# A Moving Boundary Problem Arising from the Diffusion of Oxygen in Absorbing Tissue 

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[Received 17 September 1971 and in revised form 2 November 1971]


#### Abstract

Approximate analytical and numerical solutions of a partial differential equation are obtained which describe the diffusion of oxygen in an absorbing medium. Essential mathematical difficulties are associated with the presence of a moving boundary which marks the furthest penetration of oxygen into the medium and also with the need to allow for an initial distribution of oxygen through the medium.


## 1. Introduction

The Classical moving-boundary problem in heat flow which has been most thoroughly studied is one in which a change of state occurs on the moving interface. The velocity of the boundary is determined by the physical requirement that the latent heat required for the change of phase must be supplied or removed by conduction. Such problems are often referred to as "Stefan problems" after J. Stefan who published a paper on the subject towards the end of the nineteenth century. There is an extensive literature dating from that time. An excellent survey is given by Muehlbauer \& Sunderland (1965).

The present paper concerns a problem arising from the diffusion of oxygen in a medium which simultaneously consumes the oxygen. A moving boundary is an essential feature of this problem also, but the conditions which determine its movement are different. Not only is the concentration of oxygen always zero at the boundary but, in addition, no oxygen diffuses across the boundary at any time. There is thus no relationship which contains the velocity of the moving boundary explicitly. A combination of analytical and numerical methods are applied to this problem and the results are finally expressed in the form of an approximate polynomial expression.
The work is of immediate interest in medical research concerning the uptake of oxygen by tissue and the problem was suggested to us by Dr N. T. S. Evans at the Medical Research Council's Experimental Radiopathology Unit, Hammersmith Hospital.

## 2. Statement of the Problem

First, oxygen is allowed to diffuse into a medium, and some of the oxygen is absorbed by the medium, thereby being removed from the diffusion process. The concentration of oxygen at the surface of the medium is maintained constant. This first phase of the problem continues until a steady-state is reached in which the oxygen does not penetrate any further into the medium. The supply of oxygen is then cut off
and the surface is sealed so that no further oxygen passes in or out. The medium continues to absorb the available oxygen already in it and as a consequence the boundary marking the furthest depth of penetration in the steady-state, recedes towards the sealed surface. The major problem is that of tracing the movement of the boundary during this phase and of determining the distribution of oxygen through the medium as a function of time. A secondary problem in the application of numerical techniques is associated with the discontinuity in the derivative boundary condition which results from the abrupt sealing of the surface.
The diffusion-with-absorption process is represented by the partial differential equation

$$
\begin{equation*}
\frac{\partial C}{\partial T}=D \frac{\partial^{2} C}{\partial X^{2}}-m \tag{2.1}
\end{equation*}
$$

where $C(X, T)$ denotes the concentration of the oxygen free to diffuse at a distance $X$ from the outer surface of the medium at time $T, D$ is a constant diffusion coefficient and $m$, the rate of consumption of oxygen per unit volume of the medium, is also assumed constant.

The problem has two parts:-

## (a) Steady-state solution

During the initial phase, when the oxygen is entering through the surface, the following boundary condition is satisfied,

$$
\begin{equation*}
C=C_{0}, \quad X=0, \quad T \geqslant 0 \tag{2.2}
\end{equation*}
$$

where $C_{0}$ is a constant.
A steady-state is achieved in which the concentration at every point in the medium becomes independent of time, i.e. $\partial C / \partial T=0$ everywhere, when the gradient of concentration becomes zero at the point, $X_{0}$, in the medium where the concentration itself is zero. No oxygen can then diffuse beyond this point and we have the conditions,

$$
\begin{align*}
C & =0, \quad X \geqslant X_{0},  \tag{2.3}\\
\frac{\partial C}{\partial X} & =0, \quad X \geqslant X_{0}, \tag{2.4}
\end{align*}
$$

for $T \geqslant 0$.
The steady-state is defined by a solution of

$$
\begin{equation*}
D \frac{\partial^{2} C}{\partial X^{2}}-m=0 \tag{2.5}
\end{equation*}
$$

which satisfies the boundary conditions (2.2), (2.3) and (2.4). This solution is readily seen to be

$$
\begin{equation*}
C=\frac{m}{2 D}\left(X-X_{0}\right)^{2}, \tag{2.6}
\end{equation*}
$$

where

$$
\begin{equation*}
X_{0}=\sqrt{\left(\frac{2 D C_{0}}{m}\right)} \tag{2.7}
\end{equation*}
$$

