

Help

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#include "cir1d_std.h"
#include "enums.h"

#if defined(PremiaCurrentVersion) && PremiaCurrentVersion <
    (2009+2) //The "#else" part of the code will be freely av
    ailable after the (year of creation of this file + 2)
static int CHK_OPT(MC_TEICHMANNBAYER)(void *Opt, void *Mod)
{
    return NONACTIVE;
}
int CALC(MC_TEICHMANNBAYER)(void*Opt,void *Mod,Pricing
    Method *Met)
{
    return AVAILABLE_IN_FULL_PREMIA;
}
#else

/* linear uniform interpolation of [0,T] of size N*/
/* return value = dt*/
static double linspace1(double T0, double T1, int N,
    double* t )
{
    double dt;
    int i;

    dt=(T1 - T0) / (double)( N - 1 );
    t[0] = T0;
    for (i=1; i<N; i++)
        t[i] = t[i-1] + dt;
    return dt;
}
/* linear interpolation using stepsize dt; return T */
static double linspace2( double dt, int N, double* t ){
    double T = dt * (double)( N-1 );
    int i;
    t[0] = 0.0;
    for (i=1; i<N; i++)
        t[i] = t[i-1] + dt;
    return T;
}

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}

/* extrapolate a CIR-HJM-forward rate curve from a short ra
   te */
/* l is assumed to be the length of x */
static void CIR2HJM( double r0, double kappa, double theta,
    double sigma, double gamma_tb, const double* x, int l,
    double* r ){
    double g0, g1, G1;
    int i;
    for (i=0; i<l; i++ ) {
        G1 = 2.0 * (exp( gamma_tb * x[i] ) - 1.0) / (( gamma_tb
            + kappa ) * (exp( gamma_tb * x[i] ) - 1.0) + 2.0 * gamma_
            tb );
        g0 = kappa * theta * G1;
        g1 = 4.0 * (gamma_tb * gamma_tb) * exp( gamma_tb * x[i]
            ) / (( ( gamma_tb + kappa ) * exp( gamma_tb * x[i] ) +
            gamma_tb - kappa ) * ( ( gamma_tb + kappa ) * exp( gamma_tb
            * x[i] ) + gamma_tb - kappa ));
        r[i] = g0 + r0 * g1;
    }
}

/* Generate an Nxd-array of iid. Bernoullis with p = 0.5 */
/* actually, it would be enough to consider boolean variab
   les here */
static void GenBernoulli1( int* J, int N,int generator)
{
    int i;
    for( i=0; i<N; i++)
        if (pnl_rand_uni(generator)< 0.5)
            J[i] = 0;
        else J[i] = 1;
}

/* generate a vector of "brownian increments" given the mu
   lti-index J */
static void omegadot1( int N, double dt, int NCub, const
    int* J, int n, double* dB ){
    int i,k;

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double tempd1 = sqrt(dt) / sqrt((double)(n));
for(i=0; i<(NCub-1); i++ ){
    for (k = 0; k<n; k++ )
        dB[i*n+k] = (J[i] == 0) ? tempd1 : (- tempd1);
}
for ( k = (n*(NCub-1)); k<N; k++ )
    dB[k] = (J[NCub-1] == 0) ? tempd1 : (-tempd1);
}

/* apply the shift semi-group */
/* i_shift is the integer part of the shift in terms of dx-
   units, r_shift
   the remainder */
static double Shift( const double* r, const double* x,
    double dx, double dt, int m, int k, int i_shift, double r_shift
){
    /*double ret;*/
    if (k < m - i_shift - 1 )
        return (1.0-r_shift) * r[k+i_shift] + r_shift * r[k+i_
            shift+1];
    else
        return r[m-1];
}

/* the vector field alpha0 */
static double alpha0( const double* r, double kappa,
    double theta, double sigma, double gamma_tb, const double* x,
    int m, int k, double expg){
    double g1, G1;
    G1 = 2.0 * (expg - 1.0) / (( gamma_tb + kappa ) * (expg -
        1.0) + 2.0 * gamma_tb );
    g1 = 4.0 * (gamma_tb * gamma_tb) * expg / (( ( gamma_tb +
        kappa ) * expg + gamma_tb - kappa ) * ( ( gamma_tb + kapp
        a ) * expg + gamma_tb - kappa ));
    return (sigma * sigma) * (g1) * ( r[0] * (G1) - 0.25 );
}

/* the volatility vector field(s) */
double HJMSigma( const double* r, double kappa, double thet
    a, double sigma, double gamma, const double* x, int m, int
    k, double expg, double sqrtr ){

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    return sigma * sqrtr * ( 4.0 * (gamma * gamma) * expg / (
        ( ( gamma + kappa ) * expg + gamma - kappa ) * ( ( gamma +
            kappa ) * expg + gamma - kappa )));
}

