Numerical Optimal Control of HIV Transmission in Octave/MATLAB

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Abstract: We provide easy and readable GNU Octave/MATLAB code for the simulation of mathematical models described by ordinary differential equations and for the solution of optimal control problems through Pontryagin’s maximum principle. For that, we consider a normalized HIV/AIDS transmission dynamics model based on the one proposed in our recent contribution (Silva, C.J.; Torres, D.F.M. A SICA compartmental model in epidemiology with application to HIV/AIDS in Cape Verde. *Ecol. Complex.* 2017, 30, 70–75), given by a system of four ordinary differential equations. An HIV initial value problem is solved numerically using the *ode45* GNU Octave function and three standard methods implemented by us in Octave/MATLAB: Euler method and second-order and fourth-order Runge–Kutta methods. Afterwards, a control function is introduced into the normalized HIV model and an optimal control problem is formulated, where the goal is to find the optimal HIV prevention strategy that maximizes the fraction of uninfected HIV individuals with the least HIV new infections and cost associated with the control measures. The optimal control problem is characterized analytically using the Pontryagin Maximum Principle, and the extremals are computed numerically by implementing a forward-backward fourth-order Runge–Kutta method. Complete algorithms, for both uncontrolled initial value and optimal control problems, developed under the free GNU Octave software and compatible with MATLAB are provided along the article.

Keywords: numerical algorithms; optimal control; HIV/AIDS model; GNU Octave; open source code for optimal control through Pontryagin Maximum Principle

MSC: 34K28; 49N90; 92D30

1. Introduction

In recent years, mathematical modeling of processes in biology and medicine, in particular in epidemiology, has led to significant scientific advances both in mathematics and biosciences [1,2]. Applications of mathematics in biology are completely opening new pathways of interactions, and this is certainly true in the area of optimal control: a branch of applied mathematics that deals with finding control laws for dynamical systems over a period of time such that an objective functional is optimized [3,4]. It has numerous applications in both biology and medicine [5–8].

To find the best possible control for taking a dynamical system from one state to another, one uses, in optimal control theory, the celebrated Pontryagin’s maximum principle (PMP), formulated in 1956 by the Russian mathematician Lev Pontryagin and his collaborators [3]. Roughly speaking, PMP states that it is necessary for any optimal control along with the optimal state trajectory to satisfy the so-called Hamiltonian system, which is a two-point boundary value problem, plus a maximality
condition on the Hamiltonian. Although a classical result, PMP is usually not taught to biologists and mathematicians working on mathematical biology. Here, we show how such scientists can easily implement the necessary optimality conditions given by the PMP, numerically, and can benefit from the power of optimal control theory. For that, we consider a mathematical model for HIV.

HIV modeling and optimal control is a subject under strong current research: see, e.g., Reference [9] and the references therein. Here, we consider the SICA epidemic model for HIV transmission proposed in References [10,11], formulate an optimal control problem with the goal to find the optimal HIV prevention strategy that maximizes the fraction of uninfected HIV individuals with least HIV new infections and cost associated with the control measures, and give complete algorithms in GNU Octave to solve the considered problems. We trust that our work, by providing the algorithms in an open programming language, contributes to reducing the so-called “replication crisis” (an ongoing methodological crisis in which it has been found that many scientific studies are difficult or impossible to replicate or reproduce [12]) in the area of optimal biomedical research. We trust our current work will be very useful to a practitioner from the disease control area and will become a reference in the field of epidemiology for those interested to include an optimal control component in their work.

2. A Normalized SICA HIV/AIDS Model

We consider the SICA epidemic model for HIV transmission proposed in References [10,11], which is given by the following system of ordinary differential equations:

\[
\begin{align*}
S'(t) &= bN(t) - \lambda(t)S(t) - \mu S(t) \\
I'(t) &= \lambda(t)S(t) - (\rho + \phi + \mu)I(t) + \alpha A(t) + \omega C(t) \\
C'(t) &= \phi I(t) - (\omega + \mu)C(t) \\
A'(t) &= \rho I(t) - (\alpha + \mu + d)A(t).
\end{align*}
\]

The model in Equation (1) subdivides human population into four mutually exclusive compartments: susceptible individuals (S); HIV-infected individuals with no clinical symptoms of AIDS (the virus is living or developing in the individuals but without producing symptoms or only mild ones) but able to transmit HIV to other individuals (I); HIV-infected individuals under ART treatment (the so-called chronic stage) with a viral load remaining low (C); and HIV-infected individuals with AIDS clinical symptoms (A). The total population at time \( t \), denoted by \( N(t) \), is given by \( N(t) = S(t) + I(t) + C(t) + A(t) \). Effective contact with people infected with HIV is at a rate \( \lambda(t) \), given by

\[
\lambda(t) = \frac{\beta}{N(t)}(I(t) + \eta_C C(t) + \eta_A A(t)),
\]

where \( \beta \) is the effective contact rate for HIV transmission. The modification parameter \( \eta_A \geq 1 \) accounts for the relative infectiousness of individuals with AIDS symptoms in comparison to those infected with HIV with no AIDS symptoms (individuals with AIDS symptoms are more infectious than HIV-infected individuals—pre-AIDS). On the other hand, \( \eta_C \leq 1 \) translates the partial restoration of immune function of individuals with HIV infection that use ART correctly [10]. All individuals suffer from natural death at a constant rate \( \mu \). Both HIV-infected individuals with and without AIDS symptoms have access to ART treatment: HIV-infected individuals with no AIDS symptoms \( I \) progress to the class of individuals with HIV infection under ART treatment \( C \) at a rate \( \phi \), and HIV-infected individuals with AIDS symptoms are treated for HIV at rate \( \alpha \). An HIV-infected individual with AIDS symptoms \( A \) that starts treatment moves to the class of HIV-infected individuals \( I \) and after, if the treatment is maintained, will be transferred to the chronic class \( C \). Individuals in the class \( C \) leave to the class \( I \) at a rate \( \omega \) due to a default treatment. HIV-infected individuals with no AIDS symptoms \( I \) that do not take ART treatment progress to the AIDS class \( A \) at rate \( \rho \). Finally, HIV-infected individuals with AIDS symptoms \( A \) suffer from an AIDS-induced death at a rate \( d \).
In the situation where the total population size \( N(t) \) is not constant, it is often convenient to consider the proportions of each compartment of individuals in the population, namely
\[
s = S/N, \quad i = I/N, \quad c = C/N, \quad r = R/N.
\]
The state variables \( s, i, c, \) and \( a \) satisfy the following system of differential equations:
\[
\begin{align*}
s'(t) &= b(1 - s(t)) - \beta(i(t) + \eta_Cc(t) + \eta_Aa(t))s(t) + da(t)s(t) \\
i'(t) &= \beta(i(t) + \eta_Cc(t) + \eta_Aa(t))s(t) - (\rho + \phi + b)i(t) + aa(t) + \omega c(t) + da(t)i(t) \\
c'(t) &= \phi i(t) - (\omega + b)c(t) + da(t)c(t) \\
a'(t) &= \rho i(t) - (\alpha + b + d)a(t) + da^2(t)
\end{align*}
\]
with \( s(t) + i(t) + c(t) + a(t) = 1 \) for all \( t \in [0, T] \).