## (b) Moving Boundary Problem

After the surface $X=0$ has been sealed, oxygen which is already in the medium, in the range $0 \leqslant X \leqslant X_{0}$, continues to be consumed. Consequently, the point of zeroconcentration which was initially given by (2.7) recedes towards $X=0$. Let the position of this point at any time, $T$, be represented by $X_{0}(T)$. The second phase of the problem can be expressed by the equation,

$$
\begin{equation*}
\frac{\partial C}{\partial T}=D \frac{\partial^{2} C}{\partial X^{2}}-m, \quad 0 \leqslant X \leqslant X_{0}(T) \tag{2.8}
\end{equation*}
$$

with the following conditions,

$$
\begin{align*}
& \frac{\partial C}{\partial X}=0, \quad X=0, \quad T \geqslant 0,  \tag{2.9}\\
C= & \frac{\partial C}{\partial X}=0, \quad X=X_{0}(T), \quad T \geqslant 0,  \tag{2.10}\\
C= & \frac{m}{2 D}\left(X-X_{0}\right)^{2}, \quad 0 \leqslant X \leqslant X_{0}, \quad T=0, \tag{2.11}
\end{align*}
$$

where $T=0$ is the moment when the surface is sealed. Making the changes of variables,

$$
x=\frac{X}{X_{0}}, \quad t=\frac{D}{X_{0}^{2}} T, \quad c=\frac{D}{m X_{0}^{2}}=\frac{C}{2 C_{0}},
$$

and denoting by $x_{0}(t)$ the value of $x$ corresponding to $X_{0}(T)$, the above system is reduced to the following non-dimensional form,

$$
\begin{equation*}
\frac{\partial c}{\partial t}=\frac{\partial^{2} c}{\partial x^{2}}-1, \quad 0 \leqslant x \leqslant x_{0}(t) \tag{2.12}
\end{equation*}
$$

with the corresponding boundary conditions,

$$
\begin{gather*}
\frac{\partial c}{\partial x}=0, \quad x=0, \quad t \geqslant 0,  \tag{2.13}\\
c=\frac{\partial c}{\partial x}=0, \quad x=x_{0}(t), \quad t \geqslant 0,  \tag{2.14}\\
c=\frac{1}{2}(1-x)^{2}, \quad 0 \leqslant x \leqslant 1, \quad t=0, \tag{2.15}
\end{gather*}
$$

where $x_{0}(0)=1$. The subscript $t$ in $x_{0}(t)$ is dropped in the following discussions.

## 3. Short-time Solution

The condition (2.15) shows that in the steady-state a negative unit gradient of concentration exists at the surface. When the surface is sealed a zero surface gradient is instantaneously imposed in accordance with (2.13). Because of this discontinuity in the surface-gradient numerical methods based on finite differences are liable to give inaccurate solutions in the neighbourhood of the surface for short引times. There will be an interval of time, however, before the disturbance at the surface has an effect on
the solution in the neighbourhood of $x=1$ to any specified degree of accuracy. Thus an analytical solution can be obtained which will provide a suitable approximation for small times, by assuming that the boundary, $x_{0}=1$, does not move initially.

The solution of (2.12) subject to the initial condition (2.15) and the boundary conditions (2.13) and

$$
\begin{equation*}
c=0, \quad x=1, \quad t \geqslant 0 \tag{3.1}
\end{equation*}
$$

is found by using Laplace Transforms to be

$$
\begin{align*}
& c(x, t)=\frac{1}{2}(1-x)^{2}+2 \sqrt{\left(\frac{t}{\pi}\right)} \sum_{n=0}^{\infty}(-1)^{n}\left[\exp \left\{-\left(\frac{2 n+2-x}{2 \sqrt{t}}\right)^{2}\right\}-\right. \\
& \left.\exp \left\{-\left(\frac{2 n+x}{2 \sqrt{ } t}\right)^{2}\right\}\right]-\sum_{n=0}^{\infty}(-1)^{n}\left\{(2 n+2-x) \operatorname{erfc}\left(\frac{2 n+2-x}{2 \sqrt{ } t}\right)-\right. \\
& \left.\quad(2 n+x) \operatorname{erfc}\left(\frac{2 n+x}{2 \sqrt{ } t}\right)\right\}, \quad 0 \leqslant x \leqslant 1, \quad t \geqslant 0 . \tag{3.2}
\end{align*}
$$

Values of $c(x, t)$ have been computed for $x=0(0.05) 1.0$. The typical curves of Fig. 1 demonstrate the general shape and confirm that the concentration has not changed within the accuracy of plotting near the boundary at $x=1$.


Fig. 1. Concentration distributions for $t<0.05$ before the boundary moves within the accuracy of plotting.

