

Help

```

#if defined(PremiaCurrentVersion) && PremiaCurrentVersion <
    (2008+2) //The "#else" part of the code will be freely av
    ailable after the (year of creation of this file + 2)
#else

#include "functions.h"
#include "common.h"
#include<iostream>
#include <iostream>
#include<cmath>
extern "C"{
#include "pnl/pnl_mathtools.h"
}
using namespace std;

    const long double inverselaplace::InvLF(const long double&
        T)const
{
    int n=10,k;

    std::vector<long double> v;
    long double deux=2.0;
    v.resize(n+1);
    v[1]=0.0833333333;v[2]=-32.08333333;v[3]=1279.000076;v[4]=
        -15623.66689;v[5]=84244.16946;
    v[6]=-236957.5129;v[7]=375911.6923;v[8]=-340071.6923;v[9]=
        164062.5128;v[10]=-32812.50256;
    long double s=0.0;
    for(k=1;k<=n;k++)
    {
        s=s+v[k]*(*LF).f(k*log(deux)/T);
    }

    return s*log(deux)/T;
}
std::vector<long double> newton::racine()const
{
    long double x=start, x0=start, temp;
    std::vector<long double> sol(1);
    int n=1;

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do
{
    temp=x;
    x=x0-(*F).f(x0)/(*F).Df(x0);
    x0=temp;
    n++;
}while((*F).f(x)>accuracy && n<=1000);
sol[0]=x;
return sol;
}

std::vector<long double> secante::racine()const
{
    long double x=start1, x0=start0, x1=start1, temp;
    std::vector<long double> sol(1);
    int n=1;
    do
    {
        temp=x;
        x=x1-(*F).f(x1)/(((F).f(x1)-(F).f(x0))/(x1-x0));
        x0=temp;
        x1=x;
        n++;
    }while((*F).f(x)>accuracy && n<=1000);
    sol[0]=x;
    return sol;
}

std::vector<long double> dichotomie::racine()const
{
    long double x0=a, x1=b, x,fx0=(F).f(a), fx1=(F).f(b),fx
    ;
    std::vector<long double> sol(1);

    x=(x0+x1)/2;
    fx=(F).f(x);
    sol[0]=x0;
    if(abs(fx0)<=accuracy)
        return sol;
    sol[0]=x1;
    if(abs(fx1)<=accuracy)
        return sol;
    sol[0]=x;

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    if(abs(fx)<=accuracy)
        return sol;
    while(abs(fx)>accuracy)
    {
        if(fx*fx0<0)
        {
            x1=x;
            fx1=fx;
        }
        else
        {
            x0=x;
            fx0=fx;
        }
        x=(x0+x1)/2;
        fx=(*F).f(x);
    }
    sol[0]=x;
    return sol;
}

const long double KCE::f(const long double &x)const
{
    double
    nu=param[0],
    sigma=param[1],
    lambda=param[2],
    p=param[3],
    eta1=param[4],
    eta2=param[5],
    alpha=param[6],
    q=param[7];
    return nu*x+sigma*sigma*x*x/2+lambda*(p*eta1/(eta1-x)+q*
        eta2/(eta2+x)-1)-alpha;
}

const long double KCE::Df(const long double &x)const
{
    double
    nu=param[0],
    sigma=param[1],
    lambda=param[2],
    p=param[3],

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    eta1=param[4],
    eta2=param[5],
    q=param[7];
    return nu+sigma*sigma*x+lambda*(p*eta1/((eta1-x)*(eta1-x)