3. Numerical Solution of the SICA HIV/AIDS Model

In this section, we consider Equation (2) subject to the initial conditions given by
\[
s(0) = 0.6, \quad i(0) = 0.2, \quad c(0) = 0.1, \quad a(0) = 0.1,
\]
by the fixed parameter values from Table 1, and by the final time value of \( T = 20 \) (years).

Table 1. Parameter values of the HIV/AIDS model in Equation (2) taken from Reference [11] and references cited therein.

<table>
<thead>
<tr>
<th>Symbol</th>
<th>Description</th>
<th>Value</th>
</tr>
</thead>
<tbody>
<tr>
<td>( \mu )</td>
<td>Natural death rate</td>
<td>( 1/69.54 )</td>
</tr>
<tr>
<td>( b )</td>
<td>Recruitment rate</td>
<td>( 2.1\mu )</td>
</tr>
<tr>
<td>( \beta )</td>
<td>HIV transmission rate</td>
<td>1.6</td>
</tr>
<tr>
<td>( \eta_C )</td>
<td>Modification parameter</td>
<td>0.015</td>
</tr>
<tr>
<td>( \eta_A )</td>
<td>Modification parameter</td>
<td>1.3</td>
</tr>
<tr>
<td>( \phi )</td>
<td>HIV treatment rate for ( I ) individuals</td>
<td>1</td>
</tr>
<tr>
<td>( \rho )</td>
<td>Default treatment rate for ( I ) individuals</td>
<td>0.1</td>
</tr>
<tr>
<td>( \alpha )</td>
<td>AIDS treatment rate</td>
<td>0.33</td>
</tr>
<tr>
<td>( \omega )</td>
<td>Default treatment rate for ( C ) individuals</td>
<td>0.09</td>
</tr>
<tr>
<td>( d )</td>
<td>AIDS induced death rate</td>
<td>1</td>
</tr>
</tbody>
</table>

All our algorithms, developed to solve numerically the initial value problems in Equations (2) and (3), are developed under the free GNU Octave software (version 5.1.0), a high-level programming language primarily intended for numerical computations and is the major free and mostly compatible alternative to MATLAB [13]. We implement three standard basic numerical techniques: Euler, second-order Runge–Kutta, and fourth-order Runge–Kutta. We compare the obtained solutions with the one obtained using the \texttt{ode45} GNU Octave function.

3.1. Default \texttt{ode45} routine of GNU Octave

Using the provided \texttt{ode45} function of GNU Octave, which solves a set of non-stiff ordinary differential equations with the well known explicit Dormand–Prince method of order 4, one can solve the initial value problems in Equations (2) and (3) as follows:

```octave
function dy = odeHIVsystem(t,y)
% Parameters of the model
mi = 1.0 / 69.54; b = 2.1 * mi; beta = 1.6;
etaC = 0.015; etaA = 1.3; fi = 1.0; ro = 0.1;
```
alfa = 0.33; omega = 0.09; d = 1.0;

% Differential equations of the model
dy = zeros(4,1);
aux1 = beta * (y(2) + etaC * y(3) + etaA * y(4)) * y(1);
aux2 = d * y(4);
dy(1) = b * (1 - y(1)) - aux1 + aux2 * y(1);
dy(2) = aux1 - (ro + fi + b - aux2) * y(2) + alfa * y(4) + omega * y(3);
dy(3) = fi * y(2) - (omega + b - aux2) * y(3);
dy(4) = ro * y(2) - (alfa + b + d - aux2) * y(4);

On the GNU Octave interface, one should then type the following instructions:

```octave
>> T = 20; N = 100;
>> [vT,vY] = ode45(@odeHIVsystem,[0:T/N:T],[0.6 0.2 0.1 0.1]);
```

Next, we show how such approach compares with standard numerical techniques.

### 3.2. Euler’s Method

Given a well-posed initial-value problem

\[
\frac{dy}{dt} = f(t,y) \quad \text{with} \quad y(a) = \alpha \quad \text{and} \quad a \leq t \leq b,
\]

Euler’s method constructs a sequence of approximation points \((t, w) \approx (t, y(t))\) to the exact solution of the ordinary differential equation by

\[
t_{i+1} = t_i + h \quad \text{and} \quad w_{i+1} = w_i + h f(t_i, w_i), \quad i = 0, 1, \ldots, N - 1,
\]

where \(t_0 = a, w_0 = \alpha, \) and \(h = (b - a) / N.\) Let us apply Euler’s method to approximate each one of the four state variables of the system of ordinary differential equations (Equation (2)). Our `odeEuler` GNU Octave implementation is as follows:

```octave
function dy = odeEuler(T)
    % Parameters of the model
    mi = 1.0 / 69.54; b = 2.1 * mi; beta = 1.6;
    etaC = 0.015; etaA = 1.3; fi = 1.0; ro = 0.1;
    alfa = 0.33; omega = 0.09; d = 1.0;

    % Parameters of the Euler method
    test = -1; deltaError = 0.001; M = 100;
    t = linspace(0,T,M+1); h = T / M;
    S = zeros(1,M+1); I = zeros(1,M+1);
    C = zeros(1,M+1); A = zeros(1,M+1);

    % Initial conditions of the model
    S(1) = 0.6; I(1) = 0.2; C(1) = 0.1; A(1) = 0.1;

    % Iterations of the method
    while(test < 0)
        oldS = S; oldI = I; oldC = C; oldA = A;
        for i = 1:M
            % Differential equations of the model
            aux1 = beta * (I(i) + etaC * C(i) + etaA * A(i)) * S(i);
            aux2 = d * A(i);
            auxS = b * (1 - S(i)) - aux1 + aux2 * S(i);
            auxI = aux1 - (ro + fi + b - aux2) * I(i) + alfa * A(i) + omega * C(i);
            auxC = fi * I(i) - (omega + b - aux2) * C(i);
            S(i+1) = S(i) + h * auxS;
            I(i+1) = I(i) + h * auxI;
            C(i+1) = C(i) + h * auxC;
            A(i+1) = A(i) + h * auxA;
            if abs(S(i+1) - oldS(i)) < deltaError && abs(I(i+1) - oldI(i)) < deltaError && abs(C(i+1) - oldC(i)) < deltaError && abs(A(i+1) - oldA(i)) < deltaError
                test = 0;
            end
        end
    end
```


auxA = ro * I(i) - (alfa + b + d - aux2) * A(i);

