Lawrence Berkeley National Laboratory

Recent Work

Title

ACCURACY OF THE RANDOM VORTEX METHOD FOR A PROBLEM WITH NON-SMOOTH INITIAL CONDITIONS

Permalink

https://escholarship.org/uc/item/7np1x6q9

Author

Roberts, S.

Publication Date

1983-05-01



Lawrence Berkeley Laboratory

UNIVERSITY OF CALIFORNIA

Physics, Computer Science & CEIVED Mathematics Division

JUL 21 1983

To be submitted for publication

LIBRARY AND DOCUMENTS SECTION

ACCURACY OF THE RANDOM VORTEX METHOD FOR A PROBLEM WITH NON-SMOOTH INITIAL CONDITIONS

S. Roberts

May 1983

TWO-WEEK LOAN COPY

This is a Library Circulating Copy which may be borrowed for two weeks. For a personal retention copy, call Tech. Info. Division, Ext. 6782.



DISCLAIMER

This document was prepared as an account of work sponsored by the United States Government. While this document is believed to contain correct information, neither the United States Government nor any agency thereof, nor the Regents of the University of California, nor any of their employees, makes any warranty, express or implied, or assumes any legal responsibility for the accuracy, completeness, or usefulness of any information, apparatus, product, or process disclosed, or represents that its use would not infringe privately owned rights. Reference herein to any specific commercial product, process, or service by its trade name, trademark, manufacturer, or otherwise, does not necessarily constitute or imply its endorsement, recommendation, or favoring by the United States Government or any agency thereof, or the Regents of the University of California. The views and opinions of authors expressed herein do not necessarily state or reflect those of the United States Government or any agency thereof or the Regents of the University of California.

ACCURACY OF THE RANDOM VORTEX METHOD FOR A PROBLEM WITH NON-SMOOTH INITIAL CONDITIONS¹

Stephen Roberts

Department of Mathematics and Lawrence Berkeley Laboratory University of California Berkeley, CA. 94720

May 1983

ABSTRACT

The accuracy of the random vortex method, in approximating the solution of the Navier Stokes equation, is investigated using the model problem of a circular vortex as suggested by Milinazzo and Saffman [13]. The method consists of partitioning the vorticity into "vortex blobs". These blobs are moved via two actions. First, a blob is deterministically moved under the action of the velocity field associated with the other blobs. Then to simulate viscosity a random component is added to the position of the blob. For this model problem the nonlinear terms of the Navier Stokes equation vanish. Thus the major error inherent in the deterministic component of the method vanishes. Consequently, for this model problem we concentrate on the interaction of the deterministic and random components of the method. Our results show that the accuracy of the method depends heavily on the initial distribution and strength of the computational elements i.e., the vortex blobs. With the right choice of initial conditions we find that the error is $O(R^{-\frac{N}{N}N^{-\frac{N}{N}}})$, where R is a Reynolds number and N is the number of vortex blobs.

¹This work was supported in part by the Director, Office of Energy Research, Office of Basic Energy Sciences, Engineering, Mathematical, and Geosciences Division of the U.S. Department of Energy, under Contract DE-ACO3-76SF00098.

1. INTRODUCTION

The vorticity stream formulation of the incompressible Navier-Stokes equation in the plane has the form,

$$\omega_t + u \,\omega_x + v \,\omega_y = \frac{1}{R} \Delta \omega \quad , \omega(\overline{z}, 0) = \xi(\overline{z}) \tag{1.1}$$

$$\Delta \psi = \omega \tag{1.2}$$

$$u = -\psi_u \tag{1.3}$$

$$v = \psi_x \tag{1.4}$$

where $\bar{u} = (u, v)$ is the velocity, ω the vorticity, ψ the stream function, R is the Reynolds number and $\bar{z} = (x, y)$ are the spatial coordinates.

To solve these equations numerically it is common to solve a discrete approximation of these equations on a grid. These approximations introduce an artifical viscosity term. The random vortex method (Chorin [5]) is designed to overcome this difficulty. The method involves partitioning the vorticity into a sum of "blobs". It is a fractional step method. At each time step, the convective part the Navier-Stokes equation is solved by moving each of the blobs under the action of the of the velocity field associated with the other blobs (vortex method). The diffusion part is simulated by letting the particles undertake a random walk (Lamperti [10] ch.4). It should be noted that large statistical errors can occur in the solution of vorticity by using this method. On the other hand, as velocity is an integral of vorticity, the statistical error for the velocity is much smaller. Chorin's method also allows for domains with boundaries. We shall use this method in the case in which the domain is the whole plane. Thus it is not necessary to discuss the boundary conditions.

