

**MELISSA**

**Predictive Control of biomass production  
in the Rhodobacter compartment of MELISSA**

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## NOTATIONS

$cE_b$  :  $E_b$  set point ( $W.m^{-2}$ )

$cF_R$  :  $F_R$  set point ( $W.m^{-2}$ )

$C_x$  : biomass concentration ( $kg.m^{-3}$ )

$cp_i$  : level  $i$  ( $i = 1$  or  $2$ ) production rate set point ( $g/h$ )

$cq_i$  : level  $i$  ( $i = 1$  or  $2$ ) flow rate set point ( $l/h$ )

$dil$  : dilution rate ( $h^{-1}$ )

$dq$  : flow rate ratio (dimension less)

$dt$  : control time period (h)

$E_b$  : total radiant energy absorbed by the sphere photometer ( $W.m^{-2}$ )

$F_R$  : mean incident radiant light energy flux ( $W.m^{-2}$ )

$h_c$  : coincidence horizon (h)

$p$  : biomass production rate ( $g/h$ )

$q$  : flow rate ( $l/h$ )

$TR$  : closed loop time response (h)

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# 1. INTRODUCTION

The aim of this note is to test a control law to be applied to the production of biomass in the compartment II (rhodospirillum rubrum photobioreactor) of the MELISSA project.

The process simulator is based on the model for growth of rhodospirillum rubrum under light limitation in photobioreactors (developed at LGCB in Clermont-Ferrand by J.F. Cornet ; TN 23.4).

The control law is based on the above-mentioned model and widely inspired from the one of the compartment IV (developed at ADERSA by N. Fulget ; TN 24.1) the law structure is the same, only parameters are different.

The dimensions of the simulated reactor are those of compartment IV (Spirulina photobioreactor).

The studied system (fig.1) is composed of :

- the process (photobioreactor) whose inputs are the flow rate  $q$  and the light flux  $F_R$ , and whose outputs are the concentration biomass  $C_x$  and the production  $p$  ;
- the systems A and B (detailed in fig. 2 and 3) which include, each of them, a level 0 controller associated to an actuator (valve or volumetric pump for system A ; light energy intensity for system B) and to a sensor (flow meter for system A ; integrating sphere photometer for system B). These systems are supposed to have short dynamics compared to those of the process and the level 0 controllers are supposed perfect. So they are reduced to identity function, in the simulator :

$$\begin{aligned} q &= cq1 && \text{at each instant } t \\ F_R &= cF_R \end{aligned}$$

- the level 1 controller which works on the PFC (Predictive Functional Control) principles and computes the light flux  $F_R$  to control the production  $p$  versus the level 1 production set point  $cp1$  ;
- the level 2 controller which computes the level 1 flow rate and production set points,  $cq1$  and  $cp1$ , versus the corresponding level 2 set points,  $cq2$  and  $cp2$ , taking into account the max and min concentration constraints and the flow variation  $dq$  ( $\pm 10\%$  of  $cq2$ ). The set points  $cq2$  and  $cp2$  are fixed by the operator.

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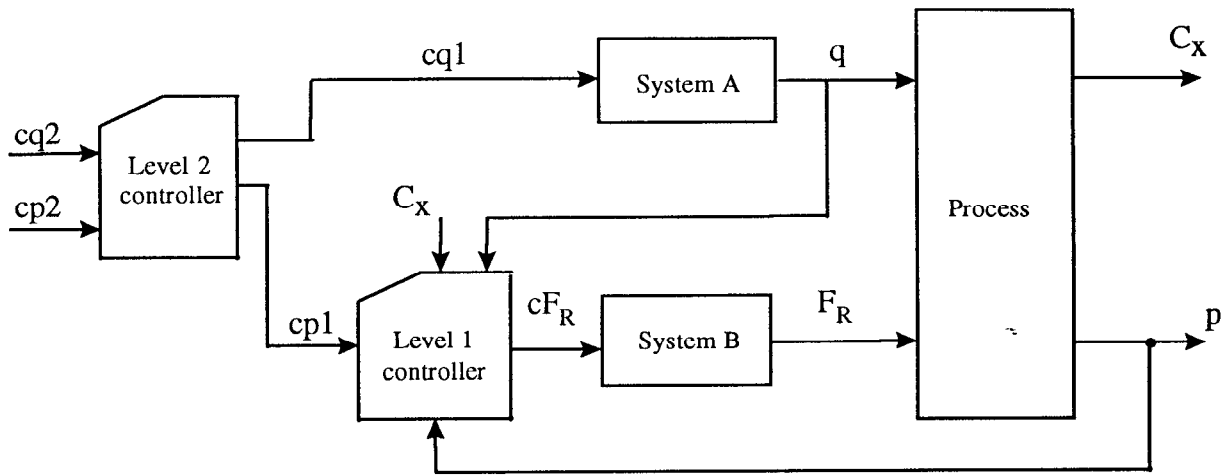


Fig 1 : Global scheme processus and levels 0 to 2 controllers

The systems A and B which contain level 0 controllers (generally PID) are detailed hereafter.

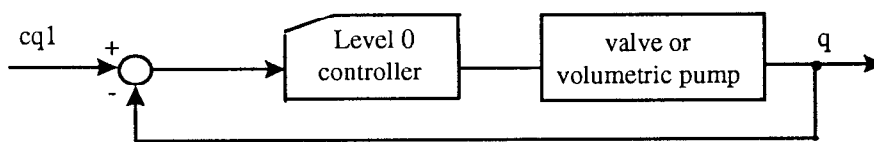


Figure 2 : System A

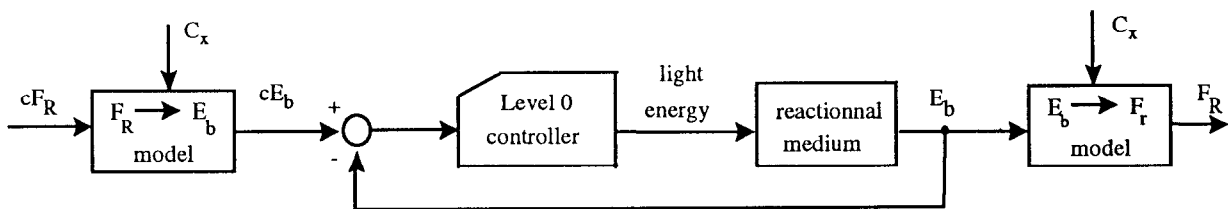


Figure 3 : System B

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## 2. CONTROL STRUCTURE

Full details are given in TN 24.1 about the levels 1 and 2 of the control. Here are a few recalls.