% Euler new approximation
S(i+1) = S(i) + h * auxS;
I(i+1) = I(i) + h * auxI;
C(i+1) = C(i) + h * auxC;
A(i+1) = A(i) + h * auxA;
end

% Absolute error for convergence
temp1 = deltaError * sum(abs(S)) - sum(abs(oldS - S));
temp2 = deltaError * sum(abs(I)) - sum(abs(oldI - I));
temp3 = deltaError * sum(abs(C)) - sum(abs(oldC - C));
temp4 = deltaError * sum(abs(A)) - sum(abs(oldA - A));
test = min(temp1,min(temp2,min(temp3,temp4)));
end
dy(1,:) = t; dy(2,:) = S; dy(3,:) = I;
dy(4,:) = C; dy(5,:) = A;

Figure 1 shows the solution of the system of ordinary differential equations (Equation (2)) with the initial conditions (Equation (3)), computed by the ode45 GNU Octave function (dashed line) versus the implemented Euler’s method (solid line). As depicted, Euler’s method, although being the simplest method, gives a very good approximation to the behaviour of each of the four system variables. Both implementations use the same discretization knots in the interval \([0, T]\) with a step size given by \(h = T/100\).
Euler’s method has a global error (total accumulated error) of $O(h)$, and therefore, the error bound depends linearly on the step size $h$, which implies that the error is expected to grow no worse than a linear manner. Consequently, diminishing the step size should give correspondingly greater accuracy to the approximations. Table 2 lists the norm of the difference vector, where each component of this vector is the absolute difference between the results obtained by the ode45 GNU Octave function and our implementation of Euler’s method, calculated by the vector norms 1, 2, and $\infty$.

### Table 2. Norms 1, 2, and $\infty$ of the difference vector between ode45 GNU Octave and Euler’s results.

<table>
<thead>
<tr>
<th>System Variables</th>
<th>$S(t)$</th>
<th>$I(t)$</th>
<th>$C(t)$</th>
<th>$A(t)$</th>
</tr>
</thead>
<tbody>
<tr>
<td>$|Octave - Euler|_1$</td>
<td>0.4495660</td>
<td>0.1646710</td>
<td>0.5255950</td>
<td>0.0443340</td>
</tr>
<tr>
<td>$|Octave - Euler|_2$</td>
<td>0.0659270</td>
<td>0.0301720</td>
<td>0.0783920</td>
<td>0.0101360</td>
</tr>
<tr>
<td>$|Octave - Euler|_{\infty}$</td>
<td>0.0161175</td>
<td>0.0113068</td>
<td>0.0190621</td>
<td>0.0041673</td>
</tr>
</tbody>
</table>

#### 3.3. Runge–Kutta of Order Two

Given a well-posed initial-value problem, the Runge–Kutta method of order two constructs a sequence of approximation points $(t, w) \approx (t, y(t))$ to the exact solution of the ordinary differential equation by $t_{i+1} = t_i + h$, $K_1 = f(t_i, w_i)$, $K_2 = f(t_i + h, w_i + hK_1)$, and $w_{i+1} = w_i + h \frac{K_1 + K_2}{2}$, for each $i = 0, 1, \ldots, N - 1$, where $t_0 = a$, $w_0 = \alpha$, and $h = (b - a) / N$. Our GNU Octave implementation of the Runge–Kutta method of order two applies the above formulation to approximate each of the four variables of the system in Equation (2). We implement the odeRungeKutta_order2 function through the following GNU Octave instructions:

```octave
function dy = odeRungeKutta_order2(T)
    % Parameters of the model
    mi = 1.0 / 69.54; b = 2.1 * mi; beta = 1.6;
    etaC = 0.015; etaA = 1.3; fi = 1.0; ro = 0.1;
    alfa = 0.33; omega = 0.09; d = 1.0;
    % Parameters of the Runge-Kutta (2nd order) method
    test = -1; deltaError = 0.001; M = 100;
    t = linspace(0,T,M+1); h = T / M; h2 = h / 2;
    S = zeros(1,M+1); I = zeros(1,M+1);
    C = zeros(1,M+1); A = zeros(1,M+1);
    % Initial conditions of the model
    S(1) = 0.6; I(1) = 0.2; C(1) = 0.1; A(1) = 0.1;
    % Iterations of the method
    while(test < 0)
        oldS = S; oldI = I; oldC = C; oldA = A;
        for i = 1:M
            % Differential equations of the model
            % First Runge-Kutta parameter
            aux1 = beta * (I(i) + etaC * C(i) + etaA * A(i)) * S(i);
            aux2 = d * A(i);
            auxS = b * (1 - S(i)) - aux1 + aux2 + S(i);
            auxI = aux1 - (ro + fi + b - aux2) * I(i) + alfa * A(i) + omega * C(i);
            auxC = fi * I(i) - (omega + b - aux2) * C(i);
            auxA = ro * I(i) - (alfa + b + d - aux2) * A(i);
            % Next Runge-Kutta parameter
            auxS = b * (1 - S(i)) - aux1 + aux2 + S(i);
            auxI = aux1 - (ro + fi + b - aux2) * I(i) + alfa * A(i) + omega * C(i);
            auxC = fi * I(i) - (omega + b - aux2) * C(i);
            auxA = ro * I(i) - (alfa + b + d - aux2) * A(i);
            % Update values
            S(i+1) = S(i) + h * auxS;
            I(i+1) = I(i) + h * auxI;
            C(i+1) = C(i) + h * auxC;
            A(i+1) = A(i) + h * auxA;
        end
        if abs(S(end) - oldS(end)) < deltaError
            test = 0;
        end
    end
end
```
%% Second Runge-Kutta parameter
auxS = S(i) + h * auxS1; auxI = I(i) + h * auxI1;
auxC = C(i) + h * auxC1; auxA = A(i) + h * auxA1;
aux1 = beta * (auxI + etaC * auxC + etaA * auxA) * auxS;
aux2 = d * auxA;
auxS2 = b * (1 - auxS) - aux1 + aux2 * auxS;
auxI2 = aux1 - (ro + fi + b - aux2) * auxI + alfa * auxA + omega * auxC;
auxC2 = fi * auxI - (omega + b - aux2) * auxC;
auxA2 = ro * auxI - (alfa + b + d - aux2) * auxA;

%% Runge-Kutta new approximation
S(i+1) = S(i) + h2 * (auxS1 + auxS2);
I(i+1) = I(i) + h2 * (auxI1 + auxI2);
C(i+1) = C(i) + h2 * (auxC1 + auxC2);
A(i+1) = A(i) + h2 * (auxA1 + auxA2);
end

%% Absolute error for convergence
temp1 = deltaError * sum(abs(S)) - sum(abs(oldS - S));
temp2 = deltaError * sum(abs(I)) - sum(abs(oldI - I));
temp3 = deltaError * sum(abs(C)) - sum(abs(oldC - C));
temp4 = deltaError * sum(abs(A)) - sum(abs(oldA - A));
test = min(temp1,min(temp2,min(temp3,temp4)));
end
dy(1,:) = t; dy(2,:) = S; dy(3,:) = I; dy(4,:) = C; dy(5,:) = A;

Figure 2 shows the solution of the system of Equation (2) with the initial value conditions in
Equation (3) computed by the ode45 GNU Octave function (dashed line) versus our implementation
of the Runge–Kutta method of order two (solid line). As we can see, Runge–Kutta’s method
produces a better approximation than Euler’s method, since both curves in each plot of Figure 2
are indistinguishable.