Chorin's method has, on the whole (see Leonard [11]), been received as a useful and important tool in the study of high Reynolds number, incompressible flows. However, Milinazzo and Saffman [13] came to the conclusion that the

method could produce significant errors, even when a large number of vortex blobs were used. They obtained 10% error with 1000 blobs for a Reynolds number of 5000. Chorin [6] has pointed out that their conclusions were based on their choice of error measurement. He also criticizes their implementation of the random vortex method, in particular their choice of cutoff parameter (see section 2).

It is our intent to reproduce Milinazzo and Saffman's results, in order to clarify their results and their disagreement with Chorin. In section 3, we present Milinazzo and Saffman's model problem. We analyse two methods of measuring the error for the method, the one proposed by Chorin in [6], and the one used by Milinazzo and Saffman in their paper. Estimates are derived for the errors expected in the numerical results. These estimates are based on the calculations of Milinazzo and Saffman [13], section 4, and agree with Chorin's [5] estimates.

In section 4 we compare the method used by Chorin and others to approximate the initial vorticity and the construction of Milinazzo and Saffman. In the first method the vortex blobs are initially placed on a uniform grid. In the second method the blobs are placed by use of a random distribution. We find that the first method produces more accurate results. For instance, using 500 blobs and a Reynolds number of 5000, we obtain a 5% error using the first method and a 12% error using the second method. These errors are calculated using Milinazzo and Saffman's method of error measurement. This result verifies Milinazzo and Saffman's result of a 10% error for 1000 blobs, since an increase in the number of blobs by a factor of 2 should decrease the error by a factor of $\sqrt{2}$. Note that if we use Chorin's error measurement for the two cases above, we obtain errors of 0.3% and 0.7% respectively.

We find that there is little difference in our results when different cutoff functions are used. In particular, we tested those cutoffs used by Beale and Majda [4], Chorin [5], Hald [7] and Milinazzo and Saffman [13] (see table 1). Since the average distribution of particles for our problem is radially symmetric it is reasonable to assume that the angular motion of the vortex particles is governed by the vortex method, whereas the radial motion is governed by the random walk algorithm. As our methods of error measurement are only sensitive to the average change in the radial position of the particles, it is not surprising that we cannot detect any difference between the various cutoff functions.

Our numerical runs show that there are two types of errors in the computations. We refer to these as the startup error and the interaction error. The startup error depends on the difference between the initial distribution of vortex blobs and the exact vorticity. The interaction error involves the errors produced through the interaction of the vortex method and the random walk algorithms at each time step. Our computations show that Milinazzo and Saffman's implementation gives 2 to 5 times larger startup errors than the standard method. For a Reynolds number greater than 1000 the startup error "swamps" the interaction error, i.e., the computed error is due to Milinazzo and Saffman's choice of initial conditions. However, for larger Reynolds numbers ($R \ge 20000$) even the standard method produces significant startup errors in comparison to the interaction error.

By measuring the error in a different way we can study the interaction term. We observe that the errors for the random vortex method and the solution of the diffusion equation using the random walk method are comparable, at least if the number of vortices are less than 500. Both methods produce errors of $O(N^{-\frac{1}{N}}R^{-\frac{1}{N}})$, where N is the number of blobs. This verifies our analysis in section 3. The error for the random vortex though, is typically larger than that of

the random walk algorithm.

With the lack of smoothness of our initial data, we could expect large errors. However the results are quite accurate and we use only moderate amounts of computing time (less than 4hr. on a Vax). Hald [7] and Beale and Majda [4] have shown that if the initial data is smooth then vortex methods can accurately solve the inviscid equations on IR². Thus, for problems involving smoother initial data, we expect more accurate results.

