

University of Southern California

EE511 Simulation Methods for Stochastic Systems

Project #6- Continuous sampling

BY

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PROBLEM 1:

BOX-MULLER TRANSFORM

- A pseudo-random number sampling method for generating pairs of independent, standard, normally distributed (zero expectation, unit variance) random numbers, given a source of uniformly distributed random numbers.
- It is expressed in two different forms
- One default form from Box and Muller, maps two samples generated from the uniform distribution on the interval $[0,1]$ to two standard normally distributed samples.
- Whereas the other, Polar form maps two samples taken from distinct interval $[-1, +1]$ to two normally distributed samples without applying sine or cosine functions.

Let U_1 and U_2 be random variables that are independent and are uniformly distributed in interval $(0,1)$.

$$Z_0 = R \cos(\Theta) = \sqrt{-2 \ln U_1} \cos(2\pi U_2)$$

$$Z_1 = R \sin(\Theta) = \sqrt{-2 \ln U_1} \sin(2\pi U_2).$$

Here in the above, Z_0 and Z_1 be random variables which are independent and follow a standard normal distribution.

POLAR MARSAGLIA METHOD

- A pseudo-random number sampling method for generating a pair of independent standard normal random variables.
- It is superior to the Box-Muller transform.
- This method works by choosing random points (x, y) in the square $-1 < x < 1, -1 < y < 1$ until

$$s = x^2 + y^2 < 1.$$

and then returning the required pair of normal random variables as.

$$x\sqrt{\frac{-2\ln(s)}{s}}, y\sqrt{\frac{-2\ln(s)}{s}},$$

Or equivalently,

$$\frac{x}{\sqrt{s}}\sqrt{-2\ln(s)}, \frac{y}{\sqrt{s}}\sqrt{-2\ln(s)},$$

where x/\sqrt{s} and y/\sqrt{s} represent the cosine and sine of the angle that the vector (x, y) makes with x axis.

ALGORITHM:

- Here U₁, U₂ are sampled from uniform distribution (-1,1)
- Here check for the condition, if $s=(U_1^2+U_2^2)$
- If the condition is not satisfied, we will be taking new U₁, U₂.
- Then value of X and Y is computed.
- $X = \sqrt{-2 \log(s)/s} * U_1$
- $Y = \sqrt{-2 \log(s)/s} * U_2$

CODE:

- Independent random variables are generated using Box-Muller method. BoxMuller Random Samples function is used to perform the required operation according to the formulas described above.
- Probability distribution is made using makedist function.
- By utilizing Matlab's inbuilt 'cov' function, covariance is calculated.
- Generation of Histogram is done, and the theoretical pdf is overlayed on the histogram as can be seen in the graph.
- Here Matlab's 'tic' and 'toc' commands computational time taken by above method like Box Muller is calculated for a million samples.
- Similarly, we calculate the computational time for Polar Marsaglia method and compare the results. We use PolarMarsaglia_RandomSamples function for the calculation.

CODE:

```
clc;
close all;
clear all;

%Box Muller
[M,N,A]=Box_Muller_method(1000);
covariance1=cov(M,N);
T = -100:100;
mu=3;
sigma=sqrt(13);
prob_dist=makedist('Normal',mu,sigma);%Make probability distribution
P=pdf(prob_dist,T);
%start timer
tic;
Box_Muller_method(1000000);
%stop timer
ET1=toc;

%Output
disp('Output for the Box-Muller Method:');
fprintf('\nThe Covariance of X and Y is: \n%f %f\n%f %f\n',covariance1);
yyaxis left
hist(A);
title('Histogram of Random Variable A');
xlabel('Value');
ylabel('Frequency');
yyaxis right
plot(T,P);
ylabel('Probability');
mx=mean(M);
my=mean(N);
vx=var(M);
vy=var(N);
fprintf('\nThe Observed mean and variance of X are:\nMean = %f\nVariance = %f\n',mx,vx);
fprintf('\nThe Observed mean and variance of Y are:\nMean = %f\nVariance = %f\n',my,vy);
%start time
```

```

%start time
tic;
[M,N]=Polarmarsaglia(1000000);
%end time
ET2=toc;

meanx1=mean(M);
meany1=mean(N);
varx1=var(M);
vary1=var(N);
cov2=cov(M,N);
disp('Polar Marsaglia method output:');
fprintf('\nObserved mean and variance of M are:\nMean = %f\nVariance = %f\n',meanx1,varx1);
fprintf('\nObserved mean and variance of N are:\nMean = %f\nVariance = %f\n',meany1,vary1);
fprintf('\nThe Covariance of M and N is: \n%f %f\n%f %f\n',cov2);
disp('Time required to generate 1,000,000 pairs of independent samples\n');
fprintf('Utilizing the Polar Marsaglia method :\n%f\n',ET2);
fprintf('Utilizing the Box Muller method:\n%f\n',ET1);

```

```

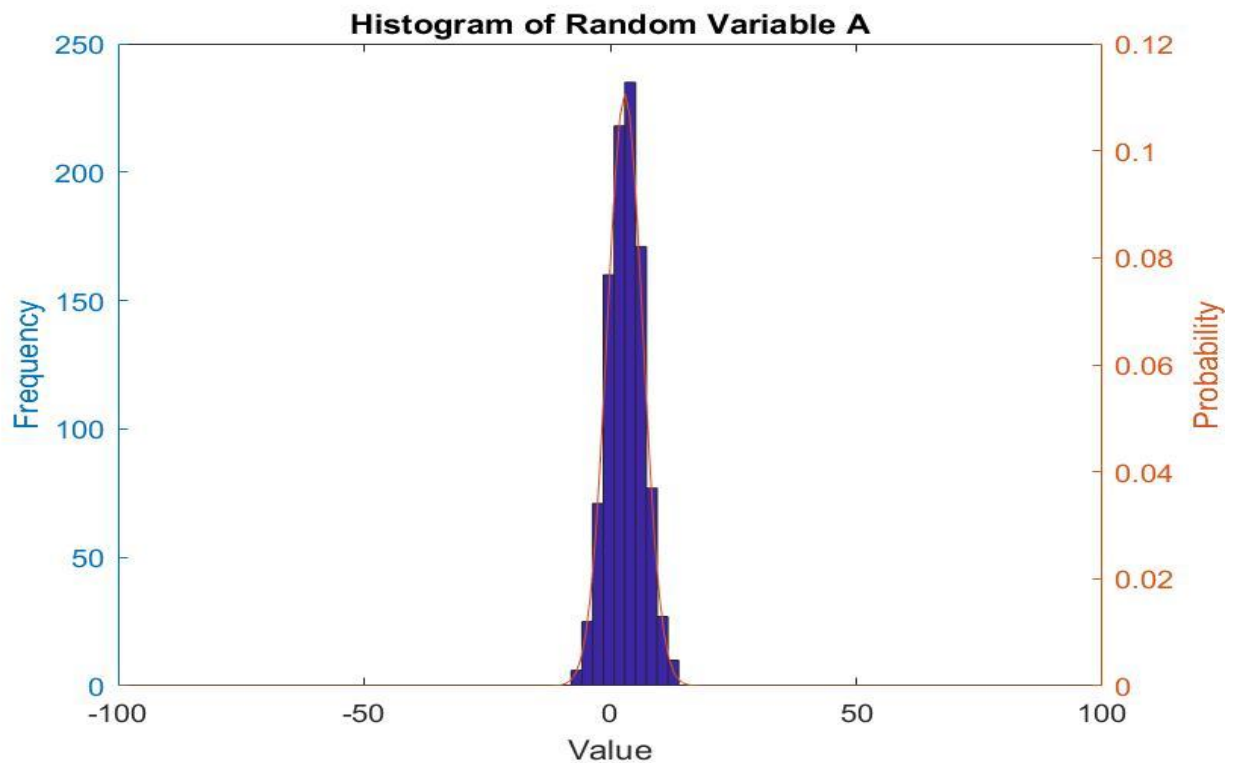
function [M,N,O] = Box_Muller_method(number)
% m-mean
% v-variance
mean2=2;
var2=9;
mean1=1;
var1=4;
for j=1:number
    %uniform random number generation
    rand2=rand();
    rand1=rand();
    x_var(j)=sqrt(-2*log(rand1))*cos(2*pi*rand2);
    y_var(j)=sqrt(-2*log(rand1))*sin(2*pi*rand2);
    M(j)=sqrt(var1)*x_var(j)+mean1;
    N(j)=sqrt(var2)*y_var(j)+mean2;
    O(j)=M(j)+N(j);
end
end

```

```

function[M,N] = Polarmarsaglia(n)
% m-mean
% v-variance
mean3 = 1;
var3 = 4;
mean4 = 2;
var4 = 9;
% algorithm generated random number
j = 0;
% Independent random variable generation
while(j<=n-1)
    rand1 = 2*rand()-1;
    rand2 = 2*rand()-1;
    sum = rand1^2 + rand2^2;
    if(sum < 1)
        j = j + 1;
        x1_var(j) = sqrt(-2*log(sum)/sum)*rand1;
        y1_var(j) = sqrt(-2*log(sum)/sum)*rand2;
    end
end
% Scaling
M = sqrt(var3)*x1_var + mean3;
N = sqrt(var4)*y1_var + mean4;
end

```



Output for the Box-Muller Method:

The Covariance of X and Y is:

3.999384 -0.034594

-0.034594 9.258359

The Observed mean and variance of X are:

Mean = 1.073860

Variance = 3.999384

The Observed mean and variance of Y are:

Mean = 1.868625

Variance = 9.258359

Polar Marsaglia method output:

Observed mean and variance of M are:

Mean = 0.996184

Variance = 4.004706

Observed mean and variance of N are:

Mean = 1.998835

Variance = 8.988067

The Covariance of M and N is:

4.004706 0.002958

0.002958 8.988067

Time required to generate 1,000,000 pairs of independent samples\n

Utilizing the Polar Marsaglia method:

0.492559

Utilizing the Box Muller method:

1.012355

Time required to generate 1,000,000 pairs of independent samples\n

Utilizing the Polar Marsaglia method :

0.461763

Utilizing the Box Muller method:

0.818044

ANALYSIS:

- Polar Marsaglia method takes lesser computational time when compared to Box Muller method. This is evident from the mentioned two trials.
- From the above results it is evident that Polar Marsaglia is superior when compared to Box Muller method.