Figure 2. HIV/AIDS system (Equation (2)): GNU Octave versus Runge–Kutta’s method of order two.
Runge–Kutta’s method of order two (RK2) has a global truncation error of order $O(h^2)$, and as it is known, this truncation error at a specified step measures the amount by which the exact solution to the differential equation fails to satisfy the difference equation being used for the approximation at that step. This might seems like an unlikely way to compare the error of various methods, since we really want to know how well the approximations generated by the methods satisfy the differential equation, not the other way around. However, we do not know the exact solution, so we cannot generally determine this, and the truncation error will serve quite well to determine not only the error of a method but also the actual approximation error. Table 3 lists the norm of the difference vector between the results from ode45 routine and Runge–Kutta’s method of order two results.

Table 3. Norms 1, 2, and $\infty$ of the difference vector between ode45 GNU Octave and RK2 results.

<table>
<thead>
<tr>
<th>System Variables</th>
<th>$S(t)$</th>
<th>$I(t)$</th>
<th>$C(t)$</th>
<th>$A(t)$</th>
</tr>
</thead>
<tbody>
<tr>
<td>$|\text{Octave} - \text{RungeKutta2}|_1$</td>
<td>0.0106530</td>
<td>0.0105505</td>
<td>0.0151705</td>
<td>0.0044304</td>
</tr>
<tr>
<td>$|\text{Octave} - \text{RungeKutta2}|_2$</td>
<td>0.0014868</td>
<td>0.0025288</td>
<td>0.0022508</td>
<td>0.0011695</td>
</tr>
<tr>
<td>$|\text{Octave} - \text{RungeKutta2}|_\infty$</td>
<td>0.0003341</td>
<td>0.0009613</td>
<td>0.0006695</td>
<td>0.0004678</td>
</tr>
</tbody>
</table>

3.4. Runge–Kutta of Order Four

The Runge–Kutta method of order four (RK4) constructs a sequence of approximation points $(t, w) \approx (t, y(t))$ to the exact solution of an ordinary differential equation by $t_{i+1} = t_i + h$, $K_1 = f(t_i, w_i)$, $K_2 = f(t_i + \frac{h}{2}, w_i + \frac{h}{2} K_1)$, $K_3 = f(t_i + \frac{h}{2}, w_i + \frac{h}{2} K_2)$, $K_4 = f(t_{i+1}, w_i + hK_3)$, and $w_{i+1} = w_i + \frac{h}{6}(K_1 + 2K_2 + 2K_3 + K_4)$, for each $i = 0, 1, \ldots, N - 1$, where $t_0 = a$, $w_0 = \alpha$, and $h = (b - a)/N$. Our GNU Octave implementation of the Runge–Kutta method of order four applies the above formulation to approximate the solution of the system in Equation (2) with the initial conditions of Equation (3) through the following instructions:

```octave
function dy = odeRungeKutta_order4(T)
% Parameters of the model
mi = 1.0 / 69.54; b = 2.1 * mi; beta = 1.6;
etaC = 0.015; etaA = 1.3; fi = 1.0; ro = 0.1;
alfa = 0.33; omega = 0.09; d = 1.0;

% Parameters of the Runge-Kutta (4th order) method
test = -1; deltaError = 0.001; M = 100;
t = linspace(0,T,M+1);
h = T / M; h2 = h / 2; h6 = h / 6;
S = zeros(1,M+1); I = zeros(1,M+1);
C = zeros(1,M+1); A = zeros(1,M+1);

% Initial conditions of the model
S(1) = 0.6; I(1) = 0.2; C(1) = 0.1; A(1) = 0.1;

% Iterations of the method
while(test < 0)
    oldS = S; oldI = I; oldC = C; oldA = A;
    for i = 1:M
        % Differential equations of the model
        % First Runge-Kutta parameter
        aux1 = beta * (I(i) + etaC * C(i) + etaA * A(i)) * S(i);
        aux2 = d * A(i);
        auxS1 = b * (1 - S(i)) - aux1 + aux2 * S(i);
        auxI1 = aux1 - (ro + fi + b - aux2) * I(i) + alfa * A(i) + omega * C(i);
```
auxC1 = fi * I(i) - (omega + b - aux2) * C(i);
auxA1 = ro * I(i) - (alfa + b + d - aux2) * A(i);

% Second Runge-Kutta parameter
auxS = S(i) + h2 * auxS1; auxI = I(i) + h2 * auxI1;
auxC = C(i) + h2 * auxC1; auxA = A(i) + h2 * auxA1;
aux1 = beta * (auxI + etaC * auxC + etaA * auxA) * auxS;
aux2 = d * auxA;

auxS2 = b * (1 - auxS) - aux1 + aux2 * auxS;
auxI2 = auxI - (ro + fi + b - aux2) * auxI + alfa * auxA + omega * auxC;
auxC2 = fi * auxI - (omega + b - aux2) * auxC;
auxA2 = ro * auxI - (alfa + b + d - aux2) * auxA;

% Fird Runge-Kutta parameter
auxS = S(i) + h2 * auxS2; auxI = I(i) + h2 * auxI2;
auxC = C(i) + h2 * auxC2; auxA = A(i) + h2 * auxA2;
aux1 = beta * (auxI + etaC * auxC + etaA * auxA) * auxS;
aux2 = d * auxA;

auxS3 = b * (1 - auxS) - aux1 + aux2 * auxS;
auxI3 = auxI - (ro + fi + b - aux2) * auxI + alfa * auxA + omega * auxC;
auxC3 = fi * auxI - (omega + b - aux2) * auxC;
auxA3 = ro * auxI - (alfa + b + d - aux2) * auxA;

% Fourth Runge-Kutta parameter
auxS = S(i) + h * auxS3; auxI = I(i) + h * auxI3;
auxC = C(i) + h * auxC3; auxA = A(i) + h * auxA3;
aux1 = beta * (auxI + etaC * auxC + etaA * auxA) * auxS;
aux2 = d * auxA;

auxS4 = b * (1 - auxS) - aux1 + aux2 * auxS;
auxI4 = auxI - (ro + fi + b - aux2) * auxI + alfa * auxA + omega * auxC;
auxC4 = fi * auxI - (omega + b - aux2) * auxC;
auxA4 = ro * auxI - (alfa + b + d - aux2) * auxA;