2. THE RANDOM VORTEX METHOD.

The random vortex method solves the Navier-Stokes equation (1.1-4) by a fractional step method, which consists of the two equations,

$$\omega_{t} = Q \omega \quad , \omega(\bar{z},0) = \xi_{1}(\bar{z})$$

$$Q \omega = -\left[u \omega_{x} + v \omega_{y}\right]$$

$$where \ \Delta \psi = \omega , \ u = -\psi_{y}, \ v = \psi_{x}$$
(2.1)

and

$$\omega_t = \frac{1}{R} \Delta \omega \quad , \omega(\bar{z}, 0) = \xi_2(\bar{z})$$
 (2.2)

At each time step, equation (2.1) is solved by following the evolution of a finite number of fluid elements (vortex blobs). The solution of equation (2.2) is then simulated by adding a gaussianly distributed random component, of appropriate variance, to the position of these fluid elements (blobs). An approximation to the Navier-Stokes equation is then obtained from the resulting vorticity distribution.

We will discuss the solution of equation (2.1) by the vortex method (Chorin [5]). The basic idea is to approximate the vorticity by a sum of "vortex blobs".

A blob can be thought of as a region of vorticity which is convected in the fluid

under the action of the other blobs. In a real fluid, the distribution of vorticity inside a blob would be distorted by the flow. In the vortex method it is assumed that the blobs are translated as a whole. The validity of this is proved in Hald [7] and Beale and Madja [4]. The problem is then reduced to following the evolution of a finite number of these blobs.

First it is necessary to choose a set of "cutoff functions", $\{\varphi_{\delta}\}_{\delta>0}$. These functions are approximate δ functions on the plane, such that for every function, $u \in L^1 \cap L^m$,

$$\varphi_{\delta} * u \rightarrow u \quad \text{as } \delta \rightarrow 0,$$
 (2.3)

Here \bullet denotes convolution. We call δ the "cutoff" for the function φ_{δ} . For any function φ with $\int \varphi = 1$, $\varphi_{\delta}(\bar{z}) = \delta^{-2} \varphi(\bar{z}/\delta)$ is a set of cutoff functions. For $\omega \in L^1 \cap L^{\bullet}$, and small δ we have the following approximation.

$$\omega(\bar{z}) \approx \varphi_{\delta} * \omega(\bar{z}) \approx \sum_{i} \gamma_{i} \varphi_{\delta}(\bar{z} - \bar{z}_{i})$$
 (2.4)

for some points \bar{z}_i and constants γ_i . The first approximation in (2.4) is due to relation (2.3), the second approximation is simply a standard Riemann sum approximation to the convolution integral. We can chose the parameters \bar{z}_i and γ_i in various ways. One approach is to allow the points \bar{z}_i to lie initially on uniform grid of mesh size h, and to let $\gamma_i = h^2 \omega(\bar{z}_i)$; another is to let $\gamma_i = \int_{\bar{B}_i} \omega$, where \bar{B}_i denotes the square centered on \bar{z}_i , of area h^2 .

The "vortex blobs" or simply the "blobs" are represented by the vorticity distributions $\omega_i(\bar{z}) = \gamma_i \varphi_\delta(\bar{z} - \bar{z}_i)$ (here we have suppressed the δ dependence of ω). We say that the center of the i'th blob is \bar{z}_i and that its strength is γ_i .

Let G be the fundamental solution to Laplace's equation and let G_{δ} denote the smoothed kernel $G * \varphi_{\delta}$. Then we have by (2.1) and (2.4) that the stream function ψ satisfies the following relation.

$$\psi(\bar{z}) = G * \omega(\bar{z}) \approx \sum_{i} \gamma_{i} G * \varphi_{\delta}(\bar{z} - \bar{z}_{i}) = \sum_{i} \gamma_{i} G_{\delta}(\bar{z} - \bar{z}_{i}). \tag{2.5}$$

Thus an approximate velocity field is given by

$$u(\bar{z}) = -\psi_y = -\sum_i \gamma_i \frac{\partial G_{\delta}}{\partial y} (\bar{z} - \bar{z}_i)$$
 (2.6)

$$v(\bar{z}) = \psi_x = \sum_i \gamma_i \frac{\partial G_{\delta}}{\partial x} (\bar{z} - \bar{z}_i)$$
 (2.7)

We use these equations to update the positions of the vortex blobs. Specifically, we suppose that the positions, $\bar{z}_i(t) = (x_i(t), y_i(t))$, of the blobs satisfy the following first order system of differential equations.