In computing $c(x, t)$ from (3.2), it is seen that the convergence of the infinite series is very rapid, so that the terms corresponding to $n=0$ are sufficient over an appreciable interval of time, when the terms less than $10^{-6}$ are neglected. Furthermore, for $0 \leqslant t \leqslant$ 0.020 , the second and the third series can be ignored to obtain an accuracy nowhere worse than $10^{-5}$. The concentration for $0 \leqslant t \leqslant 0.020$ can therefore be represented fairly accurately by the approximate expression,

$$
\begin{gather*}
\left.c(x, t)=\frac{1}{2}(1-x)^{2}-2 \sqrt{\left(\frac{t}{\pi}\right.}\right) \exp \left\{-\left(\frac{x}{2 \sqrt{ } t}\right)^{2}\right\}+x \operatorname{erfc}\left(\frac{x}{2 \sqrt{ } t}\right), \\
0 \leqslant x \leqslant 1 . \tag{3.3}
\end{gather*}
$$

## 4. Numerical Method

Once the boundary has started to move we resort to numerical methods of solution. Several methods have been proposed. Douglas \& Gallie (1955) introduced a method of variable time step, keeping the size of the space mesh fixed. Murray \& Landis (1959) used a variable space mesh and kept the time step fixed. Ehlrich (1958) employed implicit formula at the intermediate points and Taylor's expansions near the moving boundary in both time and space directions. Lotkin (1960) made use of subdivided differences while Crank (1957) suggested a three-point Lagrange interpolation formula near the moving boundary.
In the present analysis, the concentrations at the intermediate points between the two boundaries have been calculated by using simple explicit finite-difference formulae. Near the moving boundary a Lagrange-type formula has been used, as suggested by Crank (1957) because of convenience in calculation. The location of the moving point itself is determined by a Taylor's series. The method is described below in detail.

The whole region, $0 \leqslant x \leqslant 1$, is subdivided into $M$ intervals each of width $\delta x$ and we take $x_{r}=r \delta x$ where $0 \leqslant r \leqslant M(M \delta x=1)$.

### 4.1. Concentrations at the Intermediate Points

We assume that the concentrations at each of the grid points, at the $j$ th time level are known and the position of the moving boundary at that time is somewhere in the $r$ th interval between $x_{r-1}$ and $x_{r}$, given by $x_{0}=(r-1) \delta x+p^{j} \delta x$ where $p^{j}$ is positive and usually less than one, and is also known (Fig. 2). Then the concentrations at the


Fig. 2.
$(j+1)$ th time level, up to and including the mesh point $r-2$ can be calculated using the well known explicit formulae,

$$
\begin{align*}
& \frac{c_{0}^{j+1}-c_{0}^{j}}{\delta t}=\frac{2}{(\delta x)^{2}}\left(c_{1}^{j}-c_{0}^{j}\right)-1,  \tag{4.1}\\
& \frac{c_{k}^{j+1}-c_{k}^{j}}{\delta t}=\frac{1}{(\delta x)^{2}}\left(c_{k-1}^{j}-2 c_{k}^{j}+c_{k+1}^{j}\right)-1, \tag{4.2}
\end{align*}
$$

for $k=1,2, \ldots,(r-2)$, where $\delta t$ is the size of the time-step and $c_{k}^{j}$ denotes the concentration at point $k \delta x$ at time $j \delta t$.

### 4.2. Concentration in the Neighbourhood of the Moving Boundary

Let $f\left(a_{0}\right), f\left(a_{1}\right)$ and $f\left(a_{2}\right)$ be any function values corresponding to the arguments $a_{0}, a_{1}$ and $a_{2}$. A three-point Lagrangian interpolation formula can be written as

$$
f(x)=\frac{\left(x-a_{1}\right)\left(x-a_{2}\right)}{\left(a_{0}-a_{1}\right)\left(a_{0}-a_{2}\right)} f\left(a_{0}\right)+\frac{\left(x-a_{0}\right)\left(x-a_{2}\right)}{\left(a_{1}-a_{0}\right)\left(a_{1}-a_{2}\right)} f\left(a_{1}\right)+\frac{\left(x-a_{0}\right)\left(x-a_{1}\right)}{\left(a_{2}-a_{0}\right)\left(a_{2}-a_{1}\right)} f\left(a_{2}\right) .
$$

Differentiating the above twice with respect to $x$, we get

$$
\begin{equation*}
\frac{\partial^{2} f}{\partial x^{2}}=\frac{2 f\left(a_{0}\right)}{\left(a_{0}-a_{1}\right)\left(a_{0}-a_{2}\right)}+\frac{2 f\left(a_{1}\right)}{\left(a_{1}-a_{0}\right)\left(a_{1}-a_{2}\right)}+\frac{2 f\left(a_{2}\right)}{\left(a_{2}-a_{0}\right)\left(a_{2}-a_{1}\right)} . \tag{4.3}
\end{equation*}
$$