% Runge-Kutta new approximation
S(i+1) = S(i) + h6 * (auxS1 + 2 * (auxS2 + auxS3) + auxS4);
I(i+1) = I(i) + h6 * (auxI1 + 2 * (auxI2 + auxI3) + auxI4);
C(i+1) = C(i) + h6 * (auxC1 + 2 * (auxC2 + auxC3) + auxC4);
A(i+1) = A(i) + h6 * (auxA1 + 2 * (auxA2 + auxA3) + auxA4);
end

% Absolute error for convergence
temp1 = deltaError * sum(abs(S)) - sum(abs(oldS - S));
temp2 = deltaError * sum(abs(I)) - sum(abs(oldI - I));
temp3 = deltaError * sum(abs(C)) - sum(abs(oldC - C));
temp4 = deltaError * sum(abs(A)) - sum(abs(oldA - A));
test = min(temp1,min(temp2,min(temp3,temp4))); end

dy(1,:) = t; dy(2,:) = S; dy(3,:) = I;
dy(4,:) = C; dy(5,:) = A;

Figure 3 shows the solution of the initial value problem in Equations (2) and (3) computed by the ode45 GNU Octave function (dashed line) versus our implementation of the Runge–Kutta method.
of order four (solid line). The results of the Runge–Kutta method of order four are extremely good. Moreover, this method requires four evaluations per step and its global truncation error is $O(h^4)$.

Figure 3. HIV/AIDS system (Equation (2)): GNU Octave versus Runge–Kutta’s method of order four.

Table 4 lists the norm of the difference vector between results obtained by the Octave routine ode45 and the 4th order Runge–Kutta method.

Table 4. Norms 1, 2, and $\infty$ of the difference vector between ode45 GNU Octave and RK4 results.

<table>
<thead>
<tr>
<th>System Variables</th>
<th>$S(t)$</th>
<th>$I(t)$</th>
<th>$C(t)$</th>
<th>$A(t)$</th>
</tr>
</thead>
<tbody>
<tr>
<td>$|\text{Octave} - \text{RungeKutta4}|_1$</td>
<td>0.0003193</td>
<td>0.0002733</td>
<td>0.0004841</td>
<td>0.0000579</td>
</tr>
<tr>
<td>$|\text{Octave} - \text{RungeKutta4}|_2$</td>
<td>0.0000409</td>
<td>0.0000395</td>
<td>0.0000674</td>
<td>0.0000098</td>
</tr>
<tr>
<td>$|\text{Octave} - \text{RungeKutta4}|_\infty$</td>
<td>0.0000107</td>
<td>0.0000140</td>
<td>0.0000186</td>
<td>0.0000042</td>
</tr>
</tbody>
</table>

4. Optimal Control of HIV Transmission

In this section, we propose an optimal control problem that will be solved numerically in Octave/MATLAB in Section 4.2. We introduce a control function $u(\cdot)$ in the model of Equation (2), which represents the effort on HIV prevention measures, such as condom use (used consistently and correctly during every sex act) or oral pre-exposure prophylaxis (PrEP). The control system is given by
We consider the optimal control problem of determining which considers the fraction of susceptible individuals \((u)\) where the control programming problem \([15,16]\). For such an approach, one only needs to use the Octave/MATLAB routine. Indirect methods are more sound because they are based on Pontryagin’s Maximum Principle but less widespread since they are not immediately available in Octave/MATLAB. Here, we show how one can use Octave/MATLAB to solve optimal control problems through Pontryagin’s Maximum Principle, reducing the optimal control problem to the solution of a boundary value problem.

More precisely, the system of ordinary differential equations in \(a\) is considered with the set of admissible controls given by

\[
\Omega = \{ u(\cdot) \in L^\infty(0, T) \mid 0 \leq u(t) \leq u_{\text{max}}, \forall t \in [0, T] \}.
\]

We consider the optimal control problem of determining \((s^*(\cdot), i^*(\cdot), c^*(\cdot), a^*(\cdot))\) associated to an admissible control \(u^*(\cdot) \in \Omega\) on the time interval \([0, T]\), satisfying Equation (4) and the initial conditions of Equation (5) and maximizing the cost functional of Equation (6):

\[
J(u^*(\cdot)) = \max_{\Omega} J(u(\cdot)) = \max_{\Omega} \left( \int_0^T \left( s(t) - i(t) - u^2(t) \right) dt \right).
\]

Note that we are considering a \(L^2\)-cost function: the integrand of the cost functional \(J\) is concave with respect to the control \(u\). Moreover, the control system of Equation (4) is Lipschitz with respect to the state variables \((s, i, c, a)\). These properties ensure the existence of an optimal control \(u^*(\cdot)\) of the optimal control problem in Equations (4)–(8) (see, e.g., Reference [14]).
4.1. Pontryagin’s Maximum Principle

According to celebrated Pontryagin’s Maximum Principle (see, e.g., Reference [3]), if \( u^*(\cdot) \) is optimal for Equations (4)–(8) with fixed final time \( T \), then there exists a nontrivial absolutely continuous mapping \( \Lambda : [0, T] \to \mathbb{R}^4 \), \( \Lambda(t) = (\lambda_1(t), \lambda_2(t), \lambda_3(t), \lambda_4(t)) \), called the adjoint vector, such that

\[
\begin{align*}
  s' &= \frac{\partial H}{\partial \lambda_1}, \quad i' = \frac{\partial H}{\partial \lambda_2}, \quad c' = \frac{\partial H}{\partial \lambda_3}, \quad a' = \frac{\partial H}{\partial \lambda_4}, \\
  \lambda_1' &= -\frac{\partial H}{\partial s}, \quad \lambda_2' = -\frac{\partial H}{\partial i}, \quad \lambda_3' = -\frac{\partial H}{\partial c}, \quad \lambda_4' = -\frac{\partial H}{\partial a},
\end{align*}
\]

where

\[
H = H(s(t), i(t), c(t), a(t), \Lambda(t), u(t)) = s(t) - i(t) - u^2(t)
\]

\[
+ \lambda_1(t) \left( b(1 - s(t)) - (1 - u(t))\beta(i(t) + \eta C c(t) + \eta A a(t))s(t) + d a(t) s(t) \right)
\]

\[
+ \lambda_2(t) \left( (1 - u(t))\beta(i(t) + \eta C c(t) + \eta A a(t))s(t) - (\rho + \beta + b)i(t) + aa(t) + \omega c(t) + d a(t) i(t) \right)
\]

\[
+ \lambda_3(t) \left( \phi i(t) - (\omega + b)c(t) + d a(t) c(t) \right)
\]

\[
+ \lambda_4(t) \left( \rho i(t) - (\alpha + b + d)a(t) + d a^2(t) \right)
\]

is called the Hamiltonian and the maximality condition

\[
H(s^*(t), i^*(t), c^*(t), a^*(t), \Lambda(t), u^*(t)) = \max_{0 \leq u \leq u \max} H(s^*(t), i^*(t), c^*(t), a^*(t), \Lambda(t), u)
\]

holds almost everywhere on \([0, T]\). Moreover, the transversality conditions

\[
\lambda_i(T) = 0, \quad i = 1, \ldots, 4,
\]

hold. Applying the Pontryagin maximum principle to the optimal control problem in Equations (4)–(8), the following theorem follows.