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/* value P(0,T) of a zero coupon bond */
static double ZeroCB( const double* r, const double* x,
    double dx, double T ){
    int Tx = ceil( T / dx ); /* index of T in the x-grid */
    double integ = 0.0;
    int i;
    for (i=0; i<Tx; i++ )
        integ += 0.5 * ( r[i] + r[i+1] ) * dx;
    return exp( - integ );
}

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/* compute the empirical mean value of a vector */
static double mean( const double* X, int M ){
    double ret = 0.0;
    int i;
    for (i=0; i<M; i++ )
        ret += X[i];
    ret = ret / (double)(M);
    return ret;
}

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/* compute the empirical standard deviation of a vector */
static double stdev( const double* X, int M ){
    double mu = mean( X, M );
    double ret = 0.0;
    int i;
    for(i=0; i<M; i++ )
        ret += X[i]*X[i];
    ret = ret / (double)(M);
    ret = sqrt( ret - mu * mu );
    return ret;
}

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/* n number of time intervals on each cubature interval*/
/* N number of time intervals*/

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/* m number of space intervals*/
/* M number of paths for Monte-Carlo simulation*/
static int mc_teichmannbayer(double r0, double kappa,
    double t0, double sigma,double theta, double T_bond, double T_
    option,NumFunc_1 *p,int generator,int n,int L,int k, int M,
    double *price,double *error)
{
    double gamma_tb = 0;
    int NCub;
    double *t;
    double dt = 0;
    double dx = 0;
    int mAct = 0;
    double *x;
    double r_shift = 0;
    int i_shift = 0;
    int *J;
    double *r;
    double *rp,*rm,*rpc,*rmc;
    double *res,*dB;
    double Bp, Bm; /* savings account */
    double expg, sqrtrm, sqrtrp; /* auxiliary variable*/
    int i,j;
    int ii,k_bis;
    double zerop = 0;
    double zerom = 0;

    gamma_tb = sqrt( (kappa*kappa) + 2.0 * (sigma * sigma) );
    pnl_rand_init(generator,1,M);

    /* first generate the time and space grids */
    if (L % n == 0)
        NCub = L / n;
    else
        NCub = L / n + 1;
    /* generate the time grid */
    t=malloc((L+1)*sizeof(double));
    dt = linspace1( t0, T_option, L+1, t );
    /* generate the spacegrid */
    dx = ( T_bond - T_option ) / (double)( k );
    mAct = (int)( T_option / dx ) + k + (int)(L * (dt/dx)

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) + 1;
x=malloc((mAct+1)*sizeof(double));
dt = linspace1( t0, T_option, L+1, t );
linspace2( dx, mAct + 1, x );
/* for the shift semigroup, express dt in temrs of dx
:
    dt = i_shift * dx + r_shift * dx */
r_shift = dt / dx;
i_shift = (int)( r_shift );
r_shift = r_shift - i_shift;

/* J describes one cubature path */
J=malloc((NCub)*sizeof(int));

/* generate the initial forward rate curve */
r=malloc((mAct+1)*sizeof(double));/* saves initial fo
rward rate curve */
CIR2HJM( r0, kappa, theta, sigma, gamma_tb, x, mAct+1
, r );

rp=malloc((mAct+1)*sizeof(double));
rm=malloc((mAct+1)*sizeof(double));
rpc=malloc((mAct+1)*sizeof(double));
rmc=malloc((mAct+1)*sizeof(double));

/* the path-wise discounted payoff */
res =malloc((M)*sizeof(double));

/* the "brownian" increments (i.e. the cubature deriv
atives) */
dB =malloc((L)*sizeof(double));

/* now iterate through all paths for the MC-simulatio
n*/
for(j=0; j<M; j++ ){
    /* re-initialize r and B */
    GenBernoulli1( J, NCub,generator ); /* generate J *
/

    for (i=0; i<mAct+1; i++ )
        {

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        rp[i]=r[i];
        rm[i]=r[i];
    }

    Bp = 1.0;
    Bm = 1.0;

    /* generate dB */
    omegadot1( L, dt, NCub, J, n, dB );

    /* iterate through the time grid */
    for(i=0; i<L; i++){
        sqrtrp = (rp[0] > 0.0) ? sqrt(rp[0]) : 0.0;
        sqrtrm = (rm[0] > 0.0) ? sqrt(rm[0]) : 0.0;
        Bp += Bp * rp[0] * dt;
        Bm += Bm * rm[0] * dt;
        for (ii=0; ii<mAct+1; ii++ )
        {
            rpc[ii]=rp[ii];
            rmc[ii]=rm[ii];
        }