$$\frac{dx_j}{dt} = -\sum_i \gamma_i \frac{\partial G_{\bar{o}}}{\partial y} (\bar{z}_j - \bar{z}_i)$$
 (2.8)

$$\frac{dy_j}{dt} = \sum_{i} \gamma_i \frac{\partial G_{\delta}}{\partial x} (\bar{z}_j - \bar{z}_i)$$
 (2.9)

If the cutoff functions have radial symmetry we introduce the functions $K(\bar{z})$ and $K_{\delta}(\bar{z})$ and the associated function $f_{\delta}(r)$ defined by,

$$K(\bar{z}) = (-\partial_{y} \cdot \partial_{z})G(\bar{z}) = \frac{1}{2\pi |\bar{z}|^{2}}(-y, z)$$
 (2.10)

$$K_{\delta}(\bar{z}) = (-\partial_{\nu} \cdot \partial_{z})G_{\delta}(\bar{z}) = f_{\delta}(|\bar{z}|)K(\bar{z})$$
 (2.11)

In table 1, we have displayed the functions f_{δ} associated with some of the more common choices of cutoff functions. Notice that $f_{\delta} \to 1$ as $\delta \to 0$ or as $r \to \infty$.

Using the functions defined above we can also rewrite equations (2.6) and (2.7) in the simpler form,

$$(u(\bar{z}),v(\bar{z})) = \sum_{i} \gamma_{i} K_{\delta}(\bar{z} - \bar{z}_{i}). \qquad (2.12)$$

The convergence of the vortex method for smooth initial data, using specific cutoff functions, has been proved. These results state that h, the initial distance between the blobs, and the cutoff δ , must satisfy a relation of the form $\delta = h^q$,

where 0 < q < 1. Hald and Del Prete [9] first showed convergence using their cutoffs. Hald [7] subsequently improved those estimates. Finally Beale and Majda [4] have shown that there existed cutoff functions which can produce arbitrary orders of convergence.

The solution of equation (2.2) is performed by the following random walk method. It is well known that the probability distribution of the positions of particles undergoing a Brownian motion, satisfies equation (2.2), the diffusion equation (see Lamperti [10]). The idea is to use a gaussianly distributed sequence to add a random component to the positions of the vortex blobs at each time step. This will simulate the diffusion inherent in equation (2.2).

An exact analysis of the Random Vortex method is not available. Marchioro and Pulvirenti [12] have shown that the random vortex method approximates the Navier-Stokes equations in a weak sense, as the initial grid size and cutoff approaches 0 in an appropriate way. Their result is impractical for us as we need stronger forms of convergence. For instance we need convergence in an L^2 or L^∞ sense. Hald [8] has analysed a method similar to the random vortex method for a problem in chemical dynamics and obtained strong convergence. His results though, and the discussions found in [5], [13] and [3] indicate that random vortex method will approximate the solution of the Navier-Stokes equation with an error of order $R^{-\frac{1}{12}}$. This is of the same order as the error obtained when equation (2.2) is solved by using random walks.

3. THE MODEL PROBLEM.

Our model problem consists of solving equations (1.1-4), together with the following non-smooth initial condition.

$$\xi(\bar{z}) = \begin{cases} 1 & \text{if } |\bar{z}| \le r \\ 0 & \text{if } |\bar{z}| \ge r \end{cases}$$
 (3.1)

As this is radially symmetric, it follows that $Q\omega = 0$ (see Batchelor [1]). The Navier-Stokes equations reduce to the two following cases.

$$\omega_t = 0 \quad \text{if } R = \infty \tag{3.2}$$

$$\omega_t = \frac{1}{R} \Delta \omega \quad \text{if } 0 < R < \infty \tag{3.3}$$

