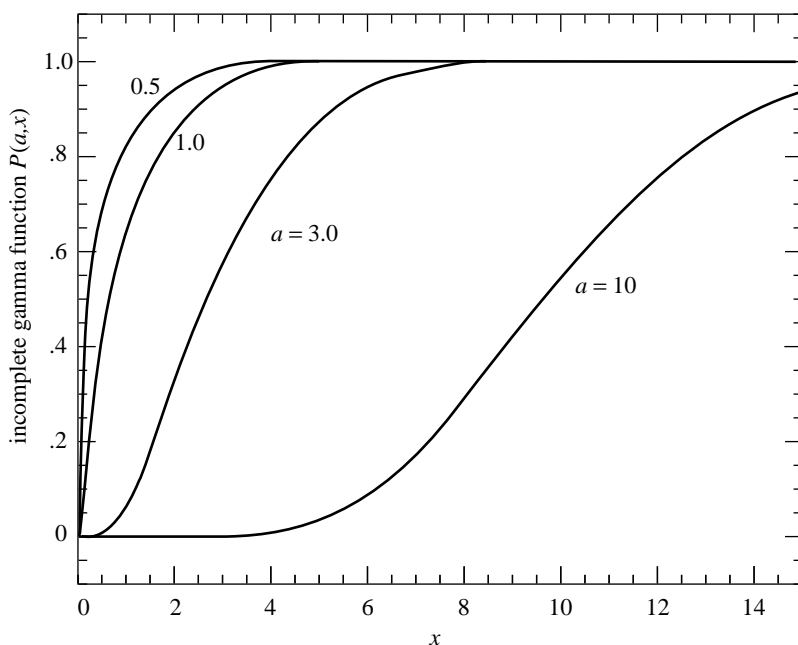


Figure 6.2.1. The incomplete gamma function  $P(a, x)$  for four values of  $a$ .

It has the limiting values

$$Q(a, 0) = 1 \quad \text{and} \quad Q(a, \infty) = 0 \quad (6.2.4)$$

The notations  $P(a, x)$ ,  $\gamma(a, x)$ , and  $\Gamma(a, x)$  are standard; the notation  $Q(a, x)$  is specific to this book.

There is a series development for  $\gamma(a, x)$  as follows:

$$\gamma(a, x) = e^{-x} x^a \sum_{n=0}^{\infty} \frac{\Gamma(a)}{\Gamma(a+1+n)} x^n \quad (6.2.5)$$

One does not actually need to compute a new  $\Gamma(a+1+n)$  for each  $n$ ; one rather uses equation (6.1.3) and the previous coefficient.

A continued fraction development for  $\Gamma(a, x)$  is

$$\Gamma(a, x) = e^{-x} x^a \left( \frac{1}{x+} \frac{1-a}{1+} \frac{1}{x+} \frac{2-a}{1+} \frac{2}{x+} \dots \right) \quad (x > 0) \quad (6.2.6)$$

It is computationally better to use the even part of (6.2.6), which converges twice as fast (see §5.2):

$$\Gamma(a, x) = e^{-x} x^a \left( \frac{1}{x+1-a-} \frac{1 \cdot (1-a)}{x+3-a-} \frac{2 \cdot (2-a)}{x+5-a-} \dots \right) \quad (x > 0) \quad (6.2.7)$$

It turns out that (6.2.5) converges rapidly for  $x$  less than about  $a+1$ , while (6.2.6) or (6.2.7) converges rapidly for  $x$  greater than about  $a+1$ . In these respective

regimes each requires at most a few times  $\sqrt{a}$  terms to converge, and this many only near  $x = a$ , where the incomplete gamma functions are varying most rapidly. Thus (6.2.5) and (6.2.7) together allow evaluation of the function for all positive  $a$  and  $x$ . An extra dividend is that we never need compute a function value near zero by subtracting two nearly equal numbers. The higher-level functions that return  $P(a, x)$  and  $Q(a, x)$  are

```
float gammq(float a, float x)
Returns the incomplete gamma function  $P(a, x)$ .
{
    void gcf(float *gammcf, float a, float x, float *gln);
    void gser(float *gamser, float a, float x, float *gln);
    void nrerror(char error_text[]);
    float gamser, gammcf, gln;

    if (x < 0.0 || a <= 0.0) nrerror("Invalid arguments in routine gammq");
    if (x < (a+1.0)) {
        gser(&gamser, a, x, &gln);
        return gamser;
    } else {
        gcf(&gammcf, a, x, &gln);
        return 1.0-gammcf;
    }
}
```

```
float gammq(float a, float x)
Returns the incomplete gamma function  $Q(a, x) \equiv 1 - P(a, x)$ .
{
    void gcf(float *gammcf, float a, float x, float *gln);
    void gser(float *gamser, float a, float x, float *gln);
    void nrerror(char error_text[]);
    float gamser, gammcf, gln;

    if (x < 0.0 || a <= 0.0) nrerror("Invalid arguments in routine gammq");
    if (x < (a+1.0)) {
        gser(&gamser, a, x, &gln);
        return 1.0-gamser;
    } else {
        gcf(&gammcf, a, x, &gln);
        return gammcf;
    }
}
```

The argument `gln` is set by both the series and continued fraction procedures to the value  $\ln \Gamma(a)$ ; the reason for this is so that it is available to you if you want to modify the above two procedures to give  $\gamma(a, x)$  and  $\Gamma(a, x)$ , in addition to  $P(a, x)$  and  $Q(a, x)$  (cf. equations 6.2.1 and 6.2.3).