### 2.1. Level 1 control

This controller is built on the principles of PFC (Predictive Functional Control) :

- internal model ;
- reference trajectory to specify the future closed loop behavior ;
- manipulated variable structuration ;
- modeling error extrapolation.

A scenario strategy is used to adapt this control law to the present process which is non linear : several (generally 2) input protocols are applied to the model from which a prediction of the model output is inferred. The computation involves the integration of a differential equation : the Euler integration method is quite convenient for this purpose.

### 2.2. Level 2 control

This level checks the compatibility flow rate and production set points, fixed by the operator (or a upper level), with regard to variation flow rate and concentration constraints.

Algorithm : at each period of time, the following computations are made :

1 - maximum and minimum flow rate

$$\begin{aligned}q_{\max} &= cq2 (1 + dq) \\q_{\min} &= cq2 (1 - dq)\end{aligned}$$

2 - maximum and minimum production

$$\begin{aligned}p_{\max} &= q_{\max} * C_{x_{\max}} \\p_{\min} &= q_{\min} * C_{x_{\min}}\end{aligned}$$

3 - level 1 production set point

$$cp1 = \max [p_{\min}, \min (p_{\max}, cp2)]$$

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4 - level 1 flow rate set point

if  $\left(\frac{cp1}{C_{x\_max}} > cq2\right)$  then

$$cq1 = \min\left(q\_max, \frac{cp1}{C_{x\_max}}\right)$$

else if  $\left(\frac{cp1}{C_{x\_min}} < cq2\right)$  then

$$cq1 = \max\left(q\_min, \frac{cp1}{C_{x\_min}}\right)$$

else

$$cq1 = cq2$$

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### 3. SIMULATION RESULTS

The present control structure works on the compartment IV (Spirulina biophotoreactor) and has already been validated. The aim is now to check its extension to the compartment II and to tune the specific parameters.

#### 3.1. Control parameter tuning

- control period :  $dt = 0,5$  hour
- closed loop response time :  $TR = 12h$
- coincidence point :  $h_c = 2.5$  hours ( $n_{hc}=5$  control periods)

#### 3.2. Parameters process

They are summed up in the file comnl2.h.

Their values are not exactly those of the TN 23.4 (CORNET J.-F., 1996. Model parameters for growth of rhodospirillum rubrum under light limitation in photobioreactors) because the present simulations have been done before its issue. Here are the differences :

Parameter	Present study	TN 23.4	Unit
Ea	600	220	$m^2/kg$
Es	1330	480	$m^2/kg$

#### 3.3. Tests results

The figures 4 to 8 show the simulations whose presentation is described in the following table.

Graph	Signal name	Type of line	Unit
1 (top of figure)	production • measure • level 2 setpoint • level 1 set point	continuous - dashed	g/h
2	light energy flux $F_R$	continuous	$W/m^2$
3	biomass concentration • measure • minimum or maximum constraint	continuous -.	g/l
4	flow rate • measure • level 2 set point	continuous -.	l/h

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A white noise (amplitude : 0.02 g/l) is added to the concentration.

- **First test (figure 4)**

The dilution rate is fixed to  $0.01 \text{ h}^{-1}$  ( $cq2 = 0.07 \text{ l/h}$ ). At time  $t=20\text{h}$ , a step (from  $0.035 \text{ g/h}$  to  $0.045 \text{ g/h}$ ) is applied to the level 2 production set point  $cp2$ . As the concentration and the flow rate do not over run their limits, the level 2 control calculates the level 1 set points equal to the level 2 set points, all over the test :

$$\begin{aligned} cq1 &= cq2 \\ cp1 &= cp2 \end{aligned}$$

Immediately after the  $cp1$  step, the max constraint of  $F_R$  is reached and slows the dynamics of the production. The time response (30 h) depends on the amplitude step of  $cp1$  and on its sense (positive or negative).

- **Second test (figure 5)**

The test is nearly the same as in the previous figure, except that the  $cp2$  step is negative, from  $0.045$  to  $0.035 \text{ g/h}$ . The production, which is not null even when  $F_R$  is minimum ( $10 \text{ W/m}^2$ ), goes down only because of dilution. The response time is 60h.

These two tests (fig 4 and 5) illustrate the high non linear aspect of the process.

- **Third test (figure 6)**

The dilution is higher :  $0.02 \text{ h}^{-1}$  ( $cq2 = 0.14 \text{ l/h}$ ). The result is opposite to the one of the two previous tests : the response time is higher for the positive step (at  $t=20\text{h}$ ) than for the negative one (at  $t=70\text{h}$ ), 30h and 20h, respectively. It is another illustration of the non linear aspect of the process.

- **Fourth test (figure 7)**

The dilution rate is lower : its initial value is  $0.004 \text{ h}^{-1}$  ( $cq2 = 0.028 \text{ l/h}$ ). At time  $t=20\text{h}$ , a production set point step  $cp2$  is applied from  $0.04 \text{ g/h}$  to  $0.06 \text{ g/h}$ .

The level 2 control detects an incompatibility between level 2 set points and the constraints. First it computes a level 1 production set point  $cp1$ , taking into account the maximum flow rate and the maximum concentration (graph 3) :  $cp1$  varies from  $0.04 \text{ g/h}$  to  $0.0308 * 1.5 = 0.0462 \text{ g/h}$  (graph 1). Then it allows a 10% increase of flow rate with regard to  $cq2$  : the flow rate  $q$  rises from  $0.028 \text{ l/h}$  to  $0.0308 \text{ l/h}$  (graph 4).