        /* iterate through the space grid */
        for (k_bis=0; k_bis<=mAct; k_bis++){
            expg = exp( gamma_tb * x[k_bis] );
            rp[k_bis] = Shift( rpc, x, dx, dt, mAct+1, k_bis,
s, i_shift, r_shift ) + alpha0( rpc, kappa, theta, sigma,
gamma_tb, x, mAct+1, k_bis, expg ) * dt;
            rp[k_bis] += HJMSigma( rpc, kappa, theta, sigma
, gamma_tb, x, mAct+1, k_bis, expg, sqrtrp ) * dB[i];
            rm[k_bis] = Shift( rmc, x, dx, dt, mAct+1, k_bis,
s, i_shift, r_shift ) + alpha0( rmc, kappa, theta, sigma,
gamma_tb, x, mAct+1, k_bis, expg ) * dt;
            rm[k_bis] -= HJMSigma( rmc, kappa, theta, sigma
, gamma_tb, x, mAct+1, k_bis, expg,sqrtrm ) * dB[i];
        }
    }

    /* compute the discounted payoff for this particul
ar path */
    zerop = ZeroCB(rp, x, dx, T_bond - T_option );
    zerom = ZeroCB(rm, x, dx, T_bond - T_option );

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        res[j]=0.5*((p->Compute)(p->Par,zerop)/Bp+(p->Compute)(p->Par,zerom)/ Bm);
    }
    /*Price*/
    *price=mean( res, M );

    /*Estimate Error*/
    *error= 1.65 * stdev( res, M ) / sqrt( (double)(M) );

    /* free memory again */
    free(t);
    free(x);
    free(J);
    free(r);
    free(rp);
    free(rpc);
    free(rm);
    free(rmc);
    free(res);
    free(dB);

    return  OK;
}

int CALC(MC_TEICHMANNBAYER)(void *Opt,void *Mod,Pricing
    Method *Met)
{
    TYPEOPT* ptOpt=(TYPEOPT*)Opt;
    TYPEMOD* ptMod=(TYPEMOD*)Mod;

    return mc_teichmannbayer(ptMod->r0.Val.V_PDOUBLE,ptMod->
        k.Val.V_DOUBLE,ptMod->T.Val.V_DATE,ptMod->Sigma.Val.V_PDOUB
        LE,
                                ptMod->theta.Val.V_PDOUBLE,pt
        Opt->BMaturity.Val.V_DATE,ptOpt->OMaturity.Val.V_DATE,ptOpt->
        PayOff.Val.V_NUMFUNC_1, Met->Par[0].Val.V_ENUM.value,Met->
        Par[1].Val.V_PINT,Met->Par[2].Val.V_PINT,Met->Par[3].Val.V_PI
        NT,Met->Par[4].Val.V_PINT,&(Met->Res[0].Val.V_DOUBLE),&(
        Met->Res[1].Val.V_DOUBLE));
}

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static int CHK_OPT(MC_TEICHMANNBAYER)(void *Opt, void *Mod)
{
    if ((strcmp(((Option*)Opt)->Name,"ZeroCouponCallBondEuro"
        )==0) || (strcmp(((Option*)Opt)->Name,"ZeroCouponPutBondEuro")
        ==0))
        return OK;
    else
        return WRONG;
}

#endif //PremiaCurrentVersion
static int MET(Init)(PricingMethod *Met,Option *Opt)
{
    if ( Met->init == 0)
    {
        Met->init=1;
        Met->Par[0].Val.V_ENUM.value=0;
        Met->Par[0].Val.V_ENUM.members=&PremiaEnumMCRNGs;
        Met->Par[1].Val.V_PINT=20;
        Met->Par[2].Val.V_PINT=400;
        Met->Par[3].Val.V_PINT=10;
        Met->Par[4].Val.V_PINT=20;
    }

    return OK;
}

PricingMethod MET(MC_TEICHMANNBAYER)=
{
    "MC_TEICHMANNBAYER",

    {"RandomGenerator",ENUM,{100},ALLOW},
    {"Number of time intervals on each cubature interval",
        INT,{100},ALLOW},
    {"Number of time intervals",INT,{100},ALLOW},
    {"Number of space intervals*",INT,{100},ALLOW},
    {"Number of paths for Monte-Carlo simulation",PINT,{100},
        ALLOW},
    {" ",PREMIA_NULLTYPE,{0},FORBID}},
    CALC(MC_TEICHMANNBAYER),

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{{"Price",DOUBLE,{100},FORBID},{ "MC Error",DOUBLE,{100},
  FORBID} ,{" ",PREMIA_NULLTYPE,{0},FORBID}},
CHK_OPT(MC_TEICHMANNBAYER),
CHK_ok,
MET(Init)
} ;
```

References