PROBLEM 2:

GAMMA RANDOM VARIABLE GENERATION

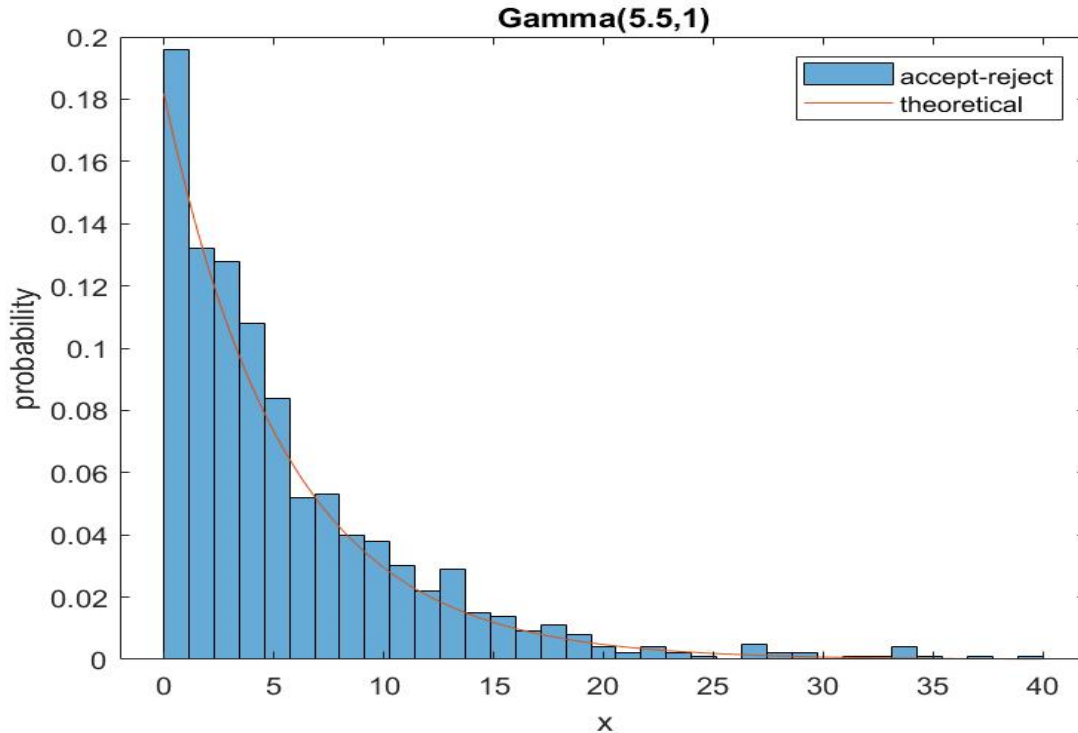
Accept-Reject Method

- Sampling values are generated using rejection sampling from a target X with arbitrary probability density function $f(x)$ by using a proposal distribution Y with probability density $g(x)$.
- The idea is that one can generate a sample value from X by instead sampling from Y and accepting the sample from Y with probability $f(x)/(Mg(x))$, repeating the draws from Y until a value is accepted.
- First, we derive the constant M , M need to be make sure that $Mg(x) > f(x)$ should be true for every x , therefore, we take the max of $(f(x)/g(x))$ to determine the M . Then we using the accept-reject method to derive $f(x)$

ALGORITHM

- Set random variable S which has distribution $g(x)$.
- Compute the constant M .
- Generate a sample s from S , generate a number u from std $U [0,1]$.
- if $u < p(s)/Mg(s)$, accept and records, else reject.
- Repeat 3,4 for trail budget times.

The exponential distribution with parameter 0.1 is used as my $g(x)$, because it is easy to use inverse sampling method to generate a sample from exponential distribution.



Generation of Gamma Accept-Reject Method

```

clc;
clear all;
close all;
trial_count = 0;
X = 2;
A = 0;

while A < 1000
    %Sample generation
    y = generateExpDis(0.1);
    if rand(1)<exponential5(y)/(X*0.1*exp(-0.1 * y))
        A = A + 1;
        sample(A) = y;
    end
    trial_count = trial_count + 1;
end

efficiency = 1000/trial_count;
histogram(sample',35,'BinLimits',[0,40],'Normalization','probability');
hold on
t = 0: 0.1: 40;

title('Gamma(5.5,1)');
plot(t,exp5_5(t));
legend('accept-reject','theoretical');
ylabel('probability');
xlabel('x');

```

ANALYSIS:

- From the above results we could see that a total of 2056 trials are taken to generate 1000 samples
- the acceptance rate is nearly every M sampling we get one sample.

PROBLEM 3:

- We call X as random variable when its characteristic function can be written as

$$\phi(\omega; \alpha, \beta, c, \mu) = \exp(i\omega\mu - |c\omega|^\alpha(1 - i\beta \operatorname{sgn}(\omega)\Phi))$$

$$\Phi = \{\tan(\pi\alpha/2), \alpha \neq 1, (-2/\pi)\log|\omega|, \alpha = 1\}$$

$\mu \in \mathbb{R}$ is a shift parameter, β belongs to range $[-1,1]$ is a skewness parameter which measures asymmetry

- In this context the usual skewness is not well defined, as for $\alpha < 2$ the distribution does not admit 2nd or higher moments, and the usual skewness definition is the 3rd central moment.

The following **program** is run for a single experiment.

```

function experiment(alpha,beta,sampleAmount)
figure

X = stblrnd(alpha,beta,1,o,sampleAmount,1);
histogram(X,'BinLimits',[-20,20],'Normalization','probability');
t = -20:0.2:20;
hold on
y = stblpdf(t,alpha,beta,1,o);
plot(t,y)
title(strcat('\alpha =', num2str(alpha) , '\beta =', num2str(beta)));
legend('histogram','theoretical')
hold off
%time series
figure
plot(X)
ylabel('sample value')
title(strcat('time series \alpha =', num2str(alpha) , '\beta =', num2str(beta)))
end

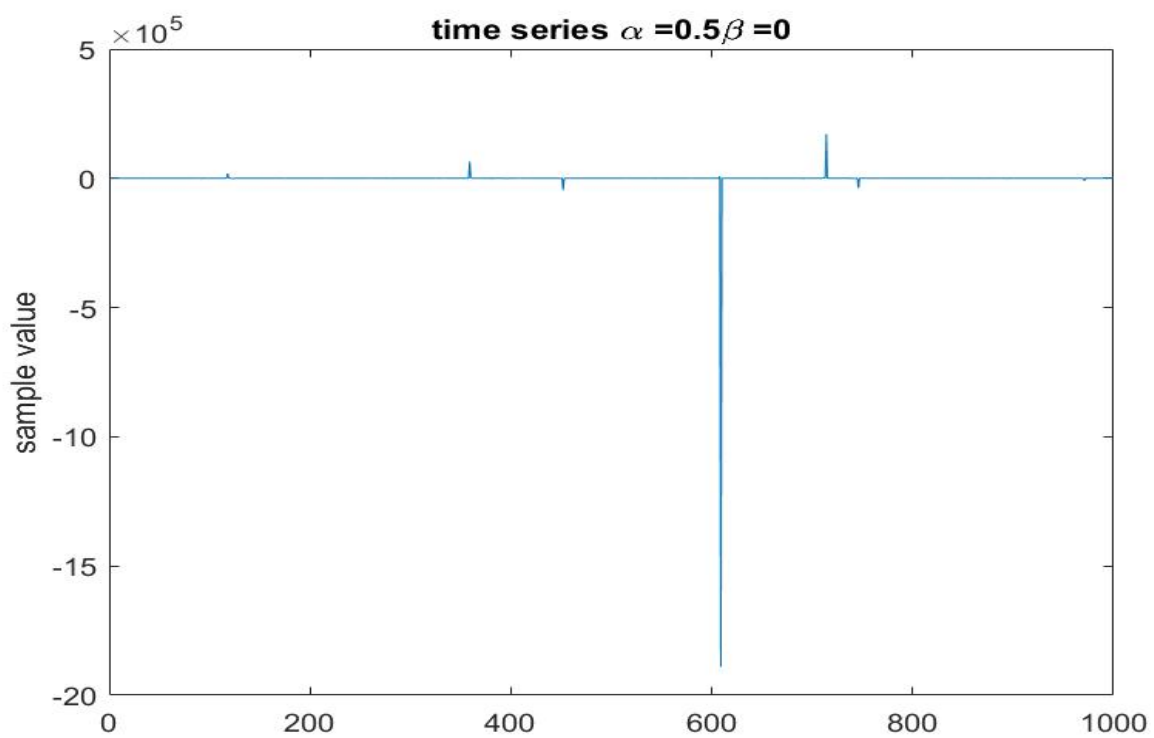
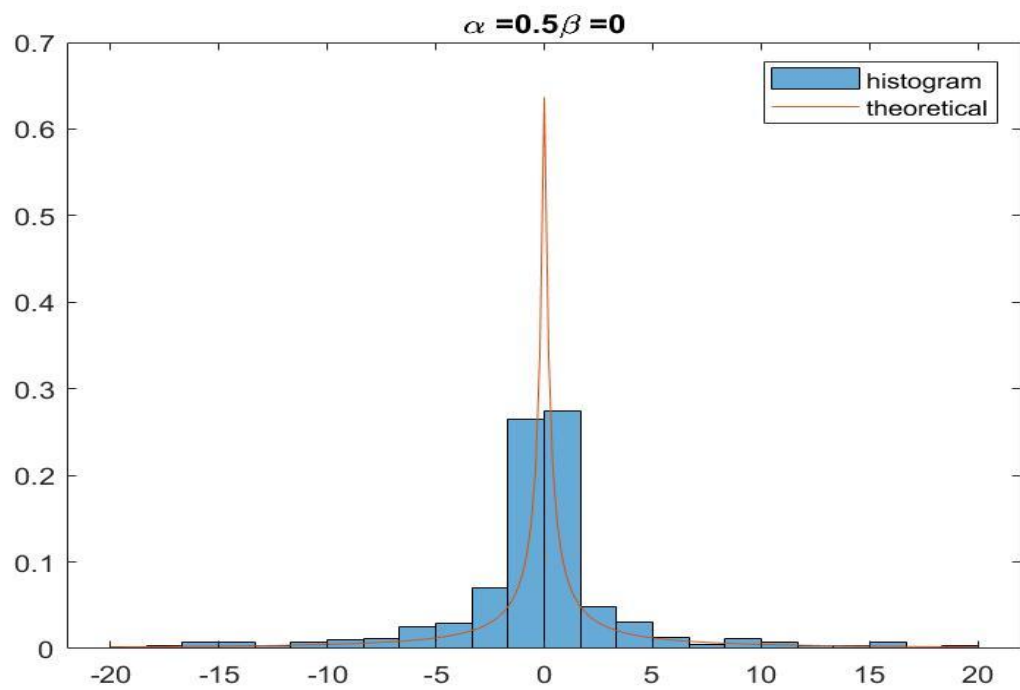
```