Application of (4.3) at the points $(r-2) \delta x,(r-1) \delta x$ and the moving point, and remembering the boundary condition (2.14), gives,

$$
\frac{\partial^{2} c}{\partial x^{2}}=\frac{2}{(\delta x)^{2}}\left(\frac{c_{r-2}}{1+p}-\frac{c_{r-1}}{p}\right)-1,
$$

and the appropriate finite-difference replacement at the point $(r-1) \delta x$ leads to

$$
\begin{equation*}
\frac{c_{r-1}^{j+1}-c_{r-1}^{j}}{\delta t}=\frac{2}{(\delta x)^{2}}\left(\frac{c_{r-2}^{j}}{1+p^{j}}-\frac{c_{r-1}^{j}}{p^{j}}\right)-1, \tag{4.4}
\end{equation*}
$$

an explicit expression for $c_{r-1}^{j+1}$.

### 4.3. Position of the Moving Boundary

In order to determine the location of the moving boundary, $x_{0}(t)$, we first derive some extra conditions there. Differentiation of (2.14) with respect to $t$, gives

$$
\begin{equation*}
\frac{d c}{d t}=\left(\frac{\partial c}{\partial x}\right)_{x=x_{0}}\left(\frac{d x_{0}}{d t}\right)+\left(\frac{\partial c}{\partial t}\right)_{x=x_{0}}=0 \tag{4.5}
\end{equation*}
$$

By using (2.12) and (2.14) in (4.5) we obtain

$$
\begin{equation*}
\frac{\partial^{2} c}{\partial x^{2}}=1, \quad x=x_{0} \tag{4.6}
\end{equation*}
$$

Differentiating (2.12) with respect to $x$, we get

$$
\begin{equation*}
\frac{\partial^{2} c}{\partial x \partial t}=\frac{\partial^{3} c}{\partial x^{3}} \tag{4.7}
\end{equation*}
$$

Again, we have from (2.14)

$$
\frac{d}{d t}\left(\frac{\partial c}{\partial x}\right)\left(\frac{\partial^{2} c}{\partial x^{2}}\right)_{x=x_{0}} \frac{d x_{0}}{d t}+\left(\frac{\partial^{2} c}{\partial t \partial x}\right)_{x=x_{0}}=0
$$

and hence using (4.6) and (4.7) in the above and assuming that order of differentiation by $x$ and $t$ can be interchanged we obtain

$$
\frac{\partial^{3} c}{\partial x^{3}}=-\frac{d x_{0}}{d t}, \quad x=x_{0}
$$

Similarly

$$
\frac{\partial^{4} c}{\partial x^{4}}=\left(\frac{d x_{0}}{d t}\right)^{2}, \quad \frac{\partial^{5} c}{\partial x^{5}}=-\frac{d^{2} x_{0}}{d t^{2}}-\left(\frac{d x_{0}}{d t}\right)^{3} \text { etc. }
$$



Now, the Taylor's series for $c_{r-1}$ obtained by expanding about the moving point can be written as

$$
\begin{align*}
c_{r-1} & =c\left(x_{0}\right)-p \delta x\left(\frac{\partial c}{\partial x}\right)_{x=x_{0}}+\frac{1}{2}(p \delta x)^{2}\left(\frac{\partial^{2} c}{\partial x^{2}}\right)_{x=x_{0}}-\frac{1}{\delta}(p \delta x)^{3} \frac{\partial^{3} c}{\partial x^{3}}+\ldots \\
& =\frac{1}{2}(p \delta x)^{2}+\frac{1}{6}(p \delta x)^{3} \frac{d x_{0}}{d t}+\ldots \tag{4.8}
\end{align*}
$$

Provided the boundary is not moving too quickly the first term of the series provides a reasonable approximation and gives

$$
\begin{equation*}
p=\frac{\sqrt{ }\left(2 c_{r-1}\right)}{\delta x} \tag{4.9}
\end{equation*}
$$

We shall see later that the boundary moves faster towards the end of the process and we then replace the finite-difference solution by an analytical expression.

When $c_{r-1}^{j+1}$ has been calculated from (4.4), the relation (4.9) gives the position of the moving point at the $(j+1)$ th time level.