        )-q*eta2/((eta2+x)*(eta2+x)));
}
const long double LPLB::f(const long double& alpha) const
{
    long double
    nu=param[0],
    sigma=param[1],
    lambda=param[2],
    p=param[3],
    eta1=param[4],
    eta2=param[5],
    S0=param[6],
    r=param[7],
    M=param[9],
        delta=param[10],
        eps=0.0000000000000001;
    long double beta1, beta2,x,x0;
        long double y[8]={nu,sigma,lambda,p,eta1,eta2,alpha
+r,1.-p};
        KCE G(y);
            x=eta1-0.1;
            while(G.f(x)<0)
                x=x+1e-2;
            dichotomie solG(G,eps,x,0.00001);
            beta1=solG.racine()[0];
            newton solG1(G,beta1,eps);
            beta1=solG1.racine()[0];

            x=eta1+10;
            x0=eta1+0.1;
            while(G.f(x0)>0)
                x0=x0-1e-2;
            while(G.f(x)<0)
                x=x+10;
            dichotomie solG2(G,x0,x,0.00001);
            beta2=solG2.racine()[0];
            newton solG3(G,beta2,eps);

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        beta2=solG3.racine()[0];

    long double A=(eta1-beta1)*beta2/(beta1-1);
    long double B=(beta2-eta1)*beta1/(beta2-1);
    long double C=(alpha+r)*eta1*(beta2-beta1);

    return S0*A/C*pow(S0/M,beta1-1)+S0*B/C*pow(S0/M,beta2-1)
        +M/(alpha+r)-S0/(alpha+delta);
}
const long double DLPLB::f(const long double& alpha) const
{
    long double
    nu=param[0],
    sigma=param[1],
    lambda=param[2],
    p=param[3],
    eta1=param[4],
    eta2=param[5],
    S0=param[6],
    r=param[7],
    M=param[9],
        delta=param[10],
        eps=0.0000000000000001;
    long double beta1, beta2,x,x0;
        long double y[8]={nu,sigma,lambda,p,eta1,eta2,alpha
+r,1.-p};
        KCE G(y);
        x=eta1-0.1;
        while(G.f(x)<0)
            x=x+1e-2;
        dichotomie solG(G,eps,x,0.00001);
        beta1=solG.racine()[0];
        newton solG1(G,beta1,eps);
        beta1=solG1.racine()[0];

        x=eta1+10;
        x0=eta1+0.1;
        while(G.f(x0)>0)
            x0=x0-1e-2;
        while(G.f(x)<0)
            x=x+10;

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        dichotomie solG2(G,x0,x,0.00001);
        beta2=solG2.racine()[0];
        newton solG3(G,beta2,eps);
        beta2=solG3.racine()[0];

    long double A=(eta1-beta1)*beta2/(beta1-1);
    long double B=(beta2-eta1)*beta1/(beta2-1);
    long double C=(alpha+r)*eta1*(beta2-beta1);

    return beta1*A/C*pow(S0/M,beta1-1)+beta2*B/C*pow(S0/M,
        beta2-1)-1/(alpha+delta);
}

const long double LCLB::f(const long double& alpha) const
{
    long double
    nu=param[0],
    sigma=param[1],
    lambda=param[2],
    p=param[3],
    eta1=param[4],
    eta2=param[5],
    S0=param[6],
    r=param[7],
    m=param[9],
        delta=param[10],
        eps=1e-16;
    long double beta1, beta2, x, x0;