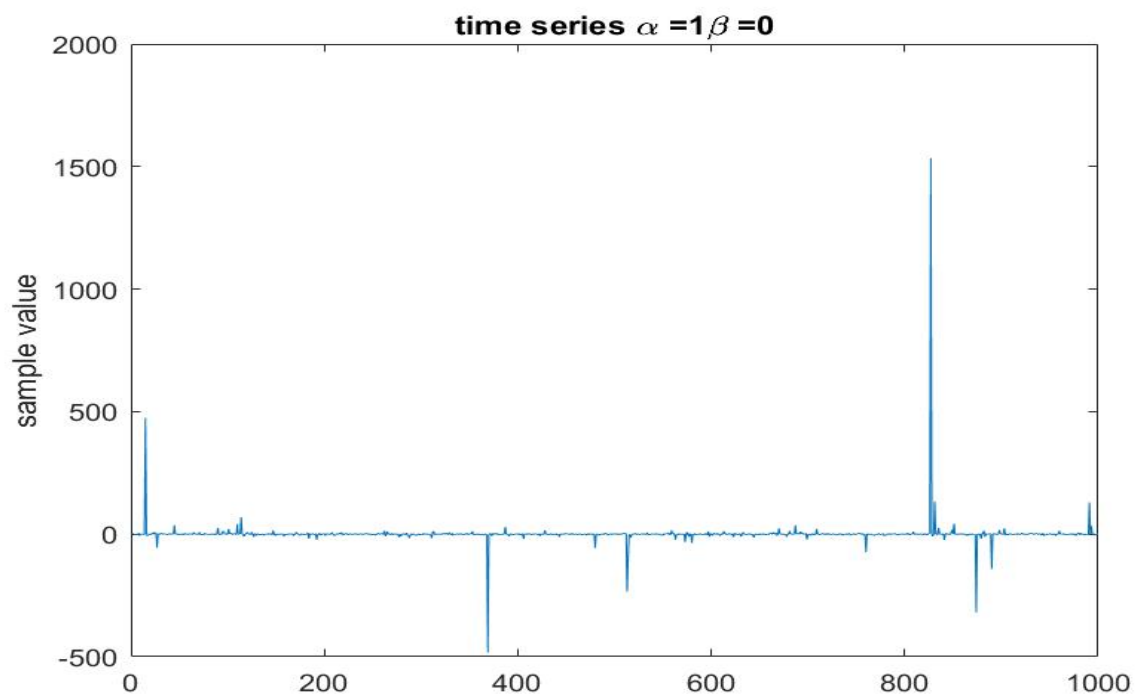
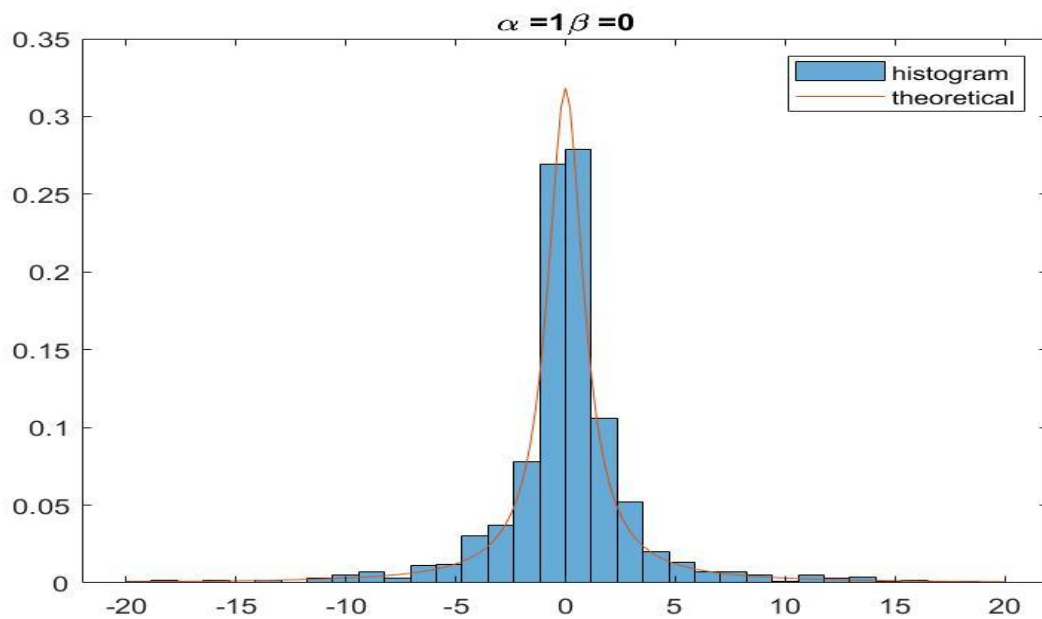
Program:

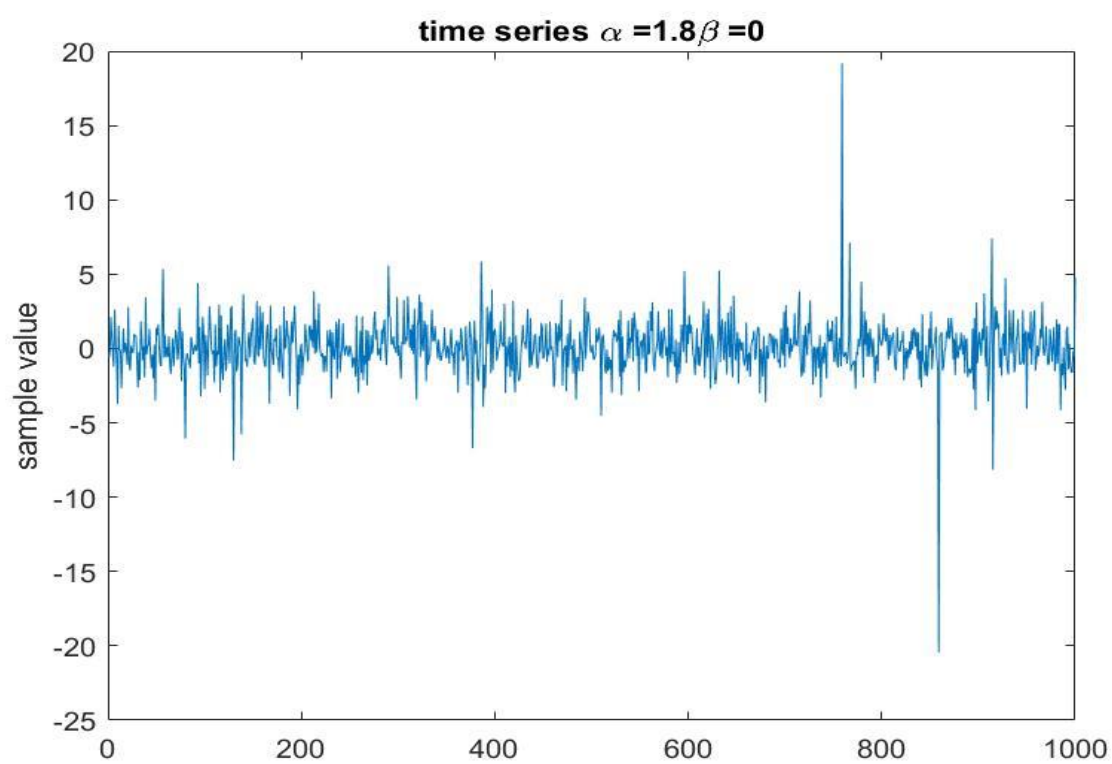
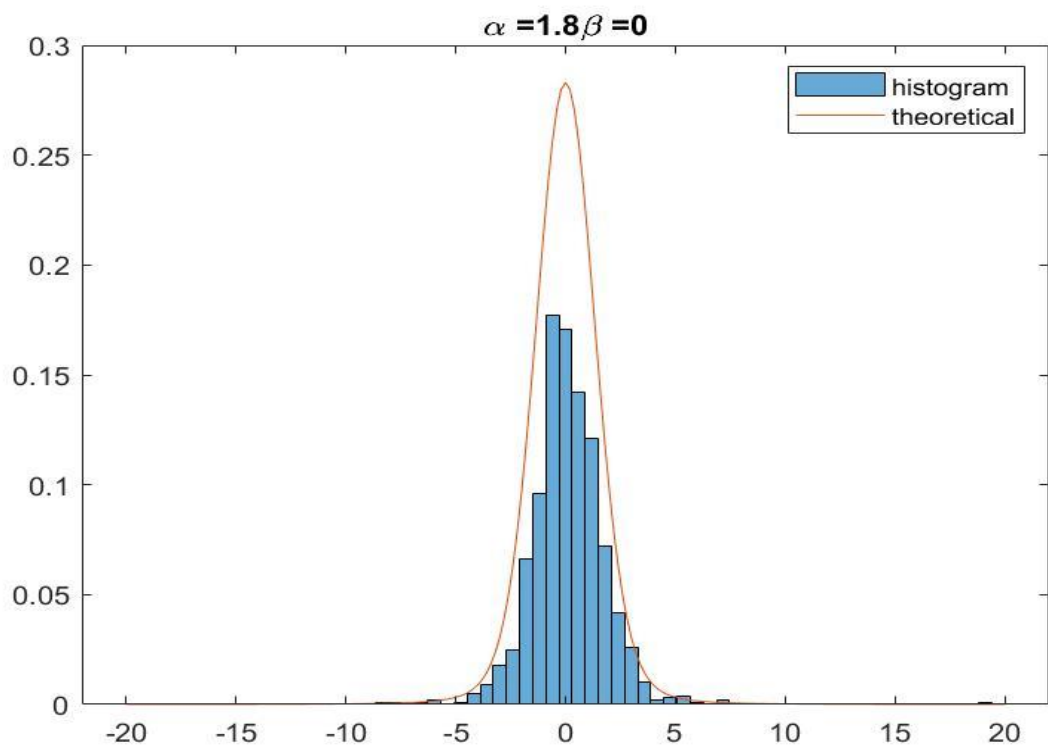
```

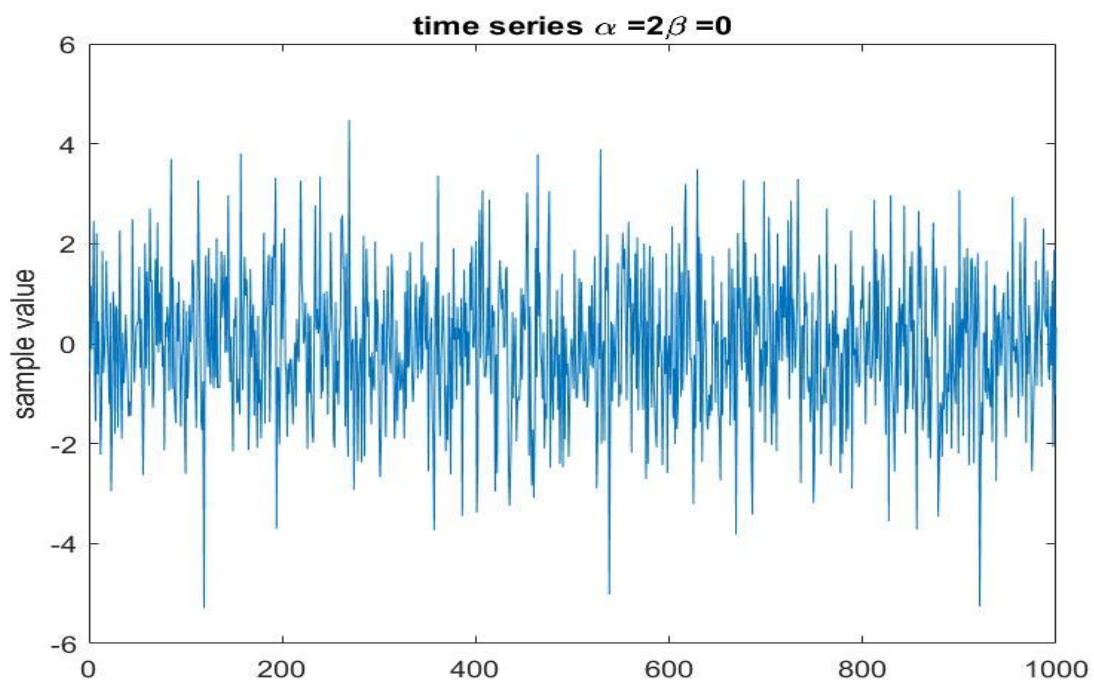
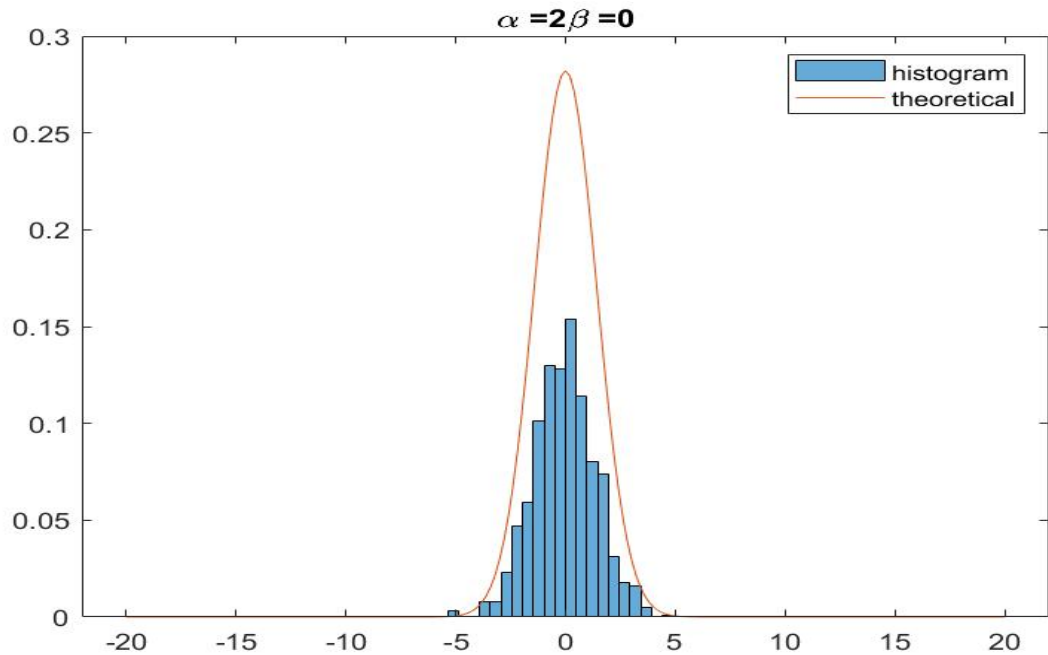
experiment(0.5,0,1000);
experiment(1,0,1000);
experiment(1.8,0,1000);
experiment(2.0,0,1000);
experiment(0.5,0.75,1000);
experiment(1.0,0.75,1000);
experiment(1.8,0.75,1000);
experiment(2.0,0.75,1000);