### 4.4. Moving Boundary Crossing a Mesh Line

As $c_{r-1}$ goes on decreasing we look for either of the two possibilities (i) $c_{r-1}^{j+1} \leqslant 0$ or (ii) $c_{r-1}^{j+1}>c_{r-1}^{j}$. With regard to the first condition it is physically impossible for $c_{r-1}$ to go negative. When the second condition is detected, it shows that the numerical process has become unstable. A stability analysis is presented in the appendix to this paper. When either of the two conditions arises, the $(r-1)$ th mesh point is given up at the ( $j-1$ )th time level and onwards. The Lagrange formula is then applied to

Table 4.2
Comparison between analytical and numerical $(\delta x=0.05)$ solutions for small times. For each time the upper entry corresponds to the analytical solution and the lower entry to numerical solutions. Tabulated values are $10^{6} \mathrm{C}$

| $\boldsymbol{t}$ | 0.0 | 0.1 | 0.2 | 0.3 | 0.4 | 0.5 | 0.6 | 0.7 | 0.8 | 0.9 |
| :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: |
| 0.001 | 464318 | 404606 | 320000 | 245000 | 180000 | 125000 | 80000 | 45000 | 20000 | 5000 |
|  | 460000 | 405000 | 320000 | 245000 | 180000 | 125000 | 80000 | 45000 | 20000 | 5000 |
| 0.002 | 449538 | 401927 | 319973 | 245000 | 180000 | 125000 | 80000 | 45000 | 20000 | 5000 |
|  | 452000 | 405000 | 320000 | 245000 | 180000 | 125000 | 80000 | 45000 | 20000 | 5000 |
| 0.003 | 438197 | 397811 | 319760 | 244998 | 180000 | 125000 | 80000 | 45000 | 20000 | 5000 |
|  | 437600 | 398000 | 320000 | 245000 | 180000 | 125000 | 80000 | 45000 | 20000 | 5000 |
| 0.004 | 428636 | 393157 | 319212 | 244981 | 180000 | 125000 | 80000 | 45000 | 20000 | 5000 |
|  | 429600 | 394760 | 320000 | 245000 | 180000 | 125000 | 80000 | 45000 | 20000 | 5000 |
| 0.005 | 420213 | 388238 | 318302 | 244924 | 179999 | 125000 | 80000 | 45000 | 20000 | 5000 |
|  | 420320 | 389128 | 318976 | 245000 | 180000 | 125000 | 80000 | 45000 | 20000 | 5000 |
| 0.010 | 387164 | 365073 | 309950 | 243276 | 179804 | 124986 | 79999 | 45000 | 20000 | 5000 |
|  | 387497 | 365668 | 310719 | 243726 | 179927 | 125000 | 80000 | 45000 | 20000 | 5000 |
|  |  |  |  |  |  |  |  |  |  |  |
| 0.050 | 247691 | 240179 | 218845 | 186955 | 148992 | 109636 | 72962 | 42030 | 18856 | 4628 |
|  | 247841 | 240358 | 219089 | 187264 | 149327 | 109945 | 73208 | 42199 | 18955 | 4673 |

recalculate $c_{r-2}^{j}$ using a new value of $p$ at the $(j-1)$ th time which is taken to be the old value of $p^{j-1}$ plus one. This process is continued until there are at least two mesh points including the sealed surface. At the end, however, an approximate solution may be useful which is discussed in the next section.

Concentrations have been computed for $\delta x=0 \cdot 10,0 \cdot 05$ and $\delta t=0 \cdot 001$. A comparison is given in Table 4.1 to indicate the order of accuracy of the results. Table 4.2 shows that the values obtained by using $\delta x=0.05$, are in a very good agreement with those calculated from the Laplace solution, for small times. It should be noted that the numerical solutions involve large errors in the beginning at the surface due to discontinuity in the gradient there at zero time, but they very soon become consistent with the Laplace solutions. At $t=0.050$, the difference between the numerical and the Laplace solutions is not more than 0.0003 anywhere when the boundary $x_{0}$ has moved a distance of 0.003 from its original position $x_{0}=1$.

## 5. Integral Method

In this section we look for simple analytical expressions for the concentrationdistribution as well as for the location of the moving boundary at any given time. We shall make use of an approximate method that was introduced by Goodman (1958) and is usually referred to as the "Integral Method". A review of integral methods and their applications to a variety of transient-heat-transfer problems is to be found in Irvine \& Hartnett (1964).

### 5.1. Description of Integral Method

In applying the Integral Method to the present problem we choose a profile which satisfies all the known conditions. This profile involves the position of the moving point as a parameter to be determined. In order to find a moving point versus time relationship we integrate both sides of the differential equation (2.12) with respect to $x$ over the range for which it is valid, i.e. $0 \leqslant x \leqslant x_{0}$. This means that the differential equation is to be satisfied on average only and not at each point. Thus we obtain

$$
\begin{equation*}
\int_{0}^{x_{0}} \frac{\partial c}{\partial t} d x=\int_{0}^{x_{0}} \frac{\partial^{2} c}{\partial x^{2}} d x-\int_{0}^{x_{0}} d x \tag{5.1}
\end{equation*}
$$

Substituting the concentration profile in (5.1) and after a certain amount of manipulation we get an ordinary differential equation for the position of the moving boundary, $x_{0}$, with $t$ as the independent variable. Once the position of the moving point, $x_{0}$, is determined at any time, substitution of this value for the parameter $x_{0}$ in the profile gives the concentration distribution at that time.