        long double y[8]={nu,sigma,lambda,p,eta1,eta2,alpha
+r,1.-p};
        KCE G(y);
        x=eta1-0.1;
        while(G.f(x)<0)
            x=x+1e-2;
        dichotomie solG(G,eps,x,1e-5);
        beta1=solG.racine()[0];
        newton solG1(G,beta1,eps);
        beta1=solG1.racine()[0];

        x=eta1+10;x0=eta1+0.1;
        while(G.f(x0)>0)

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        x0=x0-1e-2;
while(G.f(x)<0)
    x=x+10;
dichotomie solG2(G,x0,x,1e-5);
beta2=solG2.racine()[0];
newton solG3(G,beta2,eps);
beta2=solG3.racine()[0];

long double A=(eta1-beta1)*beta2/(beta1+1);
long double B=(beta2-eta1)*beta1/(beta2+1);
long double C=(alpha+r)*eta1*(beta2-beta1);

return S0*A/C*pow(m/S0,beta1+1)+S0*B/C*pow(m/S0,beta2+1)
    -m/(alpha+r)+S0/(alpha+delta);
}
const long double DLCLB::f(const long double& alpha) const
{
long double
    nu=param[0],
    sigma=param[1],
    lambda=param[2],
    p=param[3],
    eta1=param[4],
    eta2=param[5],
    S0=param[6],
    r=param[7],
    m=param[9],
    delta=param[10],
    eps=1e-16;
long double beta1, beta2, x, x0;
    long double y[8]={nu,sigma,lambda,p,eta1,eta2,alpha
+r,1.-p};
    KCE G(y);
    x=eta1-0.1;
    while(G.f(x)<0)
        x=x+1e-2;
    dichotomie solG(G,eps,x,1e-5);
    beta1=solG.racine()[0];
    newton solG1(G,beta1,eps);
    beta1=solG1.racine()[0];

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        x=eta1+10;x0=eta1+0.1;
        while(G.f(x0)>0)
            x0=x0-1e-2;
        while(G.f(x)<0)
            x=x+10;
        dichotomie solG2(G,x0,x,1e-5);
        beta2=solG2.racine()[0];
        newton solG3(G,beta2,eps);
        beta2=solG3.racine()[0];

    long double A=(eta1-beta1)*beta2/(beta1+1);
    long double B=(beta2-eta1)*beta1/(beta2+1);
    long double C=(alpha+r)*eta1*(beta2-beta1);

    return -beta1*A/C*pow(m/S0,beta1+1)-beta2*B/C*pow(m/S0,
        beta2+1)+1/(alpha+delta);
}
const long double LPsiM::f(const long double& alpha)const
{
    long double
    nu=param[0],
    sigma=param[1],
    lambda=param[2],
    p=param[3],
    eta1=param[4],
    eta2=param[5],
    b=param[7],
    eps=1e-16;
    long double beta1, beta2, x, x0;
    long double y[8]={nu,sigma,lambda,p,eta1,eta2,alp
    ha,1-p};
    KCE G(y);
    x=eta1-0.1;
    while(G.f(x)<0)
        x=x+1e-2;
    dichotomie solG(G,eps,x,1e-5);
    beta1=solG.racine()[0];
    newton solG1(G,beta1,eps);
    beta1=solG1.racine()[0];

    x=eta1+10;x0=eta1+0.1;

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        while(G.f(x0)>0)
            x0=x0-1e-2;
        while(G.f(x)<0)
            x=x+10;
        dichotomie solG2(G,x0,x,1e-5);
        beta2=solG2.racine()[0];
        newton solG3(G,beta2,eps);
        beta2=solG3.racine()[0];

        return ((eta1-beta1)*beta2*expl(-b*beta1)+(beta2-eta1)*
            beta1*expl(-b*beta2))/((beta2-beta1)*eta1*alpha);
    }