```



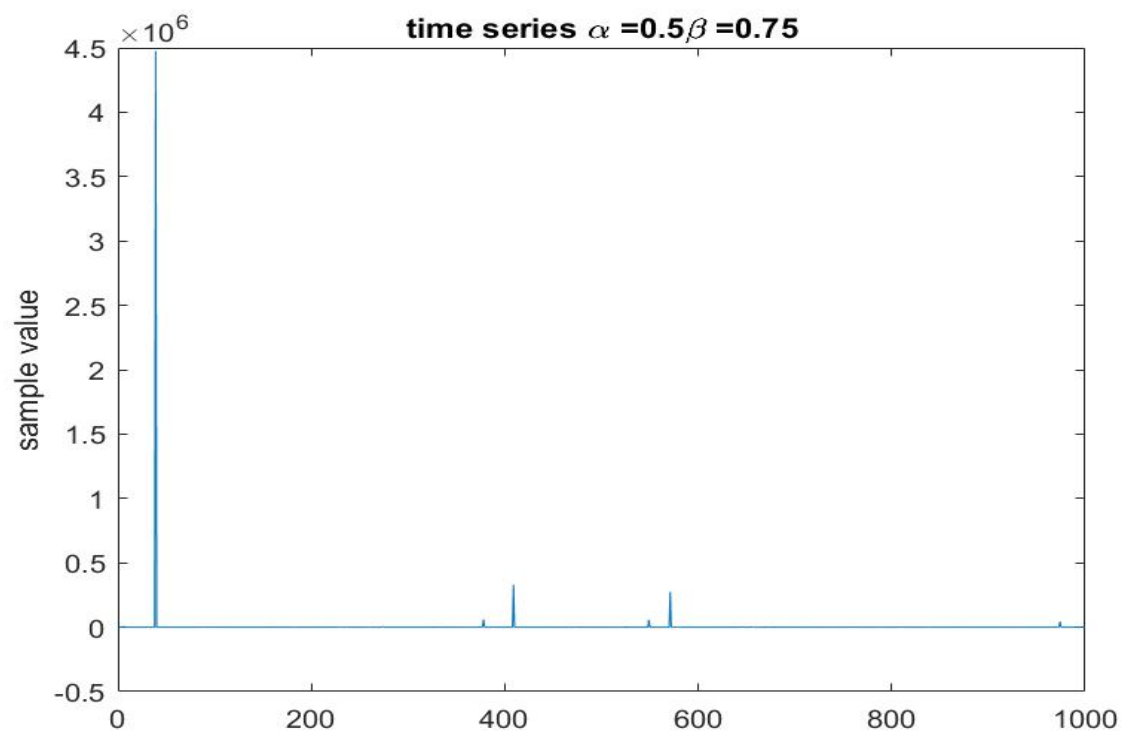
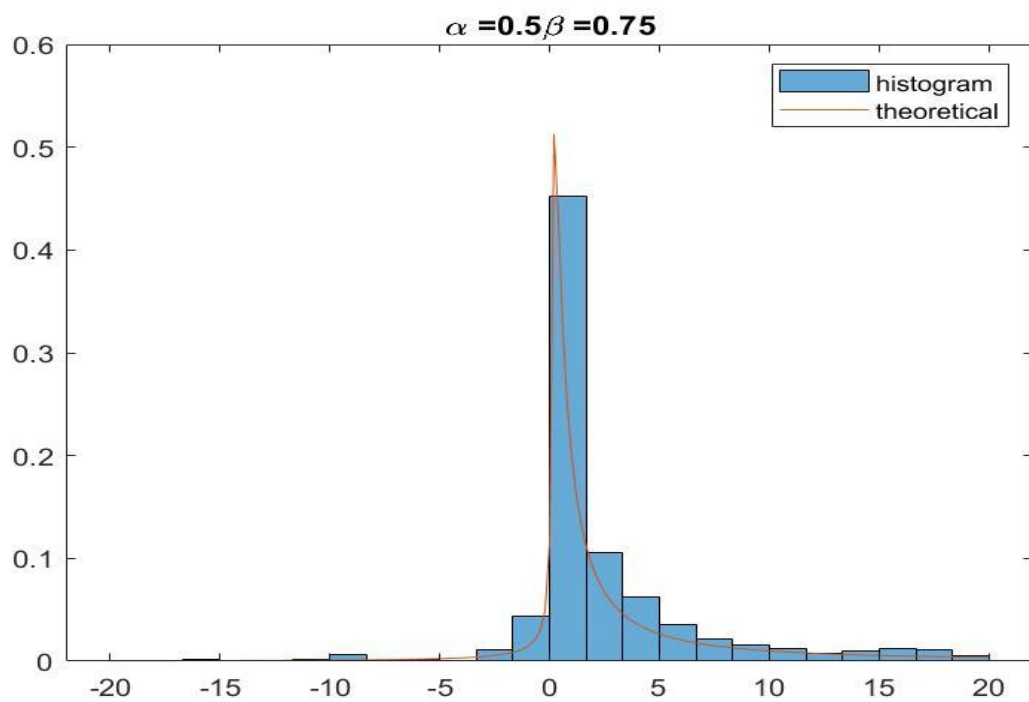


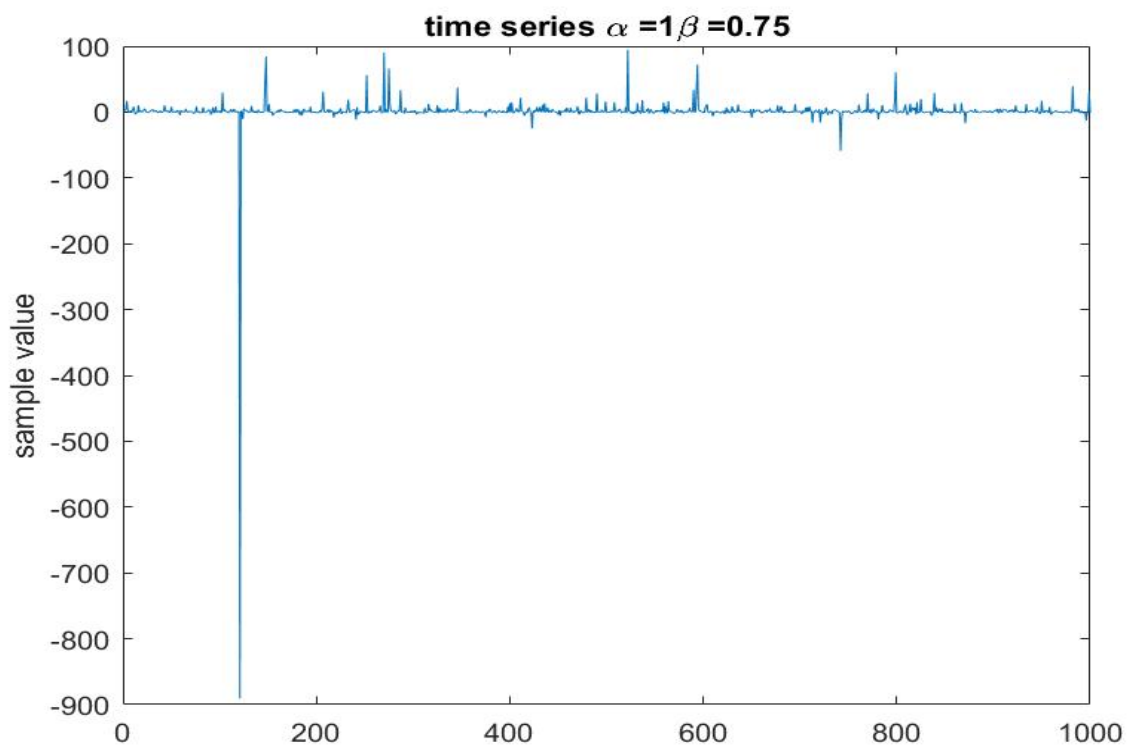
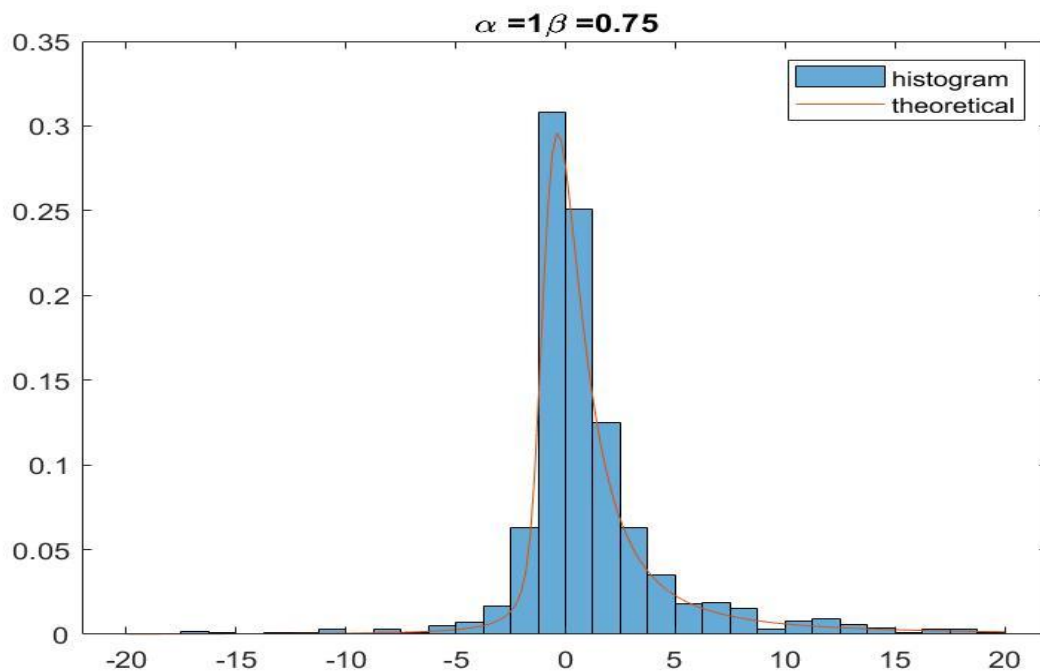