The dynamics of the production is quick : the response time is less than 10 h.

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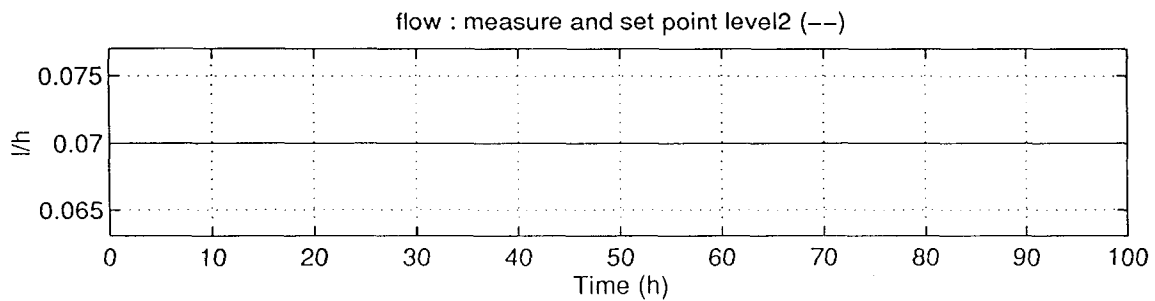
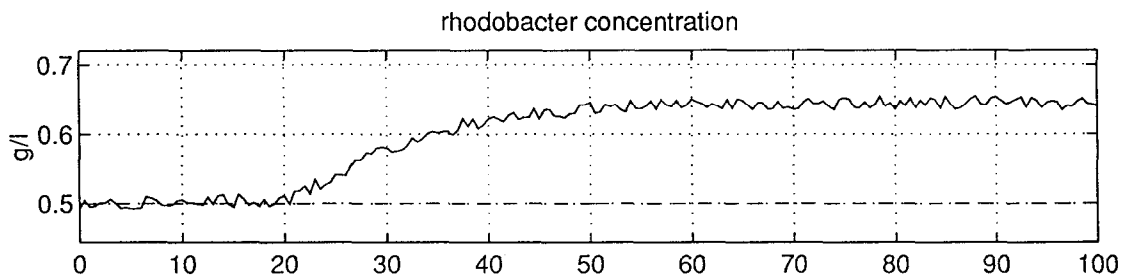
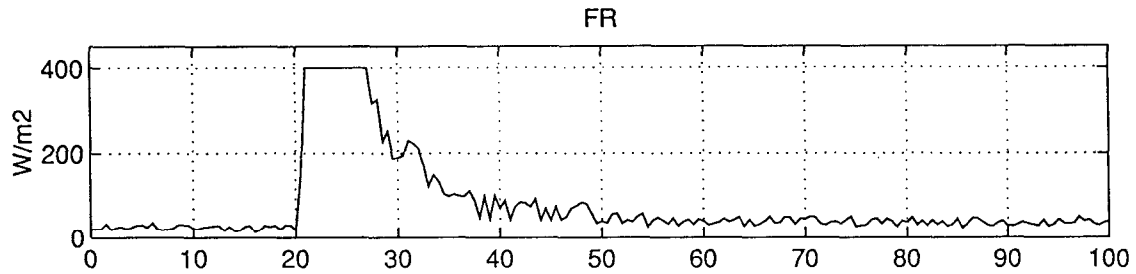
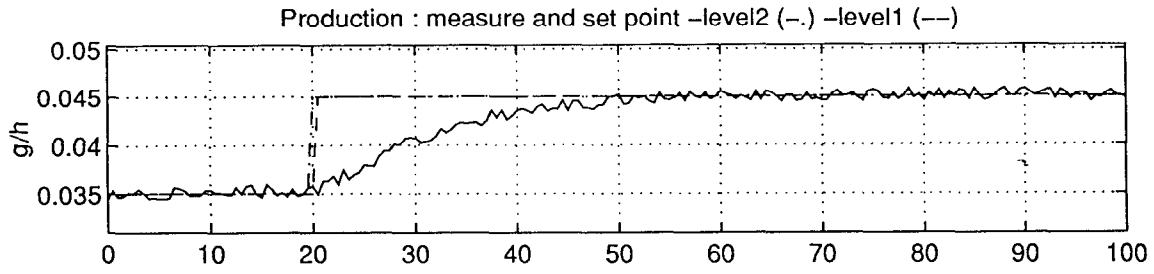
- **Fifth test (figure 8)**

A step is applied to cp2 at time  $t=20h$  and then a step is applied to cq2 at time  $t=60h$ .

At time  $t=20h$ , all happens as in the test of figure 7 : the level 2 control computer cp1 versus the maximum concentration and raises the flow rate  $q$  with a 10% increase.

At time  $t=60h$ , the level 2 control sets the flow rate  $q$  to 0.04 l/h so that cp1 remains equal to cp2 ; the concentration stays at its maximum value : 1.5 g/l.