**Theorem 1.** The optimal control problem of Equations (4)–(8) with fixed final time \( T \) admits a unique optimal solution \((s^*(\cdot), i^*(\cdot), c^*(\cdot), a^*(\cdot))\) associated to the optimal control \( u^*(\cdot) \) on \([0, T]\) described by

\[
u^*(t) = \min \left\{ \max \left\{ 0, \frac{\beta (i^*(t) + \eta C c^*(t) + \eta A a^*(t)) s^*(t) (\lambda_1(t) - \lambda_2(t))}{2} \right\}, u \max \right\}, \quad (9)
\]

where the adjoint functions satisfy

\[
\begin{align*}
\lambda_1'(t) &= -1 + \lambda_1(t) (b + (1 - u^*(t)) \beta (i^*(t) + \eta C c^*(t) + \eta A a^*(t)) - da^*(t)), \\
-\lambda_2'(t) &= (1 - u^*(t)) \beta (i^*(t) + \eta C c^*(t) + \eta A a^*(t)), \\
\lambda_3'(t) &= 1 + \lambda_1(t) (1 - u^*(t)) \beta s^*(t) - \lambda_2(t) ((1 - u^*(t)) \beta s^*(t) - (\rho + \phi + s^*(t)) + da^*(t)), \\
-\lambda_3(t) &= \lambda_4(t) \rho, \\
\lambda_4'(t) &= \lambda_1(t) (1 - u^*(t)) \beta \eta C s^*(t) - \lambda_2(t) ((1 - u^*(t)) \beta \eta C s^*(t) + \omega) + \lambda_4(t) (\omega + b - da^*(t))
\end{align*}
\]

subject to the transversality conditions \( \lambda_i(T) = 0, i = 1, \ldots, 4. \)

**Remark 1.** The uniqueness of the optimal control \( u^* \) is due to the boundedness of the state and adjoint functions and the Lipschitz property of the systems in Equations (4) and (10) (see References [17,18] and references cited therein).
We implement Theorem 1 numerically in Octave/MATLAB in Section 4.2, and the optimal solution \((s^*(\cdot), i^*(\cdot), c^*(\cdot), a^*(\cdot))\) associated to the optimal control \(u^*(\cdot)\) is computed for given parameter values and initial conditions.

4.2. Numerical Solution of the HIV Optimal Control Problem

The extremal given by Theorem 1 is now computed numerically by implementing a forward-backward fourth-order Runge–Kutta method (see, e.g., Reference [19]). This iterative method consists in solving the system in Equation (4) with a guess for the controls over the time interval \([0, T]\) using a forward fourth-order Runge–Kutta scheme and the transversality conditions \(\lambda_i(T) = 0, i = 1, \ldots, 4\). Then, the adjoint system in Equation (10) is solved by a backward fourth-order Runge–Kutta scheme using the current iteration solution of Equation (4). The controls are updated by using a convex combination of the previous controls and the values from Equation (9). The iteration is stopped if the values of unknowns at the previous iteration are very close to the ones at the present iteration.

Our `odeRungeKutta_order4_WithControl` function is implemented by the following GNU Octave instructions:

```octave
function dy = odeRungeKutta_order4_WithControl(T)
% Parameters of the model
mi = 1.0 / 69.54; b = 2.1 * mi; beta = 1.6;
etaC = 0.015; etaA = 1.3; fi = 1.0; ro = 0.1;
alfa = 0.33; omega = 0.09; d = 1.0;

% Parameters of the Runge-Kutta (4th order) method
test = -1; deltaError = 0.001; M = 1000;
t = linspace(0,T,M+1);
h = T / M; h2 = h / 2; h6 = h / 6;
S = zeros(1,M+1); I = zeros(1,M+1);
C = zeros(1,M+1); A = zeros(1,M+1);

% Initial conditions of the model
S(1) = 0.6; I(1) = 0.2; C(1) = 0.1; A(1) = 0.1;

%Vectors for system restrictions and control
Lambda1 = zeros(1,M+1); Lambda2 = zeros(1,M+1);
Lambda3 = zeros(1,M+1); Lambda4 = zeros(1,M+1);
U = zeros(1,M+1);

% Iterations of the method
while(test < 0)
    oldS = S; oldI = I; oldC = C; oldA = A;
    oldLambda1 = Lambda1; oldLambda2 = Lambda2;
    oldLambda3 = Lambda3; oldLambda4 = Lambda4;
    oldU = U;

    %Forward Runge-Kutta iterations
    for i = 1:M
        aux1 = (1 - U(i)) * beta * (I(i) + etaC * C(i) + etaA * A(i)) * S(i);
        aux2 = d * A(i);
        auxS1 = b * (1 - S(i)) - aux1 + aux2 * S(i);
        auxI1 = aux1 - (ro + fi + b - aux2) * I(i) + alfa * A(i) + omega * C(i);
        auxC1 = fi * I(i) - (omega + b - aux2) * C(i);
        auxA1 = ro * I(i) - (alfa + b + d - aux2) * A(i);
    end

    %Backward Runge-Kutta iterations
    for i = M:-1:1
        oldLambda1 = Lambda1; oldLambda2 = Lambda2;
        oldLambda3 = Lambda3; oldLambda4 = Lambda4;
        oldU = U;

        U(i) = oldU(i) + h6 * (auxS1 + auxI1 + auxC1 + auxA1);
        Lambda1(i) = oldLambda1(i) + h6 * (auxS1);
        Lambda2(i) = oldLambda2(i) + h6 * (auxI1);
        Lambda3(i) = oldLambda3(i) + h6 * (auxC1);
        Lambda4(i) = oldLambda4(i) + h6 * (auxA1);
    end

    %Check of the difference between iterations
    if abs(h * (U(2) - U(1))) < deltaError
        test = 0;
    end
end

% Final values of the model
S(end) = oldS(end); I(end) = oldI(end); C(end) = oldC(end); A(end) = oldA(end);
```

% Second Runge-Kutta parameter
auxU = 0.5 * (U(i) + U(i+1));
auxS = S(i) + h2 * auxS1; auxI = I(i) + h2 * auxI1;
auxC = C(i) + h2 * auxC1; auxA = A(i) + h2 * auxA1;
aux1 = (1 - auxU) * beta * (auxI + etaC * auxC + etaA * auxA) * auxS;
aux2 = d * auxA;
auxS2 = b * (1 - auxS) - aux1 + aux2 + auxS;
auxI2 = aux1 - (ro + fi + b - aux2) * auxI + alfa * auxA + omega * auxC;
auxC2 = fi * auxI - (omega + b - aux2) * auxC;
auxA2 = ro * auxI - (alfa + b + d - aux2) * auxA;