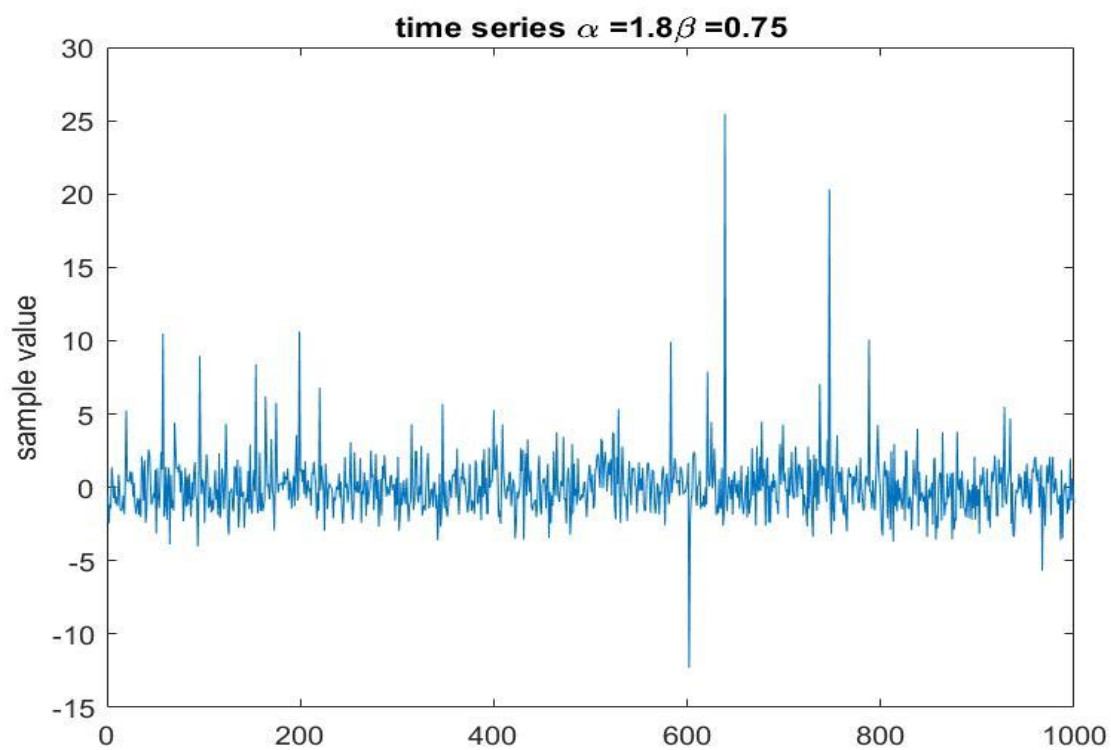
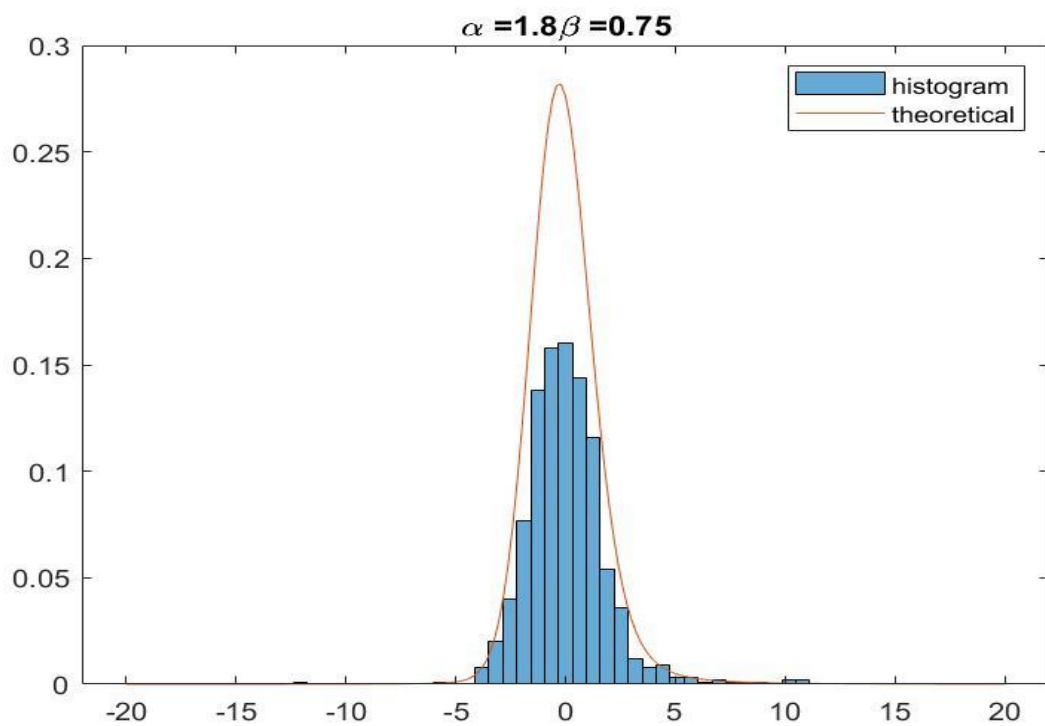


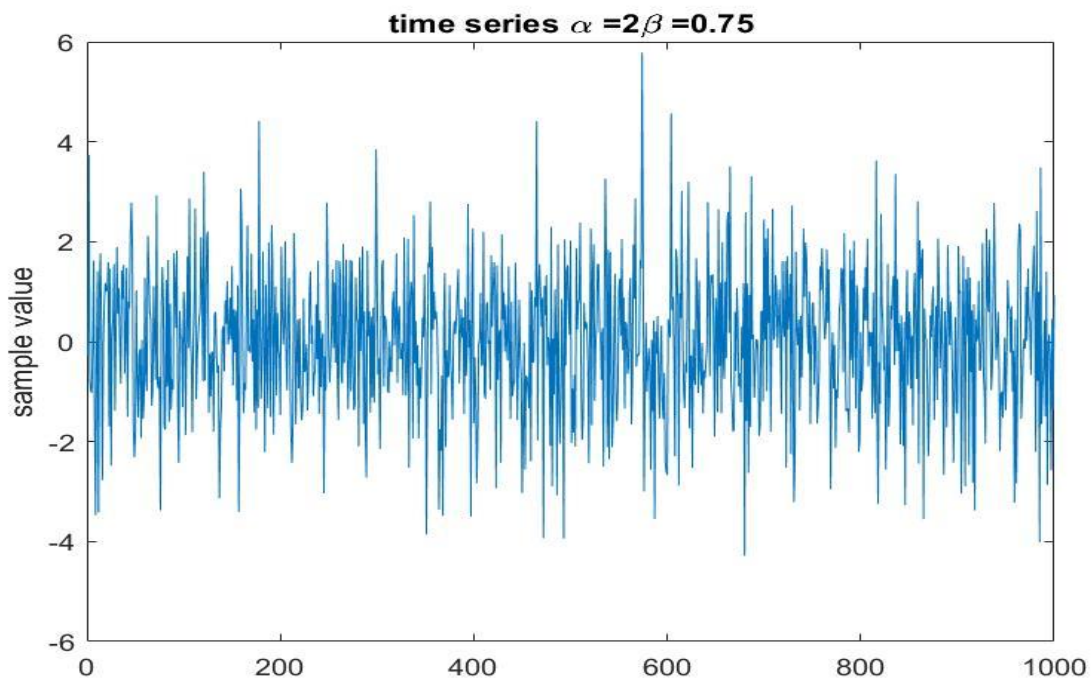
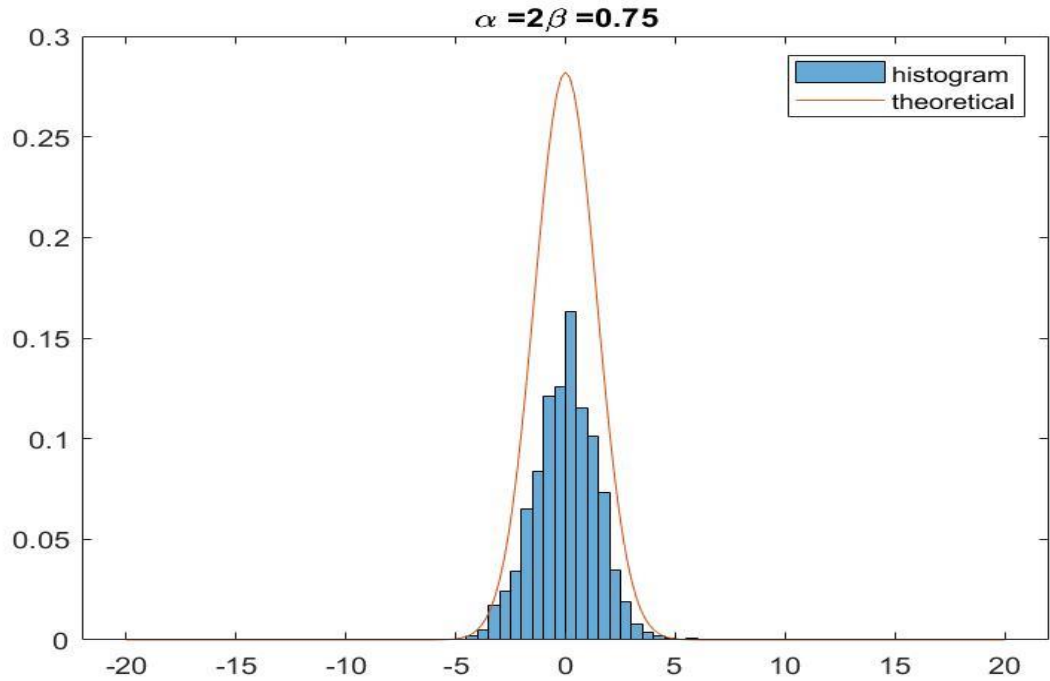
From the graphs we could see that, with the decrease in value of alpha, a less oscillating and the magnitude of the sample is large,

When the value of alpha is 2, a gaussian distribution and the white noise is simulated by samples.