    const long double LdPsiM::f(const long double& alpha)const
    {
        long double
        nu=param[0],
        sigma=param[1],
        lambda=param[2],
        p=param[3],
        eta1=param[4],
        eta2=param[5],
        b=param[7],
        eps=1e-16;
        long double beta1, beta2,x,x0;
        long double y[8]={nu,sigma,lambda,p,eta1,eta2,alpha,1-p};
        KCE G(y);
        x=eta1-0.1;
        while(G.f(x)<0)
            x=x+1e-2;
        dichotomie solG(G,eps,x,1e-5);
        beta1=solG.racine()[0];
        newton solG1(G,beta1,eps);
        beta1=solG1.racine()[0];

        x=eta1+10;x0=eta1+0.1;
        while(G.f(x0)>0)
            x0=x0-1e-2;
        while(G.f(x)<0)
            x=x+10;
        dichotomie solG2(G,x0,x,1e-5);

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        beta2=solG2.racine()[0];
        newton solG3(G,beta2,eps);
        beta2=solG3.racine()[0];

    return ((eta1-beta1)*beta2*expl(-b*beta1)+(beta2-eta1)*
        beta1*expl(-b*beta2))/((beta2-beta1)*eta1);
}

const long double LPsiMA::f(const long double& alpha)const
{
    long double
        nu=param[0],
        sigma=param[1],
        lambda=param[2],
        p=param[3],
        eta1=param[4],
        eta2=param[5],
        b=param[7],
        eps=1e-16;
    long double beta1, beta2, x, x0;
        long double y[8]={nu,sigma,lambda,p,eta1,eta2,alpha,1-p};
        KCE G(y);
        x=eta1-0.1;
        while(G.f(x)<0)
            x=x+1e-2;
        dichotomie solG(G,eps,x,1e-5);
        beta1=solG.racine()[0];
        newton solG1(G,beta1,eps);
        beta1=solG1.racine()[0];

        x=eta1+10;x0=eta1+0.1;
        while(G.f(x0)>0)
            x0=x0-1e-2;
        while(G.f(x)<0)
            x=x+10;
        dichotomie solG2(G,x0,x,1e-5);
        beta2=solG2.racine()[0];
        newton solG3(G,beta2,eps);
        beta2=solG3.racine()[0];

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    return ((eta1-beta1)*expl(-b*beta1)+(beta2-eta1)*expl(-
        b*beta2))/(beta2-beta1);
}
const long double LPsiMB::f(const long double& alpha) const
{
    long double
    nu=param[0],
    sigma=param[1],
    lambda=param[2],
    p=param[3],
    eta1=param[4],
    eta2=param[5],
    b=param[7],
    eps=1e-16;
    long double beta1, beta2, x, x0;
    long double y[8]={nu,sigma,lambda,p,eta1,eta2,alpha,1-p};
    KCE G(y);
    x=eta1-0.1;
    while(G.f(x)<0)
        x=x+1e-2;
    dichotomie solG(G,eps,x,1e-5);
    beta1=solG.racine()[0];
    newton solG1(G,beta1,eps);
    beta1=solG1.racine()[0];

    x=eta1+10;x0=eta1+0.1;
    while(G.f(x0)>0)
        x0=x0-1e-2;
    while(G.f(x)<0)
        x=x+10;
    dichotomie solG2(G,x0,x,1e-5);
    beta2=solG2.racine()[0];
    newton solG3(G,beta2,eps);
    beta2=solG3.racine()[0];

    return ((eta1-beta1)*(beta2-eta1)*(expl(-b*beta1)-expl(
        -b*beta2)))/((beta2-beta1)*eta1);
}