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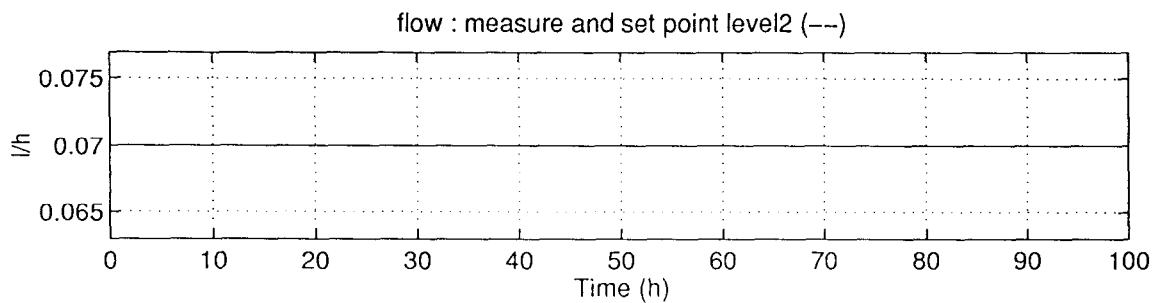
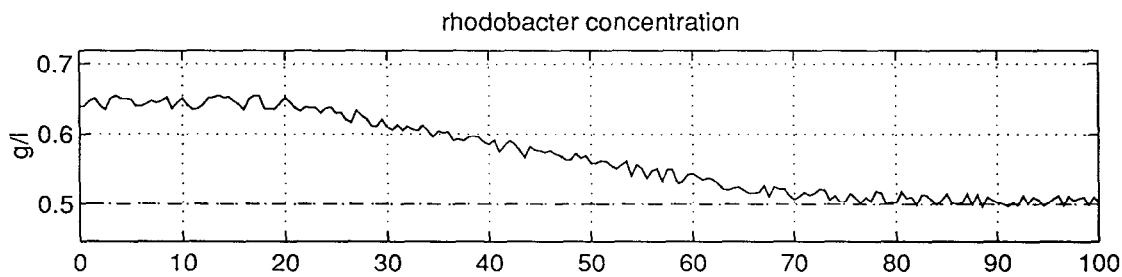
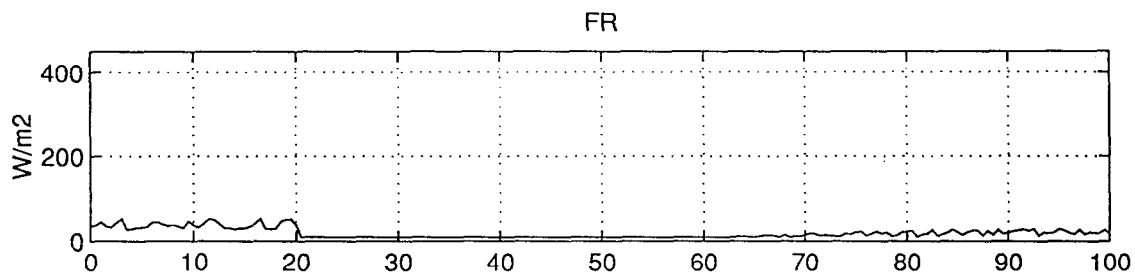
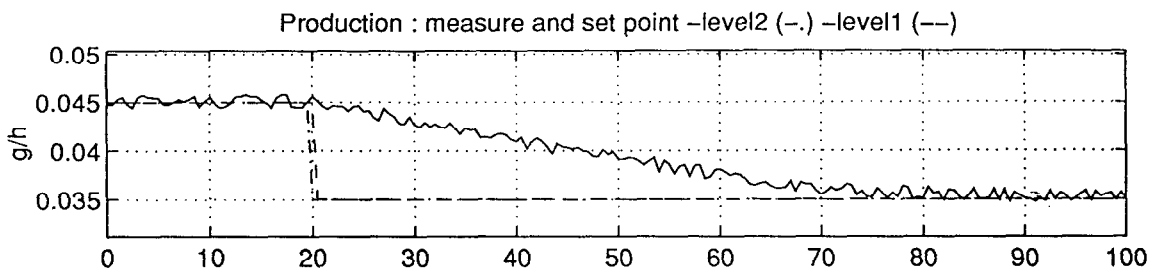


Production control

dil = .01 (1/h)

Figure 4

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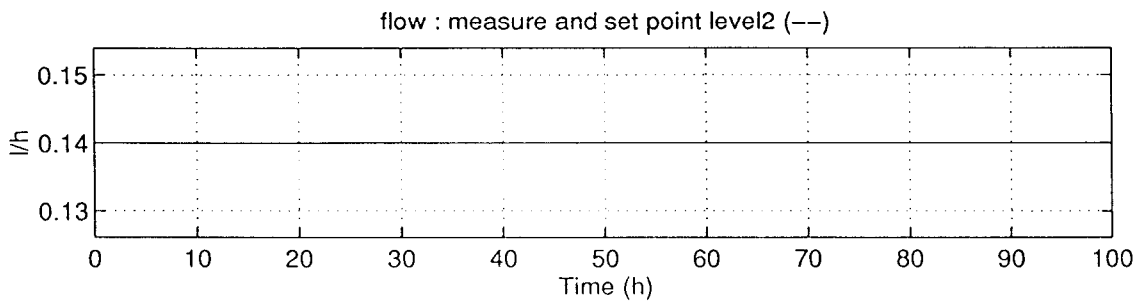
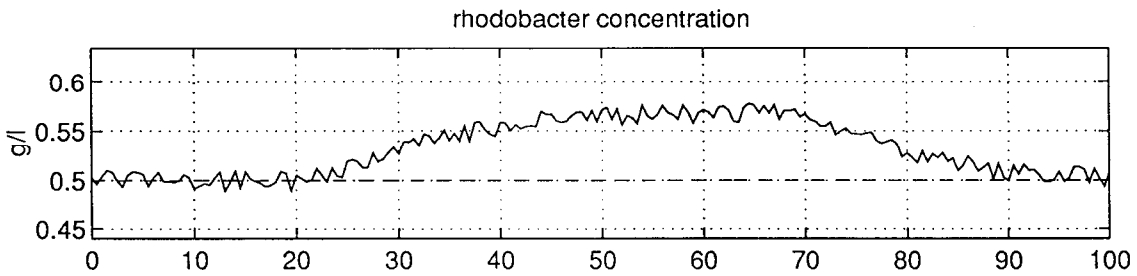
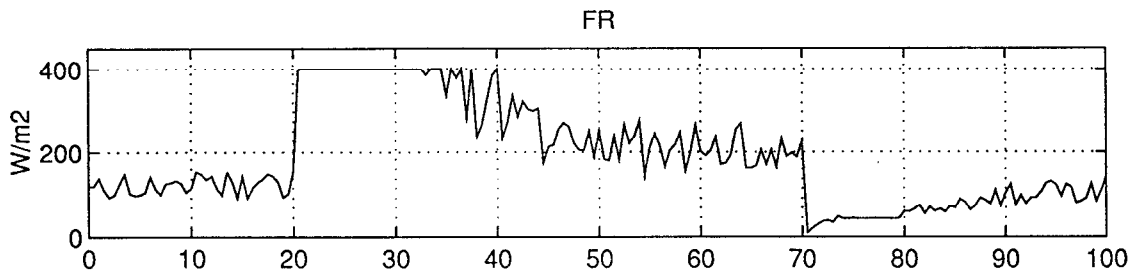
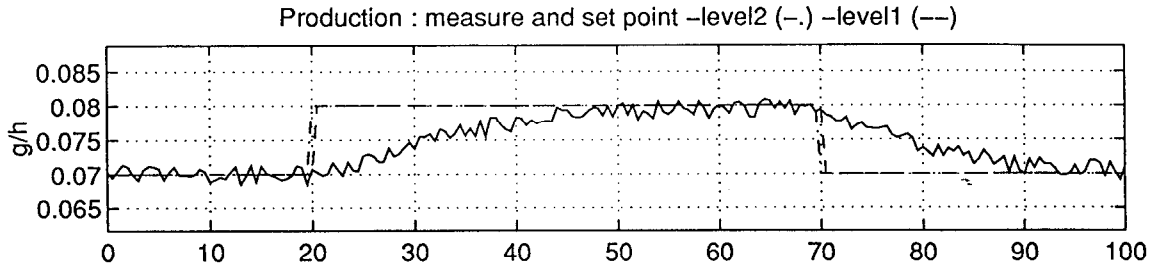


