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Apêndice A: Programa para Obtenção de Projeto Ótimo de Esquema Vp

O programa em linguagem C, mostrado a seguir, fornece, para $c_0 = 0,5, 1,0, 1,5, 2,0, 3,0, e 4,0$, para $h_s = 0,10, 0,25 e 0,50$, e para $\gamma^* = 1,5, 2,0 e 3,0$, o projeto ótimo de esquemas Vp, com obtenção das medidas de desempenho, em cada caso.

```

/*****
/*
/* Programa para projeto ótimo de esquemas Vp
/*
/*
/*****/

#include <stdlib.h>
#include <time.h>
#include <stdio.h>
#include <math.h>
#include <stdarg.h>
#include <assert.h>

/*****/
/* Funções auxiliares
*/
/*****/

double FatTable[20];
static FILE* fout = NULL;

/*
Calcula o fatorial de n. Utiliza uma tabela de constantes
pré-calculada para números menores que 20.
*/
double fat(int n)
{
    double val;

```

```
    if(n<20)
        return FatTable[n];

    val = FatTable[19];
    for(int i=20; i<=n; i++)
        val = val * i;
    return val;
}

/*
Inicializa a tabela de fatoriais pré-calculados
*/
void init_fat(void)
{
    int i;

    FatTable[0] = 1;
    for(i=1; i<20; i++)
        FatTable[i] = i * FatTable[i-1];
}

/*
Retorna a probabilidade de uma variável aleatória de Poisson
assumir um valor menor ou igual a "k".

k - Argumento para o qual a distribuição de Poisson será
avaliada.

theta - Media da distribuição de Poisson
*/
double dpoidf(int k, double theta)
{
    int i;
    double val = 0.0;

    for(i=0; i<=k; i++)
        val += (exp(-theta)*pow(theta,i)) / fat(i);

    return val;
}
```

```
/*
Imprime os parâmetros recebidos no arquivo de saída e na tela
*/
void print(char* format, ...)
{
    va_list list;

    if(!fout)
    {
        fout = fopen("resultado_vp.txt", "w");
        if(!fout)
        {
            printf("Erro abrindo arquivo de saida\n");
            exit(-1);
        }
    }

    va_start(list, format);
    vfprintf(fout, format, list);
    fflush(fout);
    vprintf(format, list);
    va_end(list);
}

/*****
/*  Funções de cálculo
*****/

#define P(a,b) p[(a)-1][(b)-1]

/*
Calcula as probabilidades de transição
*/
void calprob(double lsc1, double lsa1, double lsc2, double
             lsa2, double mS, double mL, double c0,
             double*r1,
             double p[3][3])
{
    double theta, a, b, c;
```

```
int k;

theta = mS * c0;
k      = (int)floor(lsa1);
P(1,1) = dpoidf(k,theta);

k      = (int)floor(lsc1);
P(1,2) = dpoidf(k,theta) - P(1,1);

theta = mL * c0;
k      = (int)floor(lsa2);
P(2,1) = dpoidf(k,theta);
k      = (int)floor(lsc2);
P(2,2) = dpoidf(k,theta) - P(2,1);

P(1,3) = 1 - P(1,1) - P(1,2);
P(2,3) = 1 - P(2,1) - P(2,2);
a      = P(2,3) - P(1,3);
b      = P(1,2) + P(2,1) + P(1,3) - P(2,3);
c      = -P(2,1);

assert(a);

*r1 = (-b + sqrt(b*b-4*a*c))/(2*a);
}

/*
Calcula o valor de TMAF
*/
void tmaf(double p[3][3], double r1, double hL, double hS,
          double* vtmaf)
{
double den, num;

den = (1-P(1,1))*(1-P(2,2))-P(1,2)*P(2,1);