const long double amer_eq::f(const long double& v) const
{

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long double
    ksi=param[2]*param[3]/(param[3]-1)+(1-param[2])*
    param[4]/(param[4]+1)-1,
nu=(param[7]-param[9])-param[0]*param[0]/2-param[1]*ksi,

sigma=param[0],
lambda=param[1],
p=param[2],
eta1=param[3],
eta2=param[4],
K=param[6],
r=param[7],
    delta=param[9],
T=param[8],
    C=param[10],
    D=param[11];

    long double x[8]={nu,sigma,lambda,p,eta1,eta2,log(
K/v),T};
    long double cst1=psiVN(x,7);
    x[0]=r-delta+sigma*sigma/2-lambda*ksi;
x[2]=lambda*(ksi+1);
x[3]=p*eta1/((1+ksi)*(eta1-1));
x[4]=eta1-1;
x[5]=eta2+1;

    long double cst2=psiVN(x,7);
    long double EuP=K*exp(-r*T)*(1-cst1)-v*exp(-delta*
T)*(1-cst2);
    long double proba=1-cst1;

return C*K-D*(v*exp(-delta*T)+EuP)-(C-D)*K*exp(-r*T)*
    proba;
}
////////////////////////////////////
////////////////////////////////////
////////
//fonction psi=P[ZT>=a] ou Z levy poisson compose long
//double exponentielle
long double psiVN(long double *x,int Nb)
{

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long double
nu=x[0],
sigma=x[1],
lambda=x[2],
p=x[3],
eta1=x[4],
eta2=x[5],
a=x[6],
T=x[7];

if(T<1e-16)
    return (a>0? 0.0:1.0);
std::vector<std::vector<long double> > pm,qm;
std::vector<long double> pi;
long double zr=0.0;
int i,k,n;

pm.resize(Nb+1);
for(i=0;i<=Nb;i++)
    pm[i].resize(i+1);
qm.resize(Nb+1);
for(i=0;i<=Nb;i++)
    qm[i].resize(i+1);
pi.resize(Nb+1);
pi[0]=expl(-x[2]*x[7]);
for(n=1;n<=Nb;n++)
{
    pi[n]=expl(-lambda*T)*powl(lambda*T,n)/(fact_dia(n));
    pm[n][n]=powl(p,n);
    qm[n][n]=powl(1-p,n);
    for(k=1;k<n;k++)
    {
        pm[n][k]=zr;
        qm[n][k]=zr;
        for(i=k;i<n;i++)
        {
            pm[n][k]=pm[n][k]+bin_dia(n,i)*powl(x[3],i)*powl(x[4]/(x[4]
                +x[5]),i-k)*powl((1-x[3])*x[5]/(x[4]+x[5]),n-i)*bin_dia(n-
                k-1,i-k);
        }
    }
}

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qm[n][k]=qm[n][k]+bin_dia(n,i)*powl(1-x[3],i)*powl(x[3]*x[4]
    ]/(x[4]+x[5]),n-i)*powl(x[5]/(x[4]+x[5]),i-k)*bin_dia(n-k-
    1,i-k);
    }
}
}
long double s1=zr,s2=zr;
long double c1=expl(sigma*eta1*sigma*eta1*T/2)/(sigma*sq
    rtl(2*M_PI*T));
long double c2=expl(sigma*eta2*sigma*eta2*T/2)/(sigma*sq
    rtl(2*M_PI*T));

long double c3, c4;
std::vector<long double> Inv1(Nb),Inv2(Nb);

Inv1=In(a-nu*T,-eta1,-1./(sigma*sqrtl(T)),-sigma*eta1*sq
    rtl(T),Nb);
Inv2=In(a-nu*T,eta2,1./(sigma*sqrtl(T)),-sigma*eta2*sqrtl(
    T),Nb);
c3=sigma*sqrtl(T)*eta1;
c4=sigma*sqrtl(T)*eta2;
for(n=1;n<=Nb;n++)
{
    for(k=1;k<=n;k++)
    {
        s1=s1+pi[n]*pm[n][k]*powl(c3,k)*Inv1[k-1];
        s2=s2+pi[n]*qm[n][k]*powl(c4,k)*Inv2[k-1];
    }
}
if(c1>1e+100)
{
    c1=1.;
    s1=(1-pi[0])*p*Hh0((a-nu*T)/(sigma*sqrtl(T)))/sqrtl(2*M_
        PI);
}
if(c2>1e+100)
{
    c2=1.;