Production control

dil = .01 (1/h)

Figure 5

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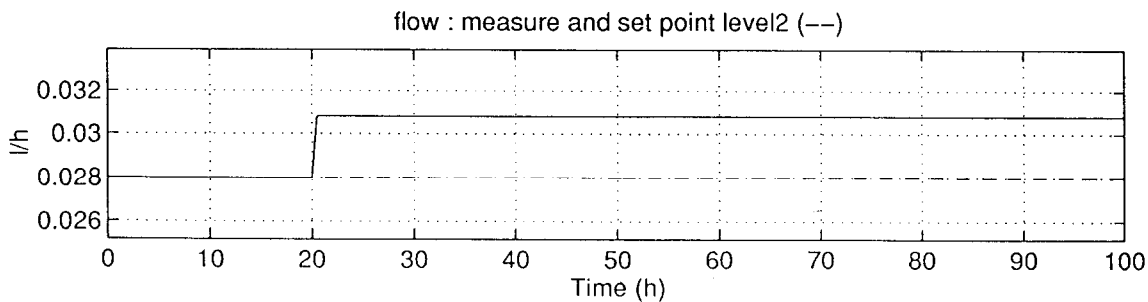
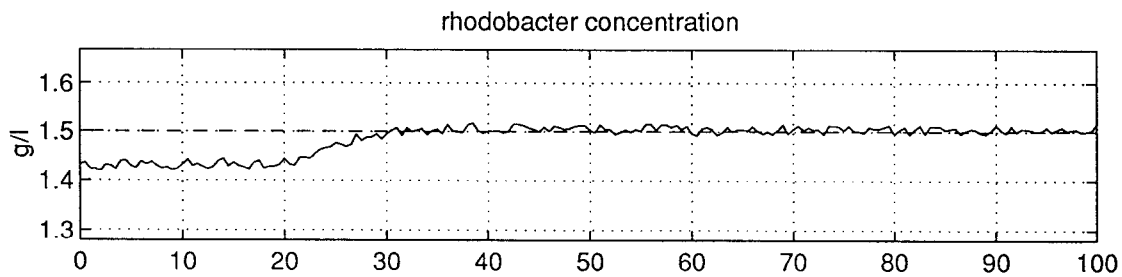
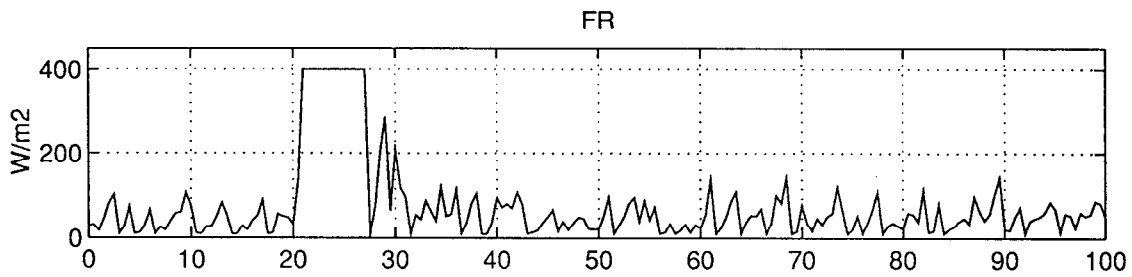
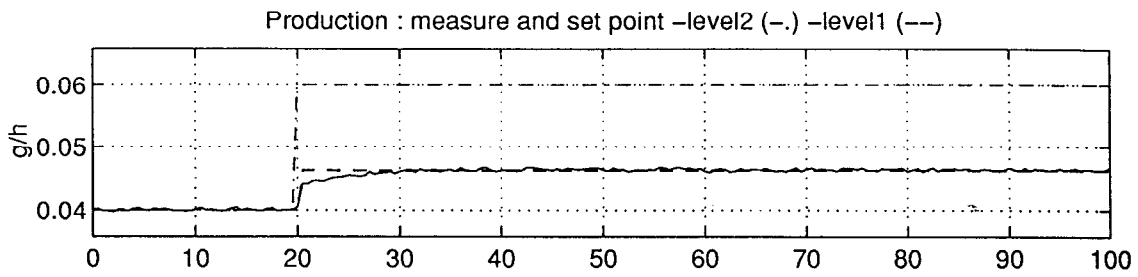


Production control

dil = .02 (1/h)