From the above figures, we could observe that when the value of α is 0.5 and α is 1 there is a thick tail because of skewness. We could see that, with the decrease in α , we got less oscillating, and the sample magnitude can be really large. Moreover, because of the change of skewness, it seems that the oscillation starts earlier.

Sample generation for Alpha-Stable

```
function r = stblrnd(alpha,beta,gamma,delta,varargin)
%STBLRND alpha-stable random number generator.
% R = STBLRND(ALPHA,BETA,GAMMA,DELTA) draws a sample from the Levy
% alpha-stable distribution with characteristic exponent ALPHA,
% skewness BETA, scale parameter GAMMA and location parameter DELTA.
% ALPHA,BETA,GAMMA and DELTA must be scalars which fall in the following
% ranges :
%   0 < ALPHA <= 2
%  -1 <= BETA <= 1
%   0 < GAMMA < inf
%  -inf < DELTA < inf
%
%
% R = STBLRND(ALPHA,BETA,GAMMA,DELTA,M,N,...) or
% R = STBLRND(ALPHA,BETA,GAMMA,DELTA,[M,N,...]) returns an M-by-N-by-...
% array.
%
%
% References:
% [1] J.M. Chambers, C.L. Mallows and B.W. Stuck (1976)
%   "A Method for Simulating Stable Random Variables"
%   JASA, Vol. 71, No. 354. pages 340-344
%
% [2] Aleksander Weron and Rafal Weron (1995)
%   "Computer Simulation of Levy alpha-Stable Variables and Processes"
%   Lec. Notes in Physics, 457, pages 379-392
%

if nargin < 4
    error('stats:stblrnd:TooFewInputs','Requires at least four input arguments.');
```

```
end

% Check parameters
if alpha <= 0 || alpha > 2 || ~isscalar(alpha)
    error('stats:stblrnd:BadInputs','"alpha" must be a scalar which lies in the interval (0,2]');
```

```
end
if abs(beta) > 1 || ~isscalar(beta)
    error('stats:stblrnd:BadInputs','"beta" must be a scalar which lies in the interval [-1,1]');
```

```
end
if gamma < 0 || ~isscalar(gamma)
    error('stats:stblrnd:BadInputs','"gamma" must be a non-negative scalar');
```

```
end
if ~isscalar(delta)
    error('stats:stblrnd:BadInputs','"delta" must be a scalar');
```

```
end

% Get output size
[err, sizeOut] = genOutsize(4,alpha,beta,gamma,delta,varargin{:});
if err > 0
    error('stats:stblrnd:InputSizeMismatch','Size information is inconsistent.');
```

```
end
```

```
%---Generate sample---
```

```
% See if parameters reduce to a special case, if so be quick, if not  
% perform general algorithm
```

```
if alpha == 2 % Gaussian distribution
```

```
    r = sqrt(2) * randn(sizeOut);
```

```
elseif alpha==1 && beta == 0 % Cauchy distribution
```

```
    r = tan( pi/2 * (2*rand(sizeOut) - 1) );
```

```
elseif alpha == .5 && abs(beta) == 1 % Levy distribution (a.k.a. Pearson V)
```

```
    r = beta ./ randn(sizeOut).^2;
```

```
elseif beta == 0 % Symmetric alpha-stable
```

```
    V = pi/2 * (2*rand(sizeOut) - 1);
```

```
    W = -log(rand(sizeOut));
```

```
    r = sin(alpha * V) ./ ( cos(V).^(1/alpha) ) .* ...  
        ( cos( V.*(1-alpha) ) ./ W ).^((1-alpha)/alpha);
```

```
elseif alpha ~= 1 % General case, alpha not 1
```

```
    V = pi/2 * (2*rand(sizeOut) - 1);
```

```
    W = - log( rand(sizeOut) );
```

```
    const = beta * tan(pi*alpha/2);
```

```
    B = atan( const );
```

```
    S = (1 + const * const).^(1/(2*alpha));
```

```
    r = S * sin( alpha*V + B ) ./ ( cos(V) ).^(1/alpha) .* ...  
        ( cos( (1-alpha) * V - B ) ./ W ).^((1-alpha)/alpha);
```

```
else % General case, alpha = 1
```

```
    V = pi/2 * (2*rand(sizeOut) - 1);
```

```
    W = - log( rand(sizeOut) );
```

```
    piover2 = pi/2;
```

```
    sclshftV = piover2 + beta * V;
```

```
    r = 1/piover2 * ( sclshftV .* tan(V) - ...  
        beta * log( (piover2 * W .* cos(V) ) ./ sclshftV ) );
```

```
end
```

```
% Scale and shift
```

```
if alpha ~= 1
```

```
    r = gamma * r + delta;
```

```
else
```

```
    r = gamma * r + (2/pi) * beta * gamma * log(gamma) + delta;
```

```
end
```

```
end
```

PDF Generation for alpha stable

```
function p = stblpdf(x,alpha,beta,gam,delta,varargin)
%P = STBLPDF(X,ALPHA,BETA,GAM,DELTA) returns the pdf of the stable
% distribution with characteristic exponent ALPHA, skewness BETA, scale
% parameter GAM, and location parameter DELTA, at the values in X. We use
% the parameterization of stable distributions used in [2] - The
% characteristic function phi(t) of a S(ALPHA,BETA,GAM,DELTA)
% random variable has the form
%
% 
$$\phi(t) = \exp(-GAM^ALPHA |t|^ALPHA [1 - i BETA (\tan(\pi ALPHA/2) \text{sign}(t)) \\ + i DELTA t]) \text{ if } \alpha \neq 1$$

%
% 
$$\phi(t) = \exp(-GAM |t| [1 + i BETA (2/\pi) (\text{sign}(t)) \log|t|] + i DELTA t) \\ \text{ if } \alpha = 1$$

%
% The size of P is the size of X. ALPHA,BETA,GAM and DELTA must be scalars
%
%P = STBLPDF(X,ALPHA,BETA,GAM,DELTA,TOL) computes the pdf to within an
% absolute error of TOL.
%
% The algorithm works by computing the numerical integrals in Theorem
% 1 in [1] using MATLAB's QUADV function. The integrands
% are smooth non-negative functions, but for certain parameter values
% can have sharp peaks which might be missed. To avoid this, STBLPDF
% locates the maximum of this integrand and breaks the integral into two
% pieces centered around this maximum (this is exactly the idea suggested
% in [1]).
%
% If abs(ALPHA - 1) < 1e-5, ALPHA is rounded to 1.
%
%P = STBLPDF(...,'quick') skips the step of locating the peak in the
% integrand, and thus is faster, but is less accurate deep into the tails
% of the pdf. This option is useful for plotting. In place of 'quick',
% STBLPDF also accepts a logical true or false (for quick or not quick)
%
% See also: STBLRND, STBLCDF, STBLINV, STBLFIT
%
% References:
%
% [1] J. P. Nolan (1997)
% "Numerical Calculation of Stable Densities and Distribution
% Functions" Commun. Statist. - Stochastic Models, 13(4), 759-774
%
% [2] G Samorodnitsky, MS Taqqu (1994)
% "Stable non-Gaussian random processes: stochastic models with
% infinite variance" CRC Press
%

if nargin < 5
    error('stblpdf:TooFewInputs','Requires at least five input arguments.');
```

end

% Check parameters

```
if alpha <= 0 || alpha > 2 || ~isscalar(alpha)
```

```

    error('stblpdf:BadInputs', ' "alpha" must be a scalar which lies in the interval (0,2]');
end
if abs(beta) > 1 || ~isscalar(beta)
    error('stblpdf:BadInputs', ' "beta" must be a scalar which lies in the interval [-1,1]');
end
if gam < 0 || ~isscalar(gam)
    error('stblpdf:BadInputs', ' "gam" must be a non-negative scalar');
end
if ~isscalar(delta)
    error('stblpdf:BadInputs', ' "delta" must be a scalar');
end

```

```

% Warn if alpha is very close to 1 or 0
if ( 1e-5 < abs(1 - alpha) && abs(1 - alpha) < .02) || alpha < .02
    warning('stblpdf:ScaryAlpha',...
        'Difficult to approximate pdf for alpha close to 0 or 1')
end

```

```

% warnings will happen during call to QUADV, and it's okay
warning('off');