### 5.2. Determination of Surface Concentration

Integral methods are not very amenable in cases of non-uniform initial distributions. In the present problem the discontinuity in the surface gradient is an additional difficulty. In order to apply an integral method we first get an expression for the surface concentration and use it as an additional condition to obtain the profile. We refer to the analytical solution (3.2) which has been obtained assuming the boundary, $x_{0}$, fixed at $x=1$. As described in Section 3, this solution is true everywhere for
small times i.e. until the boundary has not moved within the range of working accuracy. However, it is observed that the concentrations near the sealed surface have a close agreement with those obtained from the numerical solutions for $\delta x=0.05$ for all times. Therefore, an expression for surface concentration can be obtained by putting $x=0$ in (3.2). A closer examination of that expression reveals that the concentration varies linearly with the square-root of the time to an accuracy of $5 \times 10^{-4}$, as compared with the numerical solutions, and is given by

$$
\begin{equation*}
c(0, t)=\frac{1}{2}-2 \sqrt{\frac{t}{\pi}} \tag{5.2}
\end{equation*}
$$

Comparative figures are given in the following table for (i) analytical solution (3.2); (ii) numerical solution for $\delta x=0.05$ and (iii) approximate solution given by (5.2).

Table 5.1
Comparisons of $10^{6} \mathrm{c}$ at the sealed surface

| Time | 0.04 | 0.08 | 0.12 | 0.16 | 0.18 | 0.19 |
| :---: | :---: | :---: | :---: | :---: | :---: | :---: |
| Solutions |  |  |  |  |  |  |
| Analytical | 274328 | 180852 | 109134 | 48771 | 21546 | 8546 |
| Numerical | 274496 | 180969 | 109228 | 48893 | 21834 | 9039 |
| Approximate | 274324 | 180846 | 109118 | 48648 | 21269 | 8151 |

It may be mentioned here that the total time, $t_{1}$, for the concentration everywhere to become zero is given by $c\left(0, t_{1}\right)=0$ and is equal to $\pi / 16$ from (5.2).

### 5.3. Choosing a Polynomial Profile

A polynomial profile of fourth degree is now chosen containing five unknown parameters which might be functions of time and which are determined using (2.13), (2.14), (4.6) and (5.2). On writing $c_{0}$ for $c(0, t)$ the equation for the polynomial becomes

$$
\begin{equation*}
c\left(x, x_{0}\right)=\left(1-x / x_{0}\right)^{2}\left\{\frac{1}{2} x^{2}+4 c_{0}\left(1-x / x_{0}\right)-3 c_{0}\left(1-x / x_{0}\right)^{2}\right\} . \tag{5.3}
\end{equation*}
$$

This contains the position of the moving point, $x_{0}$, which still has to be determined.

### 5.4. Determination of the Moving Boundary

To obtain $x_{0}$ as a function of time we refer back to the equation (5.1) which gives

$$
\begin{equation*}
\int_{0}^{x_{0}} \frac{\partial c}{\partial t} d x=\frac{\partial}{\partial t} \int_{0}^{x_{0}} c d x=-x_{0} \tag{5.4}
\end{equation*}
$$

since $\partial c / \partial x=0$ at $x=0, x_{0}$.
Writing $c\left(x, x_{0}\right)$ from (5.3) in (5.4) and using (5.2) we get, after some manipulation

$$
\begin{equation*}
\frac{d x_{0}}{d t}=-\frac{\{20-8 / \sqrt{ }(\pi t)\} x_{0}}{x_{0}^{2}+4-16 \sqrt{ }(t / \pi)} . \tag{5.5}
\end{equation*}
$$

We know that $d x_{0} / d t \leqslant 0$. This condition will not be true until

$$
\begin{equation*}
20-\frac{8}{\sqrt{(\pi t)}} \geqslant 0 \tag{5.6}
\end{equation*}
$$

since the term in the denominator of (5.5) is positive for $0 \leqslant x_{0} \leqslant 1$ and $t \leqslant t_{1}$, where $t_{1}$ is obtained from (5.2).

The inequality (5.6) gives the minimum time $t_{0}$ for the condition $d x_{0} / d t \leqslant 0$ to hold as $4 / 25 \pi$. It should be noted that as the moving point $x_{0}$ approaches the sealed surface, its speed $d x_{0} / d t$ tends to infinity as $t$ tends to $t_{1}$.