    s2=(1-pi[0])*(1-p)*Hh0((a-nu*T)/(sigma*sqrtl(T)))/sqrtl(2
        *M_PI);
}

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}
return c1*s1+c2*s2+pi[0]*Hh0((a-nu*T)/(sigma*sqrtl(T)))/sq
    rtl(2*M_PI);
}
////////////////////////////////////
////////////////////////////////////
////////////////////////////////////
//derivÃ de psi=P[ZT>=a] % Ã a
////////////////////////////////////
////////////////////////////////////
////////////////////////////////////
//derivÃ de psi=P[ZT>=a] % Ã a
long double dpsivN(long double *x,int Nb)
{
    long double
    nu=x[0],
    sigma=x[1],
    lambda=x[2],
    p=x[3],
    eta1=x[4],
    eta2=x[5],
    a=x[6],
    T=x[7];

    std::vector<std::vector<long double> > pm,qm;
    std::vector<long double> pi;
    long double zr=0;
    int i,k,N;

    pm.resize(Nb+1);
    for(i=0;i<=Nb;i++)
        pm[i].resize(i+1);
    qm.resize(Nb+1);
    for(i=0;i<=Nb;i++)
        qm[i].resize(i+1);
    pi.resize(Nb+1);
    pi[0]=exp(-x[2]*x[7]);
    for(N=1;N<=Nb;N++)
    {
        pi[N]=exp(-lambda*T)*pow(lambda*T,N)/fact_dia(N);
        pm[N][N]=pow(p,N);
    }
}

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```

qm[N][N]=pow(1-p,N);
for(k=1;k<N;k++)
{
    pm[N][k]=zr;
    qm[N][k]=zr;
    for(i=k;i<N;i++)
    {

pm[N][k]=pm[N][k]+bin_dia(N,i)*pow(x[3],i)*pow(x[4]/(x[4]+x
    [5]),i-k)*pow((1-x[3])*x[5]/(x[4]+x[5]),N-i)*bin_dia(N-k-1
    ,i-k);

qm[N][k]=qm[N][k]+bin_dia(N,i)*pow(1-x[3],i)*pow(x[3]*x[4]/
    (x[4]+x[5]),N-i)*pow(x[5]/(x[4]+x[5]),i-k)*bin_dia(N-k-1,
    i-k);
    }
}
}
long double s1=zr,s2=zr;
long double c1=exp(sigma*eta1*sigma*eta1*T/2)/(sigma*sqrt(
    2*M_PI*T));
long double c2=exp(sigma*eta2*sigma*eta2*T/2)/(sigma*sqrt(
    2*M_PI*T));
long double c3, c4;
std::vector<long double> Inv1(Nb),Inv2(Nb);
Inv1=dIn(a-nu*T,-eta1,-1/(sigma*sqrt(T)),-sigma*eta1*sqrt(
    T),Nb);
Inv2=dIn(a-nu*T,eta2,1/(sigma*sqrt(T)),-sigma*eta2*sqrt(T)
    ,Nb);
for(N=1;N<=Nb;N++)
{
    c3=sigma*sqrt(T)*eta1;
    c4=sigma*sqrt(T)*eta2;
    for(k=1;k<=N;k++)
    {
        s1=s1+pi[N]*pm[N][k]*c3*Inv1[k-1];
        s2=s2+pi[N]*qm[N][k]*c4*Inv2[k-1];
        c3=c3*sigma*sqrt(T)*eta1;
        c4=c4*sigma*sqrt(T)*eta2;
    }
}

```



```

if(c1>1e+100)
{
    c1=1.;
    s1=-(1-pi[0])*p*dnorm(-(a-nu*T)/(sigma*sqrt(T)))/(sigma*
        sqrt(T));
}
if(c2>1e+100)
{
    c2=1.;
    s2=-(1-pi[0])*(1-p)*dnorm(-(a-nu*T)/(sigma*sqrt(T)))/(si
        gma*sqrt(T));
}
return c1*s1+c2*s2-pi[0]*dnorm(-(a-nu*T)/(sigma*sqrt(T)))/
    (sigma*sqrt(T));
}
////////////////////////////////////
////////////////////////////////////
////////////////////////////////////
//fonction psi=P[ZT+ksi+>=a] ou Z levy poisson compose long
//double exponentielle
std::vector<long double> psiVNB(long double *x,int Nb)
{
    long double
    nu=x[0],
    sigma=x[1],
    lambda=x[2],
    p=x[3],
    eta1=x[4],
    eta2=x[5],
    a=x[6],