num = ((1-P(2,2))*hL+P(1,2)*hS)*r1;
num = num + (P(2,1)*hL+(1-P(1,1))*hS)*(1-r1);

assert(den);
```

```
*vtmaf = num/den;
}

/*
Calcula o valor de TES
*/
void tes(double lsc1, double lsa1, double lsc2, double lsa2,
         double hL, double hS, double mS, double mL,
         double c1, double r1, double* vtes)
{
    double p[3][3];
    double r2, s, s1, s2, num, den, theta;
    int k;

    theta = mS * c1;
    k      = (int)floor(lsa1);
    P(1,1) = dpoidf(k,theta);
    k      = (int)floor(lsc1);
    P(1,2) = dpoidf(k,theta)-P(1,1);

    theta = mL*c1;
    k      = (int)floor(lsa2);
    P(2,1) = dpoidf(k,theta);
    k      = (int)floor(lsc2);
    P(2,2) = dpoidf(k,theta)-P(2,1);
    r2     = 1-r1;
    s      = r1*hL+r2*hS;
    s1     = r1*hL/s;
    s2     = r2*hS/s;
    den    = (1-P(1,1))*(1-P(2,2))-P(1,2)*P(2,1);
    num    = ((1-P(2,2))*hL+P(1,2)*hS)*s1 +
             (P(2,1)*hL+(1-P(1,1))*hS)*s2;

    assert(den);

    *vtes = num/den-(s1*hL+s2*hS)/2.0;
}

/*
Efetua os cálculos de TMAF e TES para os parâmetros correntes
*/
```

```
*/
int calcula(double c0, double gama, double lsc1, double lsa1,
            double lsc2, double lsa2, double mS, double mL,
            double hS, double* hL, double* vtes, double*
            vtmaf, double tmaffp)
{
    double p[3][3];
    double r1, mbarra;
    double c1 = c0 * gama;

    calprob(lsc1, lsa1, lsc2, lsa2, mS, mL, c0, &r1, p);

    assert(r1);

    mbarra = r1*mS+(1-r1)*mL;
    if (mbarra > 1.0)
        return 0;

    *hL=(1.0-(1.0-r1)*hS)/r1;

    tmaf(p, r1, *hL, hS, vtmaf);

    if (*vtmaf < tmaffp)
        return 0;

    tes(lsc1, lsa1, lsc2, lsa2, *hL, hS, mS, mL, c1, r1, vtes);
    return 1;
}

/*
Rotina principal. Contém os loops para otimização.
*/
void main(void)
{
    time_t t;

    double c0, gama, lsc1, lsa1, lsc2, lsa2, mS, mL,
           hL, hS, vtes, vtmaf;

    double lscfp, alfa, tmaffp, beta, tesfp;
```

```
double tes_minimo;

int k, lsc1i, lsc1f, lsc2i, lsc2f;

/* Melhores resultados */
double blsc1, blsa1, blsc2, blsa2, bmS, bmL,
      bhL, bhS, btes, btmaf;

/*****
/* Parâmetros de entrada -- ALTERAR SOMENTE AQUI */
*****/

/* Valores para c0, hS e gama */
double c0_vet[] = {0.5, 1.0, 1.5, 2.0, 3.0, 4.0},
      hS_vet[] = {0.1, 0.25, 0.5},
      gama_vet[] = {1.5, 2.0, 3.0};
/* Discretização de mS */
double mS_inicial = 0.1;
double mS_final   = 0.9;
double mS_inc     = 0.01;
/* Discretização de mL */
double mL_inicial = 1.05;
double mL_final   = 5.00;
double mL_inc     = 0.01;

/*****
/* FIM Parâmetros de entrada -- NAO ALTERAR MAIS */
*****/

int num_c0 = sizeof(c0_vet)/sizeof(double);
int num_hS = sizeof(hS_vet)/sizeof(double);
int num_gama = sizeof(gama_vet)/sizeof(double);
int num_mS = 1 + (int)((mS_final - mS_inicial) / mS_inc);
int num_mL = 1 + (int)((mL_final - mL_inicial) / mL_inc);

t = time(NULL);
print("Inicio: %s\n\n", ctime(&t));
```

```
init_fat();