We have found that the numerical solution of (5.5) obtained by using a Runge-Kutta algorithm can be approximated by the expression,

$$
\begin{equation*}
x_{0}=1-\exp \left\{-2\left(\frac{t_{1}-t}{t-t_{0}}\right)^{\frac{1}{2}}\right\} . \tag{5.7}
\end{equation*}
$$

Table 5.2 below provides a comparison for the position of the moving boundary as obtained from (i) numerical evaluation of (5.5) using Runge-Kutta method (ii) approximation (5.7) and (iii) numerical method of Section 4.

Table 5.2
Comparison for $10^{4} x_{0}$ at different times

| Time | 0.051 | 0.060 | 0.080 | 0.100 | 0.120 | 0.140 | 0.160 | 0.180 | 0.190 | 0.195 |
| :--- | :--- | :--- | :--- | :--- | :--- | :--- | :--- | :--- | :--- | :--- | :--- |
| Numerical <br> Solution <br> of (5.5) | 10000 | 9974 | 9750 | 9321 | 8686 | 7817 | 6634 | 4892 | 3505 | 2331 |
| Approximate <br> solution (5.7) | 10000 | 9996 | 9817 | 9393 | 8779 | 7962 | 6848 | 5092 | 3478 | 1760 |
| Numerical <br> Method of <br> section 4. | 9967 | 9922 | 9719 | 9352 | 8788 | 7975 | 6812 | 4959 | 3381 | 1618 |

It is seen from the above table that the numerical solution of (5.5) agrees with the exponential profile for the moving boundary (5.7) very well except for very large times. But the profile (5.7) has a very good agreement with the values obtained from the numerical method for all times. Therefore, (5.2) and (5.3) together with (5.7) constitute an approximate solution. It should be noted that this solution is applicable for the time interval $4 / 25 \pi \leqslant t \leqslant \pi / 16$ only. For $t \leqslant 4 / 25 \pi$, Laplace solutions (3.2) and (3.3) give analytical solutions when it has been assumed that the boundary has not moved from its original position $x_{0}=1$. Thus we have got now an analytical solution of the problem for all times.

## 6. Results and Discussion

The concentrations in the medium at various times together with the position of the moving boundary have been compared in Table 6.1 for numerical and the approximate solutions. A very close agreement is seen between the two solutions. The approximate method would specially be useful (a) to calculate the concentration and the position of the moving boundary at an arbitrary time and (b) at the end when the numerical method would not work because too few mesh points remain. Graphs have been drawn to show the concentration-distributions at various times (Fig. 3) and the progress of the moving boundary with respect to time (Fig. 4).
Note. For $t<0.050$ see Table 4.2.
Comparison of concentrations $10^{6} \mathrm{c}$ and the distances of the moving boundary from the sealed surf obtained from the approximate solution and the lower entry values from the namerical solution ( $\delta x=0.05$ )



Fig. 3. Concentration distributions for the steady-state $(t=0)$ and for $t>0.05$.


Fig. 4. Position of the moving boundary with respect to time.

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## Appendix. Stability Analysis

The set of difference equations connecting values of $c$ at two consecutive time levels can be written in the following matrix form $\left(r=\delta t /(\delta x)^{2}\right)$,

$$
\left[\begin{array}{c}
c_{0}^{j+1}  \tag{1}\\
c_{1}^{j+1} \\
\cdot \\
\cdot \\
\cdot \\
c_{N-2}^{j+1} \\
c_{N-1}^{j+1}
\end{array}\right] .\left[\begin{array}{ccccc}
1-2 r & 2 r & & & \\
r & 1-2 r & r & 0 & \\
& & \cdot & & \\
& & \cdot & & \\
& & & \cdot & \\
& 0 & r & 1-2 r & r \\
& & & \frac{2 r}{1+p^{j}} & 1-\frac{2 r}{p^{j}}
\end{array}\right]\left[\begin{array}{c}
c_{0}^{j} \\
c_{1}^{j} \\
\cdot \\
\cdot \\
\cdot \\
c_{N-2}^{j} \\
c_{N-1}^{j}
\end{array}\right]-\left[\begin{array}{c}
1 \\
1 \\
. \\
\cdot \\
. \\
1 \\
1
\end{array}\right] . \delta t
$$

or

$$
\begin{equation*}
\mathbf{c}^{j+1}=\mathbf{A}_{j} \mathbf{c}^{j}-\mathbf{U} \delta t \tag{2}
\end{equation*}
$$

where $\mathbf{A}_{j}$ is a square matrix of order $N$ which varies with $N$ and $\mathbf{U}$ is a column vector each element of which is unity. We see that elements in the last row of $\mathbf{A}_{\boldsymbol{j}}$ are depend ent on $j$ and therefore in order to make analysis possible we first replace $p^{j}$ by a constant value $p$. Later on conditions are imposed on $p$ in order to make the scheme stable. Equation (2) is then written as

$$
\begin{equation*}
\mathbf{c}^{j+1}=\mathbf{A c}^{j}-\mathbf{U} \delta t \tag{3}
\end{equation*}
$$