If $R=\infty$, the solution is a rotating blob with angular velocity $\frac{1}{2}$, and period of rotation, $T=4\pi$. For $0< R<\infty$, ω is a solution of the heat equation and can be written in the form,

$$\omega(\bar{z},t) = \int_{\mathbb{R}^2} E_t(\bar{z} - \bar{z}_1) \, \xi(\bar{z}_1) \, d\bar{z}_1 \tag{3.4}$$

where ξ is the initial distribution of vorticity and E_t is the fundamental solution for the heat equation,

$$E_t(\bar{z}) = \frac{R}{4\pi t} e^{\frac{-R|z|^2}{4t}} \tag{3.5}$$

One test of our numerical method is to calculate the exact and numerical values of vorticity (or velocity) at many points and use a discrete norm (e.g. L^2) to measure the error of the method. Another is to compare the positions of the vortex blobs at each time step with the positions of corresponding particles evolved by the exact flow. However, it is expensive to evaluate the integral (3.4) or to calculate the vorticity generated by the vortex blobs at many points. A common practice in these situations is to compare the exact and computed values of easily computable functionals. Following Milinazzo and Saffman, we chose the functional,

$$L(t) = \frac{\int\limits_{\mathbb{R}^2} |\bar{z}|^2 \omega(\bar{z}, t) d\bar{z}}{\int\limits_{\mathbb{R}^2} \omega(\bar{z}, t) d\bar{z}}$$
(3.6a)

which satisfies

$$L(t) = L(0) + \frac{4t}{R}$$
 (3.6b)

In the random vortex method L can be approximated by,

$$A(t) = \frac{1}{N} \sum_{i} \left[(x_i(t))^2 + (y_i(t))^2 \right]. \tag{3.7}$$

where the number of vortex blobs is denoted by N and their centers by $(x_i(t),y_i(t))$. The derivation of this formula supposes that all of the vortex blobs in our test problem have equal strength.

We will study the expected value, denoted by E, and the variance, denoted by σ^2 , of the quantity A. An exact analysis is not available, but we expect that the general case should be comparable to pure diffusion. It follows from section 4 of Milinazzo and Saffman [13] that,

$$E(A(t)) = A(0) + \frac{4t}{R}$$
 (3.8)

$$\sigma^{2}(A(t)) = \frac{8t}{RN}A(0) + \frac{8t^{2}}{NR^{2}}\left[5 - \frac{3\Delta t}{t}\right].$$

$$\approx \frac{8t}{RN}A(0)$$
(3.9)

for $\frac{R}{t} \gg 1$.

The analysis of the error is based on the functional L, and its numerical approximation A. We considered the relative error in A,

$$e(t) = \frac{|A(t) - L(t)|}{|L(t)|}$$
 (3.10)

Thus if the error in vorticity is small, then e(t) will be small.

The function e(t) can be decomposed into two components corresponding to "startup" and "interaction" errors. Thus we introduced the following functions.

$$e_{startup} = \frac{|A(0) - L(0)|}{|L(0)|}.$$
 (3.11)

$$e_1(t) = \frac{|A(t) - [A(0) + 4t/R]|}{|A(t)|}$$
 (3.12)

Note that A(0) + 4t / R is the expected value of A(t).

In the case of pure diffusion the standard deviaion $\sigma(e_1(t))$ can be calculated from (3.9), and saisfies

$$\sigma(e_1(t)) \approx \frac{1}{N^{\frac{1}{N}R^{\frac{1}{N}}}} \left[\frac{8t}{A(0)} \right]^{\frac{1}{N}}$$
(3.13)

for $R/t \gg 1$.

Milinazzo and Saffman [13] measure the error by comparing the functional,

$$g(t) = \frac{R}{4t} \left[L(t) - L(0) \right] \tag{3.14}$$

with its numerical approximation,

$$h(t) = \frac{R}{4t} \Big[A(t) - L(0) \Big]$$
 (3.15)

By (3.6b), g(t) = 1. Hence Milinazzo and Saffman's error is defined to be,

$$e_{MS}(t) = |g(t) - h(t)| = |1 - h(t)|.$$
 (3.16)

One can construct examples in which the vorticity field generated by a sum of vortex blobs approximates the exact vorticity for the model problem arbitrarily well, but for which the function $e_{MS}(t)$ is of arbitrary size (in particular, when R is large or t small). This is true even when t is restricted to a bounded interval 0 < t < T. Thus we prefer to estimate the error by $e_{startup}$ and e_1 .

4. NUMERICAL RESULTS.

There are many possible ways to implement the random vortex method. For example, we can make the following choices.