% Third Runge-Kutta parameter
auxS = S(i) + h2 * auxS2; auxI = I(i) + h2 * auxI2;
auxC = C(i) + h2 * auxC2; auxA = A(i) + h2 * auxA2;
aux1 = (1 - auxU) * beta * (auxI + etaC * auxC + etaA * auxA) * auxS;
aux2 = d * auxA;
auxS3 = b * (1 - auxS) - aux1 + aux2 + auxS;
auxI3 = aux1 - (ro + fi + b - aux2) * auxI + alfa * auxA + omega * auxC;
auxC3 = fi * auxI - (omega + b - aux2) * auxC;
auxA3 = ro * auxI - (alfa + b + d - aux2) * auxA;

% Fourth Runge-Kutta parameter
auxS = S(i) + h * auxS3; auxI = I(i) + h * auxI3;
auxC = C(i) + h * auxC3; auxA = A(i) + h * auxA3;
aux1 = (1 - U(i+1)) * beta * (auxI + etaC * auxC + etaA * auxA) * auxS;
aux2 = d * auxA;
auxS4 = b * (1 - auxS) - aux1 + aux2 + auxS;
auxI4 = aux1 - (ro + fi + b - aux2) * auxI + alfa * auxA + omega * auxC;
auxC4 = fi * auxI - (omega + b - aux2) * auxC;
auxA4 = ro * auxI - (alfa + b + d - aux2) * auxA;

% Runge-Kutta new approximation
S(i+1) = S(i) + h6 * (auxS1 + 2 * (auxS2 + auxS3) + auxS4);
I(i+1) = I(i) + h6 * (auxI1 + 2 * (auxI2 + auxI3) + auxI4);
C(i+1) = C(i) + h6 * (auxC1 + 2 * (auxC2 + auxC3) + auxC4);
A(i+1) = A(i) + h6 * (auxA1 + 2 * (auxA2 + auxA3) + auxA4);
end

% Backward Runge-Kutta iterations
for i = 1:M
  j = M + 2 - i;

  % Differential equations of the model
  % First Runge-Kutta parameter
  auxU = 1 - U(j);
  aux1 = auxU * beta * (I(j) + etaC * C(j) + etaA * A(j));
  aux2 = d * A(j);
  auxLambda11 = -1 + Lambda1(j) * (b + aux1 - aux2) - Lambda2(j) * aux1;
  aux1 = auxU * beta * S(j);
  auxLambda21 = 1 + Lambda1(j) * aux1 - Lambda2(j) * (aux1 - (ro + fi + b) + ...
    + aux2) - Lambda3(j) * fi - Lambda4(j) * ro;
\[ aux1 = auxU \times \beta \times \etaC \times S(j); \]
\[ aux\Lambda31 = \Lambda1(j) \times aux1 - \Lambda2(j) \times (aux1 + \ldots + \omega) + \Lambda3(j) \times (\omega + b - aux2); \]
\[ aux1 = auxU \times \beta \times \etaA \times S(j); \]
\[ aux\Lambda41 = \Lambda1(j) \times (aux1 + d \times S(j)) \ldots - \Lambda2(j) \times (aux1 + alfa + \ldots + d \times I(j)) - \Lambda3(j) \times d \times C(j) + \ldots + \Lambda4(j) \times (alfa + b + d - 2 \times aux2); \]

% Second Runge-Kutta parameter
\[ auxU = 1 - 0.5 \times \left( U(j) + U(j-1) \right); \]
\[ auxS = 0.5 \times \left( S(j) + S(j-1) \right); \]
\[ auxI = 0.5 \times \left( I(j) + I(j-1) \right); \]
\[ auxC = 0.5 \times \left( C(j) + C(j-1) \right); \]
\[ auxA = 0.5 \times \left( A(j) + A(j-1) \right); \]
\[ aux1 = auxU \times \beta \times (auxI + \etaC \times auxC + \etaA \times auxA); \]
\[ aux2 = d \times auxA; \]
\[ aux\Lambda1 = \Lambda1(j) - h2 \times aux\Lambda11; \]
\[ aux\Lambda2 = \Lambda2(j) - h2 \times aux\Lambda21; \]
\[ aux\Lambda3 = \Lambda3(j) - h2 \times aux\Lambda31; \]
\[ aux\Lambda4 = \Lambda4(j) - h2 \times aux\Lambda41; \]
\[ aux\Lambda12 = -1 + aux\Lambda1 \times (b + aux1 - aux2) - aux\Lambda2 \times aux1; \]
\[ aux1 = auxU \times \beta \times auxS; \]
\[ aux\Lambda22 = 1 + aux\Lambda1 \times aux1 - aux\Lambda2 \times (aux1 - (ro + fi + b) + \ldots + aux2) - aux\Lambda3 \times fi - aux\Lambda4 \times ro; \]
\[ aux1 = auxU \times \beta \times \etaC \times auxS; \]
\[ aux\Lambda32 = aux\Lambda1 \times aux1 - aux\Lambda2 \times (aux1 + \ldots + \omega) + aux\Lambda3 \times (omega + b - aux2); \]
\[ aux1 = auxU \times \beta \times \etaA \times auxS; \]
\[ aux\Lambda42 = aux\Lambda1 \times (aux1 + d \times auxS) \ldots - aux\Lambda2 \times (aux1 + alfa + \ldots + d \times auxI) - aux\Lambda3 \times d \times auxC + \ldots + aux\Lambda4 \times (alfa + b + d - 2 \times aux2); \]

% Third Runge-Kutta parameter
\[ aux1 = auxU \times \beta \times (auxI + \etaC \times auxC + \etaA \times auxA); \]
\[ aux\Lambda1 = \Lambda1(j) - h2 \times aux\Lambda12; \]
\[ aux\Lambda2 = \Lambda2(j) - h2 \times aux\Lambda22; \]
\[ aux\Lambda3 = \Lambda3(j) - h2 \times aux\Lambda32; \]
\[ aux\Lambda4 = \Lambda4(j) - h2 \times aux\Lambda42; \]
\[ aux\Lambda13 = -1 + aux\Lambda1 \times (b + aux1 - aux2) - aux\Lambda2 \times aux1; \]
\[ aux1 = auxU \times \beta \times auxS; \]
\[ aux\Lambda23 = 1 + aux\Lambda1 \times aux1 \ldots - aux\Lambda2 \times (aux1 - (ro + fi + b) + aux2) - aux\Lambda3 \times fi - aux\Lambda4 \times ro; \]
\[ aux1 = auxU \times \beta \times \etaC \times auxS; \]
\[ aux\Lambda33 = aux\Lambda1 \times aux1 - aux\Lambda2 \times (aux1 + \ldots + \omega) + aux\Lambda3 \times (omega + b - aux2); \]
\[ aux1 = auxU \times \beta \times \etaA \times auxS; \]
\[ aux\Lambda43 = aux\Lambda1 \times (aux1 + d \times auxS) \ldots - aux\Lambda2 \times (aux1 + alfa + d \times auxI) - aux\Lambda3 \times d \times auxC + \ldots + aux\Lambda4 \times (alfa + b + d - 2 \times aux2); \]
% Fourth Runge-Kutta parameter
auxU = 1 - U(j-1); auxS = S(j-1);
auxI = I(j-1); auxC = C(j-1); auxA = A(j-1);