The functions `gser` and `gcf` which implement (6.2.5) and (6.2.7) are

```
#include <math.h>
#define ITMAX 100
#define EPS 3.0e-7

void gser(float *gamser, float a, float x, float *gln)
Returns the incomplete gamma function  $P(a, x)$  evaluated by its series representation as gamser.
Also returns  $\ln \Gamma(a)$  as gln.
{
    float gammln(float xx);
```

```

void nrerror(char error_text[]);
int n;
float sum,del,ap;

*gln=gammln(a);
if (x <= 0.0) {
    if (x < 0.0) nrerror("x less than 0 in routine gser");
    *gamser=0.0;
    return;
} else {
    ap=a;
    del=sum=1.0/a;
    for (n=1;n<=ITMAX;n++) {
        ++ap;
        del *= x/ap;
        sum += del;
        if (fabs(del) < fabs(sum)*EPS) {
            *gamser=sum*exp(-x+a*log(x)-(*gln));
            return;
        }
    }
    nrerror("a too large, ITMAX too small in routine gser");
    return;
}
}

#include <math.h>
#define ITMAX 100           Maximum allowed number of iterations.
#define EPS 3.0e-7         Relative accuracy.
#define FPMIN 1.0e-30      Number near the smallest representable
                           floating-point number.

void gcf(float *gammcf, float a, float x, float *gln)
Returns the incomplete gamma function  $Q(a, x)$  evaluated by its continued fraction representation as gammcf. Also returns  $\ln \Gamma(a)$  as gln.
{
    float gammln(float xx);
    void nrerror(char error_text[]);
    int i;
    float an,b,c,d,del,h;

    *gln=gammln(a);
    b=x+1.0-a;
    c=1.0/FPMIN;
    d=1.0/b;
    h=d;
    for (i=1;i<=ITMAX;i++) {
        an = -i*(i-a);
        b += 2.0;
        d=an*d+b;
        if (fabs(d) < FPMIN) d=FPMIN;
        c=b+an/c;
        if (fabs(c) < FPMIN) c=FPMIN;
        d=1.0/d;
        del=d*c;
        h *= del;
        if (fabs(del-1.0) < EPS) break;
    }
    if (i > ITMAX) nrerror("a too large, ITMAX too small in gcf");
    *gammcf=exp(-x+a*log(x)-(*gln))*h;
}

```

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## Error Function

The error function and complementary error function are special cases of the incomplete gamma function, and are obtained moderately efficiently by the above procedures. Their definitions are

$$\operatorname{erf}(x) = \frac{2}{\sqrt{\pi}} \int_0^x e^{-t^2} dt \quad (6.2.8)$$

and

$$\operatorname{erfc}(x) \equiv 1 - \operatorname{erf}(x) = \frac{2}{\sqrt{\pi}} \int_x^\infty e^{-t^2} dt \quad (6.2.9)$$

The functions have the following limiting values and symmetries:

$$\operatorname{erf}(0) = 0 \quad \operatorname{erf}(\infty) = 1 \quad \operatorname{erf}(-x) = -\operatorname{erf}(x) \quad (6.2.10)$$

$$\operatorname{erfc}(0) = 1 \quad \operatorname{erfc}(\infty) = 0 \quad \operatorname{erfc}(-x) = 2 - \operatorname{erfc}(x) \quad (6.2.11)$$

They are related to the incomplete gamma functions by

$$\operatorname{erf}(x) = P\left(\frac{1}{2}, x^2\right) \quad (x \geq 0) \quad (6.2.12)$$

and

$$\operatorname{erfc}(x) = Q\left(\frac{1}{2}, x^2\right) \quad (x \geq 0) \quad (6.2.13)$$

We'll put an extra "f" into our routine names to avoid conflicts with names already in some C libraries:

```
float erf(float x)
Returns the error function erf(x).
{
    float gammp(float a, float x);

    return x < 0.0 ? -gammp(0.5,x*x) : gammp(0.5,x*x);
}

float erffc(float x)
Returns the complementary error function erfc(x).
{
    float gammp(float a, float x);
    float gammq(float a, float x);

    return x < 0.0 ? 1.0+gammp(0.5,x*x) : gammq(0.5,x*x);
}
```