    T=x[7];
    std::vector<long double> pi,y;
    y.resize(2);
    std::vector<std::vector<long double> > pm,qm;
    long double zr=0.0;
    int i,k,n,j;

    if(T<1e-16)
    {
        if(a>0)
        {

```

```

        y[0]=0.0;
        y[1]=expl(-eta1*a);
    }
    else
    {
        y[0]=1.0;
        y[1]=1.0;
    }

    return y;
}

pm.resize(Nb+1);
for(i=0;i<=Nb;i++)
    pm[i].resize(i+1);
qm.resize(Nb+1);
for(i=0;i<=Nb;i++)
    qm[i].resize(i+1);
pi.resize(Nb+1);
pi[0]=expl(-x[2]*x[7]);
for(n=1;n<=Nb;n++)
{
    pi[n]=expl(-lambda*T)*powl(lambda*T,n)/(fact_dia(n));
    pm[n][n]=powl(p,n);
    qm[n][n]=powl(1-p,n);
    for(k=1;k<n;k++)
    {
        pm[n][k]=zr;
        qm[n][k]=zr;
        for(i=k;i<n;i++)
        {
            pm[n][k]=pm[n][k]+bin_dia(n,i)*powl(x[3],i)*powl(x[4]/(x[4]
                +x[5]),i-k)*powl((1-x[3])*x[5]/(x[4]+x[5]),n-i)*bin_dia(n-
                k-1,i-k);

            qm[n][k]=qm[n][k]+bin_dia(n,i)*powl(1-x[3],i)*powl(x[3]*x[4]
                /(x[4]+x[5]),n-i)*powl(x[5]/(x[4]+x[5]),i-k)*bin_dia(n-k-
                1,i-k);
        }
    }
}
}

```

```

std::vector<std::vector<long double> > pmb,qmb;
pmb.resize(Nb+1);
for(i=0;i<=Nb;i++)
    pmb[i].resize(i+2);
qmb.resize(Nb+1);
for(i=0;i<=Nb;i++)
    qmb[i].resize(i+1);

for(n=1;n<=Nb;n++)
{
    pmb[n][1]=0;
    for(i=1;i<=n;i++)
        pmb[n][1]=pmb[n][1]+qm[n][i]*pow(eta2/(eta2+eta1),i);
    for(i=2;i<=n+1;i++)
    {
        pmb[n][i]=pm[n][i-1];
    }
    for(i=1;i<=n;i++)
    {
        qmb[n][i]=0;
        for(j=i;j<=n;j++)
            qmb[n][i]=qmb[n][i]+(eta1/(eta1+eta2))*pow(eta2/(eta1+eta2),j-i)*qm[n][j];
    }
}

long double s1=zs,s2=zs,s3=zs,s4=zs;
long double c1=expl(sigma*eta1*sigma*eta1*T/2)/(sigma*sqrtl(2*M_PI*T));
long double c2=expl(sigma*eta2*sigma*eta2*T/2)/(sigma*sqrtl(2*M_PI*T));
long double c=expl(sigma*eta1*sigma*eta1*T/2)*eta1/sqrtl(2*M_PI);
long double c3, c4;
std::vector<long double> Inv1(Nb+1),Inv2(Nb);

Inv1=In(a-nu*T,-eta1,-1./(sigma*sqrtl(T)),-sigma*eta1*sqrtl(T),Nb+1);
Inv2=In(a-nu*T,eta2,1./(sigma*sqrtl(T)),-sigma*eta2*sqrtl(T),Nb);
c3=sigma*sqrtl(T)*eta1;

```

```

    c4=sigma*sqrtl(T)*eta2;
    for(n=1;n<=Nb;n++)
    {
        for(k=1;k<=n;k++)
        {
            s1+=pi[n]*pm[n][k]*powl(c3,k)*Inv1[k-1];
            s3+=pi[n]*pmb[n][k]*powl(c3,k)*Inv1[k-1];
            s2+=pi[n]*qm[n][k]*powl(c4,k)*Inv2[k-1];
            s4+=pi[n]*qmb[n][k]*powl(c4,k)*Inv2[k-1];
        }
        s3+=pi[n]*pmb[n][n+1]*powl(c3,n+1)*Inv1[n];
    }
    if(c>1e+100)
    {
        c=Hh0((a-nu*T)/(sigma*sqrtl(T)))/(sqrtl(2*M_PI));
    }
    else
        c*=I0(a-nu*T,-eta1,-1./(sigma*sqrt(T)),-eta1*sigma*sqrt(T));
    if(c1>1e+100)
    {
        c1=1.;
        s1=(1-pi[0])*p*Hh0((a-nu*T)/(sigma*sqrtl(T)))/sqrtl(2*M_PI);
        s3=s1;
    }
    if(c2>1e+100)
    {
        c2=1.;