We denote the computed values by $\psi$ so that we have actually solved the equations

$$
\begin{equation*}
\psi^{j+1}=\mathbf{A} \psi^{j}-\mathbf{U} \delta \mathbf{t} . \tag{4}
\end{equation*}
$$

The computational error is then given by subtracting (4) from (3). If the error introduced at the $k$ th step is denoted by the vector $\mathrm{e}^{k}$, then

$$
\mathbf{c}^{j+1}-\psi^{j+1}=\mathbf{A}\left(\mathbf{c}^{j}-\psi^{j}\right)
$$

i.e.

$$
\begin{equation*}
\mathbf{e}^{j+1}=\mathbf{A} \mathbf{e}^{j} \tag{5}
\end{equation*}
$$

The recurrence relation (5) gives

$$
\begin{equation*}
\mathbf{e}^{j+1}=(\mathbf{A})^{j+1} \mathbf{e}^{0} \tag{6}
\end{equation*}
$$

where $\mathbf{e}^{0}$ is an error vector for the starting values.
Let us express $\mathbf{e}^{0}$ as the linear combination of the eigenvectors of $\mathbf{A}$, such that

$$
\mathrm{e}^{0}=\sum_{s=1}^{N} a_{s} \mathrm{v}_{s}
$$

where $\nabla_{s}$ is an eigenvector of $A$ corresponding to the eigenvalue $\lambda_{s}$ and $a$ 's are constants. It is easy to show that

$$
\mathbf{e}^{n}=\sum_{s=1}^{N} a_{s} \lambda_{s}^{\lambda_{s}} \mathbf{v}_{s}
$$

For $\mathbf{e}^{n}$ to tend to zero, as $n$ increases, it follows that the largest of $\left|\lambda_{1}\right|,\left|\lambda_{2}\right|, \ldots,\left|\lambda_{N}\right|$ must be less than unity. If $Q_{s}$ is the sum of the moduli of the terms along the sth row excluding the diagonal term $a_{s s}$ in matrix $A$ then by Brauer's theorem every eigenvalue of $\mathbf{A}$ lies inside or on the boundary of at least one of the circles $\left|\lambda-a_{s s}\right|=Q_{s}$.

As we are interested in the bounds of $p$, applying Brauer's theorem to the last row of A that contains $p$, we have

$$
Q_{s}=\frac{2 r}{1+p}, \quad a_{s s}=1-\frac{2 r}{p},
$$

so that

$$
\left|\lambda-\left(1-\frac{2 r}{p}\right)\right| \leqslant \frac{2 r}{1+p}
$$

The bounds for $\lambda$ are given by

$$
\lambda_{1}=1-\frac{2 r}{p(1+p)} ; \quad \lambda_{2}=\frac{2 r(1+2 p)}{p(1+p)}-1 .
$$

For stability we require $\left|\lambda_{1}\right| \leqslant 1,\left|\lambda_{2}\right| \leqslant 1$, and hence

$$
-1 \leqslant 1-\frac{2 r}{p(1+p)} \leqslant 1 \text { giving } \frac{r}{p(1+p)} \leqslant 1
$$

and

$$
-1 \leqslant \frac{2 r(1+2 p)}{p(1+p)} \leqslant 1 \quad \text { giving } \quad \frac{r(1+2 p)}{p(1+p)} \leqslant 1 .
$$

Since $p$ is always positive, the condition for stability is given by the second inequality because the first one is then satisfied automatically. Therefore, for overall stability

$$
p^{2}+(1-2 r) p-r \geqslant 0 .
$$

Since $r \leqslant \frac{1}{2}$ for the stability of the simple explicit scheme used at the intermediate points, it can be shown that

$$
\begin{equation*}
p \geqslant r-\frac{1}{2}+\sqrt{\left(\frac{1}{4}+r^{2}\right) .} \tag{7}
\end{equation*}
$$

For $\delta x=0.1, \delta t=0.001$ we get the stability condition $p \geqslant 0.11$ and for $\delta x=0.05$, $\delta t=0.001$ we have $p \geqslant 0.54$. This suggests that an instability may arise when the moving point is nearer than 0.011 to the neighbouring mesh point in the first case and 0.027 in the second case ( $\delta x=0.05$ ). This confirms the need for the stability check described in Section 4.4.

One of us (R. S. G.) has been supported during the course of this work by a studentship awarded by the Science Research Council.