for(int i_c0=0; i_c0 < num_c0; i_c0++) // c0
{
    print(" C0  GAMA*  LSC1  LSA1  LSC2  LSA2  MS"
          " ML   HS      HL   TES   TMAF\n");
    print("-----"
          " -----"
          " -----"
          " -----"
          " -----\n");

    c0 = c0_vet[i_c0];

    for(int i_hS=0; i_hS<num_hS; i_hS++) // hS
    {
        hS = hS_vet[i_hS];

        for(int i_gama=0; i_gama<num_gama; i_gama++) // gama
        {
            gama = gama_vet[i_gama];

            lscfp = 0.6195 + 1.00523*c0 + 2.983*sqrt(c0);
            k      = (int)floor(lscfp);
            alfa   = 1.0 - dpoidf(k, c0);
            tmaffp = 1.0 / alfa;
            beta   = dpoidf(k, c0*gama);
            tesfp  = 1.0 / (1-beta) - 0.5;

            tes_minimo = tesfp;

            for(int i_mS=0; i_mS<num_mS; i_mS++) // mS
            {
                mS = mS_inicial + i_mS*mS_inc;

                lscli =
                    (int)(0.6195+1.00523*mS*c0+2.983*sqrt(mS*c0));
                lsc1f = lscli+2;

                for(int i_lsc1=lscli; i_lsc1<=lsc1f; i_lsc1++) //
                    lsc1
                {
                    lsc1 = i_lsc1 + 0.5;
```

```
for(int i_lsa1=0; i_lsa1<i_lsc1; i_lsa1++) //
    lsa1
{
    lsa1 = i_lsa1 + 0.5;

    for(int i_mL=0; i_mL<num_mL; i_mL++) // mL
    {
        mL = mL_inicial + i_mL*mL_inc;

        lsc2f = (int)(0.6195+1.00523*mL*c0+2.983 *
                    sqrt(mL*c0));
        lsc2i = (int)((0.6195+1.00523*mL*c0+2.983 *
                    sqrt(mL*c0))-c0);

        for(int i_lsc2=lsc2i; i_lsc2<=lsc2f;
            i_lsc2++)
        {
            lsc2 = i_lsc2 + 0.5;

            for(int i_lsa2=0; i_lsa2<i_lsc2;
                i_lsa2++)//lsa2
            {
                lsa2 = i_lsa2 + 0.5;

                if(!calcula(c0, gama, lsc1, lsa1, lsc2,
                    lsa2, mS, mL, hS, &hL, &vtes, &vtmaf,
                    tmaffp))
                    continue;

                // Se resultado obtido melhor que o mínimo
                // até agora, armazena o resultado
                if(vtes < tes_minimo)
                {
                    blsc1 = lsc1;
                    blsc2 = lsc2;
                    blsa1 = lsa1;
                    blsa2 = lsa2;
                    bhL = hL;
                    bhS = hS;
                    bmS = mS;
                    bmL = mL;
```

```
        btes = vtes;
        btmaf = vtmaf;
        tes_minimo = vtes;
    }
    } // lsa2
    } // lsc2
    } // mL
    } // lsa1
    } // lsc1
} // mS

// Imprime o melhor resultado encontrado para estes
// valores de c0, hS e gama
print("%5.2f %6.2f %7.2f %7.2f %7.2f %7.2f %7.3f
      %7.3f %7.3f %7.3f %7.2f %7.1f\n",
      c0, gama, blsc1, blsa1, blsc2, blsa2, bmS, bmL,
      bhS, bhL, btes, btmaf);

    } // gama
    print("\n");
    } // hS
    print("\n");
} // c0

t = time(NULL);
print("\n\nFim: %s\n\n", ctime(&t));
fclose(fout);
}
```