Figure 6

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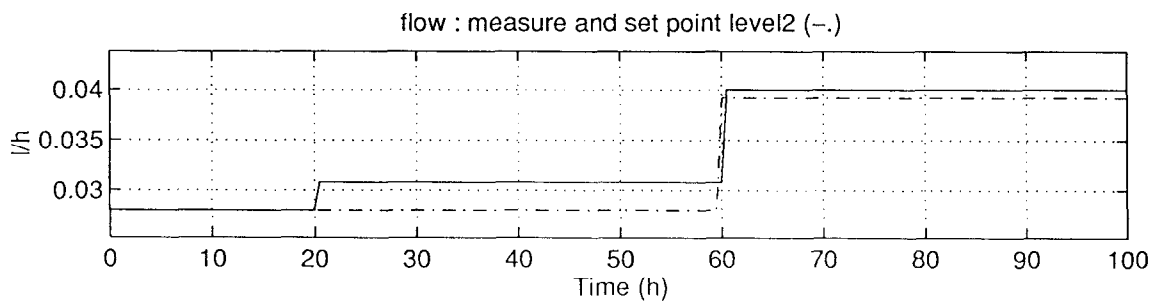
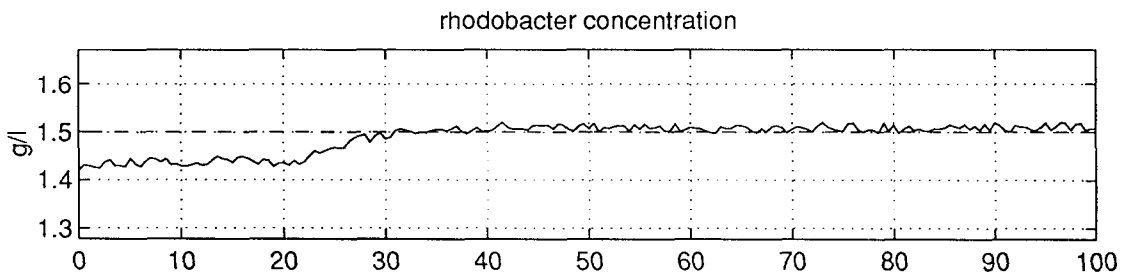
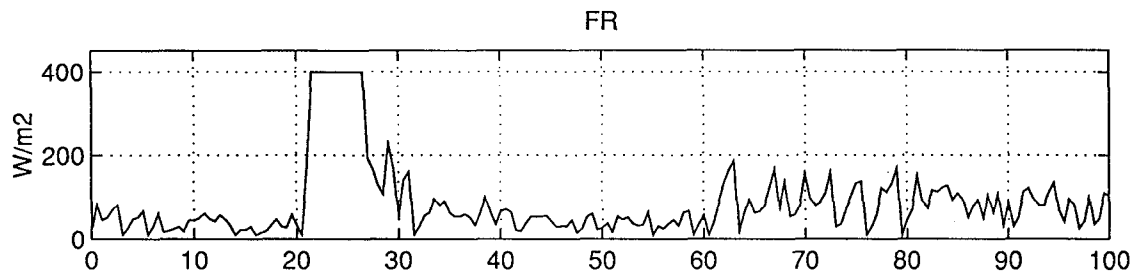
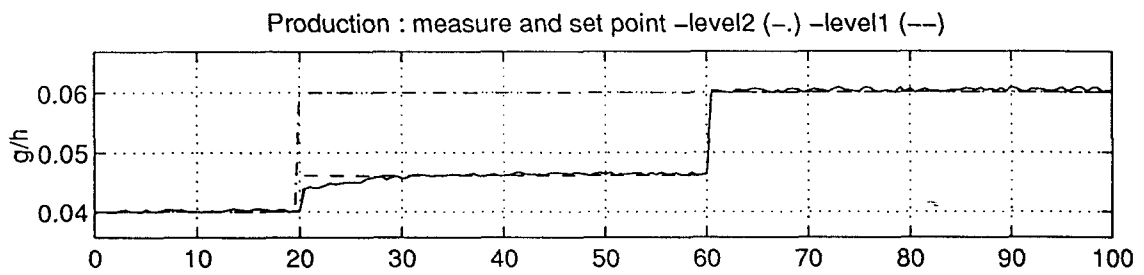


Production control

dil = .004 (1/h)

Figure 7

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Production control

dil variation at time = 60 : from .004 to .0056 (1/h)

Figure 8

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#### 4. CONCLUSION

The present control law based on the PFC principles is checked in simulation, and is quite convenient to the rhodospirillum rubrum photobioreactor, highly non linear process.

It can now be validated on the real process.

The parameters of the regulator are summed up in the file 'comnl2.h' in annex (do not forget to give to  $E_a$  and  $E_s$  their right values according to the Technical Note 23.4).

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## ANNEX

- scomnl.m : main control programme (Matlab® language)
- mx\_comnl2.c : interface programme from Matlab® to C language
- comnl2.c : levels 2 and 1 control routine (C language)
- funcalc.c : specific calculation functions
- comnl2.h : control model parameters
- proto2.h : declaration of functions

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```
function [sys,x0]=scomnl(tn,x,u,flag,dt,vol_rhodo,prod0,qe0)

% SCOMNL      S-Function d'un regulateur PFC (utilisation avec SIMULINK).
%
% Synopsis
%   [sys,x0]=scomnl(tn,x,u,flag,dt,vol_rhodo,prod0,qe0)
%
% Parameters
%   dt      periode de commande.
%   vol_rhodo  volume du reacteur 2 (rhodobacter)
%   prod0   consigne initiale de production niveau 2
%   qe0     consigne initiale de debit niveau 2
%
% Entrees
%   u(1)    consigne production niveau 2
%   u(2)    consigne debit niveau 2
%   u(3)    concentration mesuree
%   u(4)    flux lumineux mesure
%   u(5)    debit mesure
%
% Sorties
%   s(1)    FR
%   s(2)    consigne production niveau 1
%   s(3)    consigne debit niveau 1
%
%   L'utilisation de cette fonction suppose l'existence du Mex-File
%   MX_COMNL2
%
% ADERSA Imm. 6 15-DEC-94
%   modifie le 5-JUIL-96 par JJL

%> Taille des parametres et Conditions initiales -----

global flinit Fr0

if flag==0,
    sys = [
        0      % continuous states
        3      % discrete states
        3      % outputs
        5      % inputs
        0      % discontinuous ...
        0      % direct feedthrough
    ];