aux1 = auxU * beta * (auxI + etaC * auxC + etaA * auxA);
aux2 = d * auxA;
auxLambda1 = Lambda1(j) - h * auxLambda13;
auxLambda2 = Lambda2(j) - h * auxLambda23;
auxLambda3 = Lambda3(j) - h * auxLambda33;
auxLambda4 = Lambda4(j) - h * auxLambda43;

auxLambda14 = -1 + auxLambda1 * (b + aux1 - aux2) ... 
    - auxLambda2 * aux1;
aux1 = auxU * beta * auxS;
auxLambda24 = 1 + auxLambda1 * aux1 ... 
    - auxLambda2 * (aux1 - (ro + fi + b) + ... 
    + aux2) - auxLambda3 * fi - auxLambda4 * ro;
aux1 = auxU * beta * etaC * auxS;
auxLambda34 = auxLambda1 * aux1 - auxLambda2 * (aux1 + ... 
    + omega) + auxLambda3 * (omega + b - aux2);
aux1 = auxU * beta * etaA * auxS;
auxLambda44 = auxLambda1 * (aux1 + d * auxS) ... 
    - auxLambda2 * (aux1 + alfa + ... 
    + d * aux1) - auxLambda3 * d * auxC + ... 
    + auxLambda4 * (alfa + b + d - 2 * aux2);

% Runge-Kutta new approximation
Lambda1(j-1) = Lambda1(j) - h6 * (auxLambda11 + ... 
    + 2 * (auxLambda12 + auxLambda13) + auxLambda14);
Lambda2(j-1) = Lambda2(j) - h6 * (auxLambda21 + ... 
    + 2 * (auxLambda22 + auxLambda23) + auxLambda24);
Lambda3(j-1) = Lambda3(j) - h6 * (auxLambda31 + ... 
    + 2 * (auxLambda32 + auxLambda33) + auxLambda34);
Lambda4(j-1) = Lambda4(j) - h6 * (auxLambda41 + ... 
    + 2 * (auxLambda42 + auxLambda43) + auxLambda44);
end

% New vector control
for i = 1:M+1
vAux(i) = 0.5 * beta * (I(i) + etaC * C(i) + ... 
    + etaA * A(i)) * S(i) * (Lambda1(i) - Lambda2(i));
auxU = min(max([0.0 vAux(i)]) 0.5));
U(i) = 0.5 * (auxU + oldU(i));
end

% Absolute error for convergence
temp1 = deltaError * sum(abs(S)) - sum(abs(oldS - S));
temp2 = deltaError * sum(abs(I)) - sum(abs(oldI - I));
temp3 = deltaError * sum(abs(C)) - sum(abs(oldC - C));
temp4 = deltaError * sum(abs(A)) - sum(abs(oldA - A));
temp5 = deltaError * sum(abs(U)) - sum(abs(oldU - U));
temp6 = deltaError * sum(abs(Lambda1)) - sum(abs(oldLambda1 - Lambda1));
temp7 = deltaError * sum(abs(Lambda2)) - sum(abs(oldLambda2 - Lambda2));
temp8 = deltaError * sum(abs(Lambda3)) - sum(abs(oldLambda3 - Lambda3));
temp9 = deltaError * sum(abs(Lambda4)) - sum(abs(oldLambda4 - Lambda4));
test = min(temp1,min(temp2,min(temp3,min(temp4, ...
   min(temp5,min(temp6,min(temp7,min(temp8,temp9))))))));
end
dy(1,:) = t; dy(2,:) = S; dy(3,:) = I;
dy(4,:) = C; dy(5,:) = A; dy(6,:) = U;
disp("Value of LAMBDA at FINAL TIME");
disp([Lambda1(M+1) Lambda2(M+1) Lambda3(M+1) Lambda4(M+1)]);

For the numerical simulations, we consider \( u_{\text{max}} = 0.5 \), representing a lack of resources or misuse of the preventive HIV measures \( u(\cdot) \), that is, the set of admissible controls is given by

\[
\Omega = \{ u(\cdot) \in L^\infty(0,T) \mid 0 \leq u(t) \leq 0.5, \ \forall t \in [0,T] \} \tag{11}
\]

with \( T = 20 \) (years). Figures 4 and 5 show the numerical solution to the optimal control problem of Equations (4)–(8) with the initial conditions of Equation (3) and the admissible control set in Equation (11) computed by our odeRungeKutta_order4_WithControl function. Figure 6 depicts the extremal control behaviour of \( u^* \).

![Figure 4](image1.png)
![Figure 4](image2.png)

**Figure 4.** Optimal state variables for the control problem in Equations (4)–(8) subject to the initial conditions in Equation (3) and the admissible control set in Equation (11) *versus* trajectories without control measures.
Figure 5. Comparison: solutions to the initial value problem in Equations (2)–(3) versus solutions to the optimal control problem in Equations (4)–(8) subject to initial conditions in Equation (3) and the admissible control set in Equation (11).

Figure 6. Optimal control $u^*$ for the HIV optimal control problem in Equations (4)–(8) subject to the initial conditions in Equation (3) and the admissible control set in Equation (11).

5. Conclusion

The paper provides a study on numerical methods to deal with modelling and optimal control of epidemic problems. Simple but effective Octave/MATLAB code is fully provided for a recent model proposed in Reference [10]. The given numerical procedures are robust with respect to the parameters: we have used the same values as the ones in Reference [10], but the code is valid for other values of the parameters and easily modified to other models. The results show the effectiveness of optimal control theory in medicine and the usefulness of a scientific computing system such as GNU Octave: using the control measure as predicted by Pontryagin’s maximum principle and numerically computed by our Octave code, one sees that the number of HIV/AIDS-infected and -chronic individuals diminish and,
as a consequence, the number of susceptible (not ill) individuals increase. We trust our paper will be very useful to a practitioner from the disease control area. Indeed, this work has been motivated by many emails we received and continue to receive, asking us to provide the software code associated to our research papers on applications of optimal control theory in epidemiology, e.g., References [20,21].

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**References**


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