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Apêndice B: Programa para Simulação dos Tempos até o Sinal obtidos pelo Esquema VSI EWMA

O programa em linguagem C, mostrado a seguir, implementa as simulações para obtenção da média e do desvio-padrão do tempo até o sinal, usando como dados de entrada os valores de c_0 , c_1 , λ , LSA, LSC, h_S e h_L de esquemas VSI EWMA.

```
#include <assert.h>
#include <time.h>
#include <stdio.h>
#include <stdlib.h>
#include <math.h>

#include "random.h"

long  SimulationSeed = 0;
long  N1;

    /******
    /*  Parâmetros de entrada -- ALTERAR SOMENTE AQUI          */
    /******

long  RUNS    = 100000;

double LAMBDA = 0.80;
long  N2      = 10000;

double C0    = 2.00;
double C1    = 3.00;    /* gama = C1/C0 */

double LSA   = 2.08;
double LSC   = 5.79;

double HL    = 1.345;
double HS    = 0.50;

    /******
    /*  FIM Parâmetros de entrada -- NAO ALTERAR MAIS          */
    /******
```

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Programa para Simulação dos Tempos até o Sinal obtidos pelo Esquema VSI EWMA 135

```
double CalculaTs(double* Z, double* h)
{
    long    i, j;
    long    c;
    double  Tsf = 0.0;
    double  Q;
    int     serie_invalida = 1;

    /* Gerar N1 valores de Ci, onde C ~ Poisson(C0) e calcular Zi;
       i = 1, ..., N1; se algum Zi > LSC, descarta a série
    */
    Z[0] = C0;
    while(serie_invalida)
    {
        serie_invalida = 0;
        for(i=1; i<=N1; i++)
        {
            c = RandomPoisson(C0);
            Z[i] = (1.0-LAMBDA)*Z[i-1] + LAMBDA*c;
            if(Z[i] > LSC)
                serie_invalida = 1;
        }
    }

    /* Cálculo de h[1..N1] */
    for(j=1; j<=N1; j++)
    {
        if (Z[j-1] <= LSA)
            h[j] = HL;
        else
            h[j] = HS;
    }

    /* Gerar valores de Ci agora com Ci ~ Poisson (C1) e calcular
       Zi; i = N1+1 até sinal; até ocorrência de um sinal (ou até
       gerar N2 valores); Calcula também os valores para hi
       correspondentes
    */
    for(j=N1+1; j<=N1+N2; j++)
    {
        c = RandomPoisson(C1);
        Z[j] = (1.0-LAMBDA)*Z[j-1] + LAMBDA*c;
        if(Z[j-1] <= LSA)
            h[j] = HL;
        else if(Z[j-1] < LSC)
            h[j] = HS;
        else
        {
            h[j] = -1; /* SINAL */
            break;
        }
    }

    printf("%d - ", j-N1);
    assert(j<=N1+N2); /* Passou pelos N2 elementos sem sinal */

    /* Cálculo de TSf*/
    for(j=N1+1; h[j]!=-1; j++)
        Tsf += h[j];

    /* Calcula Ts */
    Q = RandomUniform() * h[N1+1];
}
```

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Programa para Simulação dos Tempos até o Sinal obtidos pelo Esquema VSI EWMA 136

```
    return Tsf - Q;
}

void RunSimulation(void)
{
    double* Z;
    double* h;
    double* Ts;
    long    i;
    double  mean;
    double  stddev;

    N1 = (long)ceil(log(0.01)/(2*log(1.0-LAMBDA)));

    /* Aloca vetores (vetores com um elemento a mais para serem
       indexados a partir de 1 e não de 0)
    */
    Z = (double*)malloc((N1+N2+1)*sizeof(double));
    assert(Z);

    h = (double*)malloc((N1+N2+1)*sizeof(double));
    assert(h);

    Ts = (double*)malloc((RUNS+1)*sizeof(double));
    assert(Ts);

    /* Loop de simulação para cálculo de TS */
    for(i=1; i<=RUNS; i++)
    {
        Ts[i] = CalculaTs(Z, h);
        printf("Ts[%d] = %g\n", i, Ts[i]);
    }

    /* Calcula média e desvio-padrão dos TS */
    mean = 0.0;
    for(i=1; i<=RUNS; i++)
        mean += Ts[i];
    mean = mean / RUNS;

    stddev = 0.0;
    for(i=1; i<=RUNS; i++)
        stddev += (Ts[i]-mean)*(Ts[i]-mean);
    stddev = sqrt(stddev / RUNS);

    printf("N1 = %d\n",N1);
    printf("Média = %g\n", mean);
    printf("Desvio-padrão = %g\n", stddev);

    /* Libera memória */
    free(Z);
    free(h);
    free(Ts);
}

int main(int argc, char* argv[])
{
    time_t  t;

    RandomInit(SimulationSeed);
```

9 Apêndice B:

Programa para Simulação dos Tempos até o Sinal obtidos pelo Esquema VSI EWMA 137

```
t = time(NULL);
printf("Início da Simulação: %s\n", ctime(&t));

RunSimulation();

t = time(NULL);
printf("Final da Simulação: %s\n\n", ctime(&t));

system("pause");
return 0;
}
```