If you care to do so, you can easily remedy the minor inefficiency in `erff` and `erffc`, namely that  $\Gamma(0.5) = \sqrt{\pi}$  is computed unnecessarily when `gammp` or `gammq` is called. Before you do that, however, you might wish to consider the following routine, based on Chebyshev fitting to an inspired guess as to the functional form:

```
#include <math.h>

float erfcc(float x)
Returns the complementary error function erfc(x) with fractional error everywhere less than
1.2 × 10-7.
{
    float t,z,ans;

    z=fabs(x);
    t=1.0/(1.0+0.5*z);
    ans=t*exp(-z*z-1.26551223+t*(1.00002368+t*(0.37409196+t*(0.09678418+
        t*(-0.18628806+t*(0.27886807+t*(-1.13520398+t*(1.48851587+
            t*(-0.82215223+t*0.17087277)))))))));
    return x >= 0.0 ? ans : 2.0-ans;
}
```

There are also some functions of *two* variables that are special cases of the incomplete gamma function:

### Cumulative Poisson Probability Function

$P_x(< k)$ , for positive  $x$  and integer  $k \geq 1$ , denotes the *cumulative Poisson probability* function. It is defined as the probability that the number of Poisson random events occurring will be between 0 and  $k - 1$  *inclusive*, if the expected mean number is  $x$ . It has the limiting values

$$P_x(< 1) = e^{-x} \quad P_x(< \infty) = 1 \quad (6.2.14)$$

Its relation to the incomplete gamma function is simply

$$P_x(< k) = Q(k, x) = \text{gammq}(k, x) \quad (6.2.15)$$

### Chi-Square Probability Function

$P(\chi^2|\nu)$  is defined as the probability that the observed chi-square for a correct model should be less than a value  $\chi^2$ . (We will discuss the use of this function in Chapter 15.) Its complement  $Q(\chi^2|\nu)$  is the probability that the observed chi-square will exceed the value  $\chi^2$  by chance *even* for a correct model. In both cases  $\nu$  is an integer, the number of degrees of freedom. The functions have the limiting values

$$P(0|\nu) = 0 \quad P(\infty|\nu) = 1 \quad (6.2.16)$$

$$Q(0|\nu) = 1 \quad Q(\infty|\nu) = 0 \quad (6.2.17)$$

and the following relation to the incomplete gamma functions,

$$P(\chi^2|\nu) = P\left(\frac{\nu}{2}, \frac{\chi^2}{2}\right) = \text{gammp}\left(\frac{\nu}{2}, \frac{\chi^2}{2}\right) \quad (6.2.18)$$

$$Q(\chi^2|\nu) = Q\left(\frac{\nu}{2}, \frac{\chi^2}{2}\right) = \text{gammq}\left(\frac{\nu}{2}, \frac{\chi^2}{2}\right) \quad (6.2.19)$$

## CITED REFERENCES AND FURTHER READING:

Abramowitz, M., and Stegun, I.A. 1964, *Handbook of Mathematical Functions*, Applied Mathematics Series, vol. 55 (Washington: National Bureau of Standards; reprinted 1968 by Dover Publications, New York), Chapters 6, 7, and 26.

Pearson, K. (ed.) 1951, *Tables of the Incomplete Gamma Function* (Cambridge: Cambridge University Press).

### 6.3 Exponential Integrals

The standard definition of the exponential integral is

$$E_n(x) = \int_1^\infty \frac{e^{-xt}}{t^n} dt, \quad x > 0, \quad n = 0, 1, \dots \quad (6.3.1)$$

The function defined by the principal value of the integral

$$\text{Ei}(x) = - \int_{-x}^\infty \frac{e^{-t}}{t} dt = \int_{-\infty}^x \frac{e^t}{t} dt, \quad x > 0 \quad (6.3.2)$$

is also called an exponential integral. Note that  $\text{Ei}(-x)$  is related to  $-E_1(x)$  by analytic continuation.

The function  $E_n(x)$  is a special case of the incomplete gamma function

$$E_n(x) = x^{n-1} \Gamma(1-n, x) \quad (6.3.3)$$

We can therefore use a similar strategy for evaluating it. The continued fraction — just equation (6.2.6) rewritten — converges for all  $x > 0$ :

$$E_n(x) = e^{-x} \left( \frac{1}{x+} \frac{n}{1+} \frac{1}{x+} \frac{n+1}{1+} \frac{2}{x+} \dots \right) \quad (6.3.4)$$

We use it in its more rapidly converging even form,

$$E_n(x) = e^{-x} \left( \frac{1}{x+n-} \frac{1 \cdot n}{x+n+2-} \frac{2(n+1)}{x+n+4-} \dots \right) \quad (6.3.5)$$

The continued fraction only really converges fast enough to be useful for  $x \gtrsim 1$ . For  $0 < x \lesssim 1$ , we can use the series representation

$$E_n(x) = \frac{(-x)^{n-1}}{(n-1)!} [-\ln x + \psi(n)] - \sum_{\substack{m=0 \\ m \neq n-1}}^{\infty} \frac{(-x)^m}{(m-n+1)m!} \quad (6.3.6)$$

The quantity  $\psi(n)$  here is the digamma function, given for integer arguments by

$$\psi(1) = -\gamma, \quad \psi(n) = -\gamma + \sum_{m=1}^{n-1} \frac{1}{m} \quad (6.3.7)$$