%   Calcul de la valeur stationnaire de Fr
%   -----
%   Initialisation de la precision relative
    prec = 1e-2;
%   Initialisation du point de depart de la procedure iterative
    c0 = prod0 / qe0;
    xx0 = 40;
    yy0 = prod0 / vol_rhodo;
    epsilon = yy0 * prec;
    deltax = 10;
%   Calcul par procedure iterative
    Fr0 = caliter(c0, xx0, yy0, deltax, epsilon);
    if Fr0 == Inf
        Fr0 = 10;
    end

    x0 = [prod0 Fr0 qe0];
    flinit=0;
```

```
%> Etats Discrets x(n+1) -----
elseif abs(flag)==2,
  if abs( round(tn/dt)-(tn/dt) ) < sqrt(eps),
    %> Entree
    consp = u(1);
    consq = u(2);
    cx    = u(3);
    frmes = u(4);
    qemes = u(5);

    %> Calcul de la commande
    % Dans l'etat actuel du regulateur, l'argument 'flinit' de 'mx_comnl2'
    % est inutile.
    if flinit == 0 ,
      frmes = Fr0;
      [fr, cons_p1, cons_q1] = mx_comnl2( consp , cx , consq , ...
                                         frmes , qemes , flinit);
      flinit=1;
    else
      [fr, cons_p1, cons_q1] = mx_comnl2( consp , cx , consq , ...
                                         frmes , qemes , flinit);
    end

    %> Sortie de la fonction: vecteur d'etat x
    sys = [cons_p1 fr cons_q1];

  else
    sys = x;
  end

%> Sorties du systeme -----
elseif flag==3,
  sys = x;

%> Instant du prochain appel -----
elseif flag==4,
  ns = tn/dt;           % nombre de simulation
  sys = dt * (1 + floor(ns + 1e-13*(1+ns)));

%> -----
else
  sys = [];
end
```

## mx\_comnl2.c

```
/* --- MX_COMNL2.C -----  
Fonction:  
    mex_fonction permettant le calcul de la commande d'un regulateur PFC  
    pour le compartiment 2 (Rhodobacter)  
Synopsis:  
    [FR,CONS_PROD1,CONS_QE1]=MX_COMNL2(CONS_PROD2,CX,QE2,FRMES,QEMES,FLAG)  
Description:  
    Calcul des consignes de production et debit de niveau 1, CONS_PROD1  
    et CONS_QE1, a partir des consignes recues par le niveau 2, CONS_PROD2  
    et QE2  
    Calcul la commande courante FR a partir de  
    la mesure de concentration CX, de la consigne CONS_PROD1, du  
    debit mesure QEMES et du flux lumineux estime FRMES  
*/  
  
/* modifie le 5-JUIL-96 par JJJ */  
  
#include "mex.h"  
#include "comnl2.h"  
#include "proto2.h"  
  
/* Arguments d'entree/sortie */  
#define PC_IN    prhs[0]  
#define YP_IN    prhs[1]  
#define QC_IN    prhs[2]  
#define FR_IN    prhs[3]  
#define QM_IN    prhs[4]  
#define FL_IN    prhs[5]  
#define FR_OUT   plhs[0]  
#define CONS_P_OUT plhs[1]  
#define CONS_Q_OUT plhs[2]  
  
mexFunction(nlhs, plhs, nrhs, prhs)  
int nlhs, nrhs;  
Matrix *plhs[], *prhs[];  
{  
    double    *yp , *c , *d , *fr , *frmes , *qemes , *fl , *cons_p1 , *cons_q1 ;  
    double varsortie[3];  
  
    /* Test sur le nombre d'arguments */  
    if (nrhs<6 | nrhs>6)  
    {  
        mexErrMsgTxt("MX_COMNL2 a besoin de 6 arguments d'entree.");  
    }  
    else if (nlhs <3 | nlhs>3)  
    {  
        mexErrMsgTxt("MX_COMNL2 a besoin de 3 arguments de sortie.");  
    }  
  
    /* Creation de matrice pour les arguments de sortie */  
    FR_OUT = mxCreateFull( 1 , 1 , REAL );  
    CONS_P_OUT = mxCreateFull( 1 , 1 , REAL );  
    CONS_Q_OUT = mxCreateFull( 1 , 1 , REAL );  
  
    /* Affectation des pointeurs pour les parametres */  
    fr = mxGetPr( FR_OUT );  
    cons_p1 = mxGetPr( CONS_P_OUT );  
    cons_q1 = mxGetPr( CONS_Q_OUT );  
    yp = mxGetPr( YP_IN );  
    c = mxGetPr( PC_IN );  
    d = mxGetPr( QC_IN );  
    frmes = mxGetPr( FR_IN );  
    qemes = mxGetPr( QM_IN );  
    fl = mxGetPr( FL_IN );
```

```
/* Appel de la fonction COMNL2 */  
comnl2 ( *c , *yp , *d , *frmes , *qemes , *fl , varsortie );  
*fr = varsortie[0];  
*cons_p1 = varsortie[1];  
*cons_q1 = varsortie[2];  
}
```

## comnl2.c

```
#include "comnl2.h"
#include "math.h"
#include "proto2.h"

/*
  COMNL2.C  Algorithme du regulateur non lineaire.
  ESA - MELISSA - RHODOBACTER

  Date creation pour compartiment Spiruline : 15-DEC-94
  Modifie le 5-JUIL-96 par JJL
*/

/* --- COMNL2 -----
Fonction:
  Equations du regulateur non lineaire du compartiment 2 (Rhodobacter)
Synopsis:
  COMNL2 (CONS_PROD2, CX, CONS_QE2, FRMES, QEMES, FR, CONS_PROD1, CONS_QE1)
Description:
  Calcul des consignes de production et debit de niveau 1, CONS_PROD1
  et CONS_QE1, a partir des consignes recues par le niveau 2, CONS_PROD2
  et CONS_QE2
  Calcul la commande courante FR a partir de
  la mesure de concentration CX, de la consigne CONS_PROD1, des mesures
  de debit QEMES et du flux lumineux FRMES
*/
void comnl2(cons_prod2, cx, cons_qe2, frmes, qemes, indic, varsortie)

double  cons_prod2, cx, cons_qe2, frmes, qemes, indic , varsortie[3];
{
/* declaration des variables internes */
double  prod, dil, prod_ref, delfr;
double  fr , fr1, fr2, prod1, prod2;
double  qe_max , qe_min , prod_max , prod_min ;
double  cons_prod1 , cons_qe1;

  prod = cx*cons_qe2;