```

```

% Check and initialize additional inputs

```

```

quick = false;
tol = [];
for i=1:length(varargin)
    if strcmp(varargin{i}, 'quick')
        quick = true;
    elseif islogical(varargin{i})
        quick = varargin{end};
    elseif isscalar(varargin{i})
        tol = varargin{i};
    end
end
end

```

```

if isempty(tol)
    if quick
        tol = 1e-8;
    else
        tol = 1e-12;
    end
end
end

```

```

%===== Compute pdf =====%

```

```

% Check to see if you are in a simple case, if so be quick, if not do

```

```

% general algorithm

```

```

if alpha == 2          % Gaussian distribution
    x = (x - delta)/gam;    % Standardize
    p = 1/sqrt(4*pi) * exp( -.25 * x.^2 ); % ~ N(0,2)
    p = p/gam; %rescale

```

```

elseif alpha==1 && beta == 0 % Cauchy distribution

```



```

x = (x - delta)/gam;          % Standardize
p = (1/pi) * 1./(1 + x.^2);
p = p/gam; %rescale

elseif alpha == .5 && abs(beta) == 1 % Levy distribution
x = (x - delta)/gam;          % Standardize
p = zeros(size(x));
if beta ==1
    p( x <= 0 ) = 0;
    p( x > 0 ) = sqrt(1/(2*pi)) * exp(-.5./x(x>0)) ./...
        x(x>0).^1.5;
else
    p(x >= 0) = 0;
    p(x < 0 ) = sqrt(1/(2*pi)) * exp(.5./x(x<0) ) ./...
        ( -x(x<0) ).^1.5;
end
p = p/gam; %rescale

elseif abs(alpha - 1) > 1e-5      % Gen. Case, alpha ~ 1

xold = x; % Save for later
% Standardize in (M) parameterization ( See equation (2) in [1] )
x = (x - delta)/gam - beta * tan(alpha*pi/2);

% Compute pdf
p = zeros(size(x));
zeta = - beta * tan(pi*alpha/2);
thetao = (1/alpha) * atan(beta*tan(pi*alpha/2));
A1 = alpha*thetao;
A2 = cos(A1)^(1/(alpha-1));
exp1 = alpha/(alpha-1);
alpham1 = alpha - 1;
c2 = alpha ./ (pi * abs(alpha - 1) * ( x(x>zeta) - zeta ) );
V = @(theta) A2 * ( cos(theta) ./ sin( alpha*(theta + thetao) ) ).^exp1.*...
    cos( A1 + alpham1*theta ) ./ cos(theta);

% x > zeta, calculate integral using QUADV
if any(x(:) > zeta)
    xshift = (x(x>zeta) - zeta) .^ exp1;

    if beta == -1 && alpha < 1
        p(x > zeta) = 0;
    elseif ~quick % Locate peak in integrand and split up integral
        g = @(theta) xshift(:) .* V(theta) - 1;
        R = repmat([-thetao, pi/2 ],numel(xshift),1);
        if abs(beta) < 1
            thetaz = bisectionSolver(g,R,alpha);
        else
            thetaz = bisectionSolver(g,R,alpha,beta,xshift);
        end
        thetaz = reshape(thetaz,size(xshift));
        % change variables so the two integrals go from

```

```

% 0 to 1/2 and 1/2 to 1.
theta2shift1 = 2*(theta2 + theta0);
theta2shift2 = 2*(pi/2 - theta2);
g1 = @(theta) xshift .* ...
    V(theta2shift1 * theta - theta0);
g2 = @(theta) xshift .* ...
    V(theta2shift2 * (theta - .5) + theta2);
zexpz = @(z) max(0,z .* exp(-z)); % use max incase of NaN

p(x > zeta) = c2 .* ...
    (theta2shift1 .* quadv(@(theta) zexpz( g1(theta) ),...
        0 , .5, tol) ...
    + theta2shift2 .* quadv(@(theta) zexpz( g2(theta) ),...
        .5 , 1, tol) );

else % be quick - calculate integral without locating peak
    % Use a default tolerance of 1e-6
    g = @(theta) xshift * V(theta);
    zexpz = @(z) max(0,z .* exp(-z)); % use max incase of NaN
    p( x > zeta ) = c2 .* quadv(@(theta) zexpz( g(theta) ),...
        -theta0 , pi/2, tol );
end
p(x > zeta) = p(x>zeta)/gam; %rescale

end

% x = zeta, this is easy
if any( abs(x(:) - zeta) < 1e-8 )
    p( abs(x - zeta) < 1e-8 ) = max(0,gamma(1 + 1/alpha)*...
        cos(theta0)/(pi*(1 + zeta^2)^(1/(2*alpha))));
    p( abs(x - zeta) < 1e-8 ) = p( abs(x - zeta) < 1e-8 )/gam; %rescale
end

% x < zeta, recall function with -xold, -beta, -delta
% This doesn't need to be rescaled.
if any(x(:) < zeta)
    p( x < zeta ) = stblpdf( -xold( x<zeta ),alpha,-beta,...
        gam , -delta , tol , quick);
end

else % Gen case, alpha = 1

x = (x - (2/pi) * beta * gam * log(gam) - delta)/gam; % Standardize

% Compute pdf
piover2 = pi/2;
twooverpi = 2/pi;
oneoverb = 1/beta;
theta0 = piover2;
% Use logs to avoid overflow/underflow
logV = @(theta) log(twooverpi * ((piover2 + beta *theta)./cos(theta))) + ...
    ( oneoverb * (piover2 + beta *theta) .* tan(theta) );

```

```

c2 = 1/(2*abs(beta));
xterm = ( -pi*x/(2*beta));

if ~quick % Locate peak in integrand and split up integral
    % Use a default tolerance of 1e-12
    logg = @(theta) xterm(:) + logV(theta) ;
    R = repmat([-thetao, pi/2 ],numel(xterm),1);
    theta2 = bisectionSolver(logg,R,1-beta);
    theta2 = reshape(theta2,size(xterm));
    % change variables so the two integrals go from
    % 0 to 1/2 and 1/2 to 1.
    theta2shift1 = 2*(theta2 + thetao);
    theta2shift2 = 2*(pi/2 - theta2);
    logg1 = @(theta) xterm + ...
        logV(theta2shift1 * theta - thetao);
    logg2 = @(theta) xterm + ...
        logV(theta2shift2 * (theta - .5) + theta2);
    zexpz = @(z) max(0,exp(z) .* exp(-exp(z))); % use max incase of NaN

    p = c2 .* ...
        (theta2shift1 .* quadv(@(theta) zexpz( logg1(theta) ),...
            0 , .5, tol) ...
        + theta2shift2 .* quadv(@(theta) zexpz( logg2(theta) ),...
            .5 , 1, tol) );

else % be quick - calculate integral without locating peak
    % Use a default tolerance of 1e-6
    logg = @(theta) xterm + logV(theta);
    zexpz = @(z) max(0,exp(z) .* exp(-exp(z))); % use max incase of NaN
    p = c2 .* quadv(@(theta) zexpz( logg(theta) ),-thetao , pi/2, tol );

end

p = p/gam; %rescale

end

p = real(p); % just in case a small imaginary piece crept in
    % This might happen when (x - zeta) is really small

end

```

```

function X = bisectionSolver(f,R,alpha,varargin)
% Solves equation g(theta) - 1 = 0 in STBLPDF using a vectorized bisection
% method and a tolerance of 1e-5. The solution to this
% equation is used to increase accuracy in the calculation of a numerical
% integral.
%
% If alpha ~= 1 and 0 <= beta < 1, the equation always has a solution

```

```

%
% If alpha > 1 and beta <= 1, then g is monotone decreasing
%
% If alpha < 1 and beta < 1, then g is monotone increasing
%
% If alpha = 1, g is monotone increasing if beta > 0 and monotone
% decreasing if beta < 0. Input alpha = 1 - beta to get desired results.
%
%

if nargin < 2
    error('bisectionSolver:TooFewInputs','Requires at least two input arguments.');
```

end

```

noSolution = false(size(R,1));
% if ~isempty(varargin)
%     beta = varargin{1};
%     xshift = varargin{2};
%     if abs(beta) == 1
%         Vo=(1/alpha)^(alpha/(alpha-1))*(1-alpha)*cos(alpha*pi/2)*xshift;
%         if alpha > 1
%             noSolution = Vo - 1 %>= 0;
%         elseif alpha < 1
%             noSolution = Vo - 1 %<= 0;
%         end
%     end
% end
% end

tol = 1e-6;
maxiter = 30;

[N M] = size(R);
if M ~= 2
    error('bisectionSolver:BadInput',...
        '"R" must have 2 columns');
end

a = R(:,1);
b = R(:,2);
X = (a+b)/2;

try
    val = f(X);
catch ME
    error('bisectionSolver:BadInput',...
        'Input function inconsistent with rectangle dimension')
end

if size(val,1) ~= N
    error('bisectionSolver:BadInput',...
        'Output of function must be a column vector with dimension of input');
end
```

```

% Main loop
val = inf;
iter = 0;

while( max(abs(val)) > tol && iter < maxiter )
    X = (a + b)/2;
    val = f(X);
    l = (val > 0);
    if alpha > 1
        l = 1-l;
    end
    a = a.*l + X.*(1-l);
    b = X.*l + b.*(1-l);
    iter = iter + 1;
end

if any(noSolution(:))
    X(noSolution) = (R(1,1) + R(1,2))/2;
end

end

```