        s2=(1-pi[0])*(1-p)*Hh0((a-nu*T)/(sigma*sqrtl(T)))/sqrtl(2*M_PI);
        s4=s2;
    }
    y[0]=c1*s1+c2*s2+pi[0]*Hh0((a-nu*T)/(sigma*sqrtl(T)))/sqrtl(2*M_PI);
    y[1]=c1*s3+c2*s4+pi[0]*c;

    return y;
}
//psi=P[ZT>=a,maxZs>=b s dans[0 T]] ou Z levy poisson compose long double exponentielle

```

[illegible]

```

//prix du call loockback floating strike
long double CLB(long double *x,const long double &T)
{
    LCLB LF(x);
    inverselaplace F(LF);
    return F.InvLF(T);
};
////////////////////////////////////
////////////////////////////////////
////////////////////////////////////
//delta du put loockback floating strike
long double dPLB(long double *x,const long double &T)
{
    DLPLB LF(x);
    inverselaplace F(LF);
    return F.InvLF(T);
};
////////////////////////////////////
////////////////////////////////////
////////////////////////////////////
//delta du call loockback floating strike
long double dCLB(long double *x,const long double &T)
{
    DLCLB LF(x);
    inverselaplace F(LF);
    return F.InvLF(T);
};
////////////////////////////////////
////////////////////////////////////
////////////////////////////////////

////////////////////////////////////
////////////////////////////////////
////////////////////////////////////
//psi=P[maxZs>=b s dans[0 T]] ou Z levy poisson compose lon
g double exponentielle
long double psiM(long double *x,const long double &T)
{
    LPsiM LF(x);
    inverselaplace F(LF);
    return F.InvLF(T);
};

```

```

};
////////////////////////////////////
////////////////////////////////////
////////////////////////////////////
long double dpsim(long double *x,const long double &T)
{
    LdPsiM LF(x);
    inverselaplace F(LF);
    return F.InvLF(T);
};
////////////////////////////////////
////////////////////////////////////
////////////////////////////////////
long double rebateproba(long double *x,const long double &
    r,const long double &T)
{
    long double sm=0.0, pas=0.01, s=0.0;
    int N;
    N=(int)(T/pas);
    if(T-N*pas<1e-16)
        N-=1;
    for(int n=1;n<=N;n++)
    {
        s=n*pas;
        sm+=exp(-r*s)*dpsim(x,s)*pas;
    }
    return sm;
};
//psi=P[maxZs>=b s dans[0 T]] ou Z levy poisson compose lon
g double exponentielle
long double psiMA(long double *x,const long double &T)
{
    LPsiMA LF(x);
    inverselaplace F(LF);
    return F.InvLF(T);
};
////////////////////////////////////
////////////////////////////////////
////////////////////////////////////
//psi=P[maxZs>=b s dans[0 T]] ou Z levy poisson compose lon
g double exponentielle

```

```

long double psiMB(long double *x,const long double &T)
{
    LPsiMB LF(x);
    inverselaplace F(LF);
    return F.InvLF(T);
};
////////////////////////////////////
    //////////////////////////////////
    //////////////////////////////////
    //////////////////////////////////

#endif //PremiaCurrentVersion

```

References