/* algorithme de commande de niveau 2 */
qe_max = cons_qe2*(1+dq);
qe_min = cons_qe2*(1-dq);
prod_max = qe_max*cx_max;
prod_min = qe_min*cx_min;
cons_prod1 = max(prod_min, min(prod_max, cons_prod2));
cons_qe1 = cons_qe2;
if (cons_prod1/cx_max > cons_qe2 )
  {
  cons_qe1 = min(qe_max, cons_prod1/cx_max);
  }
if (cons_prod1/cx_min < cons_qe2 )
  {
  cons_qe1 = max(qe_min, cons_prod1/cx_min);
  }

  dil = cons_qe1/vol;
  prod = cx * qemes;

/* trajectoire de reference */
  prod_ref = cons_prod1 - pow(lambda, nhc)*(cons_prod1 - prod);

/* commande precedente mesuree */
  fr = frmes;

/* premier scenario */
```

## comnl2.c

```
fr1 = fr;
prod1 = model2(cx,fr1,dil);

/* deuxieme scenario */
delfr = dfr*sign(cons_prod1 - prod);
fr2 = fr1 + delfr;
prod2 = model2(cx,fr2,dil);

/* calcul de fr */
fr+ = (prod_ref - prod1)/(prod2 - prod1)*delfr;

/* contraintes sur fr */
fr = max(fr_min,min(fr_max,fr));

/* ecriture des arguments de sortie */
varsortie[0] = fr;
varsortie[1] = cons_prod1;
varsortie[2] = cons_qel;
return;
}
/* --- MODEL2 -----
Fonction:
    integration du modele
Synopsis:
    MODEL2 (CX,FR,DIL,PROD)

*/
double model2(cx,fr,dil)

double cx, fr, dil ;

{
    double v, dv, vout , prod;
    int k;

    v = cx;
    for (k=1 ; k <= nhc; k++)
    {
        dv = dercx2(v,fr,dil);
        vout =v + dt *dv;
        v=vout;
    }
    prod=vout*dil*vol;
    return (prod);
}

/* --- DERCX2 -----
Fonction:
    calcul de la derivee de cx
Synopsis:
    DERCX2 (cx,fr,dil,dvt);

*/
double dercx2(cx,fr,dil)

double cx, fr, dil;

{
    double dcdxdt;
    double alpha, delta, pij, pijz;
    double z, rx;
```



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comnl2.c

3

```
alpha = sqrt(Ea/(Ea+Es));
delta = (Ea+Es)*alpha*RT*cx;

pij = 0.;
for (z=jstep/2; z<=1-jstep/2; z+=jstep)
{
    pijz = fr/z*2*cosh(delta*z)/(cosh(delta)+alpha*sinh(delta));
    if (pijz>=Fmin)
    {
        pij+ = 2*z*pijz/(Kj+pijz)*jstep;
    }
}
rx = muM*pij*cx*wiv;

dcxdt = -dil*cx + rx ;
return (dcxdt);
}
```

```
#include "math.h"

/* --- MIN.C -----
Fonction:
    Minimum de deux valeurs.
Synopsis:
    X=MIN(Y,Z)
*/
double min( x1 , x2 )

double x1 , x2;
{
    double x;
    x = (x1 < x2) ? x1 : x2;
    return( x );
}

/* --- MAX.C -----
Fonction:
    Maximum de deux valeurs.
Synopsis:
    X=MAX(Y,Z)
*/
double max( x1 , x2 )

double x1 , x2;
{
    double x;
    x = (x1 > x2) ? x1 : x2;
    return( x );
}

/* --- SIGN.C -----
Fonction:
    signe d'une valeur.
Synopsis:
    X=SIGN(Y)
*/
double sign( y )

double y;
{
    double x;
    x = (y < 0) ? -1. : 1.;
    return( x );
}
```

```
/*
  Nom : comnl2.h

  Fonction : Coefficients de la commande du compartiment 2 (Rhodobacter)

  Date : 21-MAY-96
  Modified 5-JUL-96
*/

#define dt          0.5      /* control period (in h ) */
#define nhc         5.       /* coincidence point (in dt) */
#define lambda      0.88    /* reference trajectory dynamic */
#define dfr         5.       /* radiant flux increment (in W/m2) */
#define fr_min      10.      /* min constraint on FR (in W/m2) */
#define fr_max      400.     /* max constraint on FR (in W/m2) */
#define dq          .1      /* flow variation (dimensionless) */
#define cx_min      0.5      /* min constraint on CX (in g/l) */
#define cx_max      1.5      /* max constraint on CX (in g/l) */
#define vol         7.       /* reactor volume */

/* not used : */
/* #define zpc */           /* biotic mass fraction of phycocyanins */
/* #define zch */           /* biotic mass fraction of chlorophylles */
/* #define zg */            /* biotic mass fraction of glycogen */
/* #define za */           /* resultant biotic mass fraction of the
                           light harvesting antenna */

#define Ea          600.     /* global absorption mass coefficient (m2/kg) */
#define Es          1330.   /* global scattering mass coefficient (m2/kg) */
#define RT          .048    /* radius of the cylindrical bioreactor */
#define Kj          7.      /* half saturation constant for radiant energy
                           available (W/m2) */

#define muM         .17     /* maximum growth rate (1/h) */
#define wiv         .52     /* working illuminated volume */
#define Fmin        1.      /* minimum radiant energy flux (W/m2) */
#define jstep       .01     /* integrative dimensionless radius */
```

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proto2.h

1

```
void comn12( );  
double model2( );  
double dercx2( );  
double sign( );  
double min( );  
double max( );
```