- (1) Initial placement and strength of the vortex blobs.
- (2) Form of cutoff function and the appropriate cutoff size for a given number of blobs.
- (3) Choice of time integration scheme to solve the system of equations, (2.8) and (2.9).
- (4) Random number generator to be used for the simulation of the random walk.

Our implementation consisted of the following choices.

The positions \bar{z}_i of the blobs were initially given on a uniform grid of mesh size h. Let B_i be the square of area h^2 centered on \bar{z}_i and parallel to the grid Then the strengths of each blob are given by, $\gamma_i = \int_{B_i} \omega$. We call this a uniform grid distribution with averaging.

For the cutoff function we chose Chorin's (see table 1) with a cutoff $\delta = 2h$.

To integrate equations (2.8) and (2.9), we used forth order Runge-Kutta as our time integration method. Milinazzo and Saffman [13] chose Heun's method, whereas Chorin [5] used Euler's method. With a time step of $\Delta t = 0.05$ the error in the O.D.E. was less than 10^{-15} .

To simulate the random walk we needed a 2-dimensional gaussianly distributed sequence of points (ζ_1, ζ_2) , with zero mean and standard deviation ν . This was obtained from a uniformly distributed sequence of points (η_1, η_2) , on $[0,1] \times [0,1]$, by using the formula,

$$\zeta_1 = \nu \cos(2\pi\eta_1) \left[-2\log(\eta_2) \right]^{\frac{1}{2}} \tag{4.1}$$

$$\zeta_2 = \nu \sin(2\pi\eta_1) \left[-2\log(\eta_2) \right]^{\frac{1}{2}} \tag{4.2}$$

Figures 1a-b and 2a-b display typical results. These figures display the computed value of A(t), compared to the expected value of A(t) and its standard deviation as given by (3.8) and (3.9) (R=1250, N=429 and 468, $\Delta t=0.05$). Notice that the error due to the random walk algorithm seems to dominate at this relatively low Reynolds number.

We then chose to study the effect of varying the initial distribution of the vortex blobs and the choice of cutoff function.

Following Milinazzo and Saffman [13] (p.384), we investigated the effect of distributing the blobs randomly at t = 0. The positions, \bar{z}_i , of the blobs were generated by using a uniformly distributed sequence of points on $[0,1] \times [0,1]$. This type of initial distribution will be called random and h will denote the average distance between the blobs.

The strengths of the blobs were determined by either:

(1) $\gamma_i = \int_{B_i} \omega$ where B_i denotes the square of area h^2 centered on \bar{z}_i ;

(2)
$$\gamma_i = \omega(\bar{z}_i) h^2$$

Figures 3a-d display the errors obtained for the various initial startup procedures discussed. Notice the particularly "choppy" behavior of figures 3c-d. This can be explained by observing that the addition of a blob at position (x,y) will change the error by an amount of N^{-1} (x^2+y^2) , N being the total number

of blobs for a particular run. Hence the addition of one blob near the boundary of the support of the vorticity will produce a change of 1/N.

Comparison of figures 3a-d clearly show the superior accuracy of the uniform grid with blob strength given by (1), in approximating the initial angular moment. Figures 4a-d show the error e(3T) (T is the period of the exact solution for the inviscid case.) obtained when blobs with the various initial conditions were moved by the random walk algorithm (R=1250). There is a correlation between the errors shown in figures 3a-b and figures 4a-b. This shows that for the methods where the blobs have a random distribution at t=0 the initial error, as measured by $e_{startup}$, dominates the error produced by the random walks. Figures 1b and 2b show that the phenomenon is less for a uniform initial distribution. Finally the averaging gives the best results. This is expected since the initial data is not smooth.

In the remainder of this paper we will use uniform grid and averaging.

For runs with Reynolds number greater than 20000, it became evident that even with the use of a uniform grid and averaging that startup errors were distorting the errors produced by the interaction of the two fractional steps of our method (see table 2). To eliminate the effect of the startup error we have introduced the functional e_1 .

We observed that practically any choice of cutoff function and cutoff δ , produced similar results. This seems to be due to our method of error measurement. If the vortex blobs have a distribution which is radially symmetry, then the velocity given by (2.12), for any cutoff, has no mean radial component. As our modes of error measurement only detect changes in radial positions (L depends on distance of blobs from origin) we cannot accurately compare different cutoff functions. On the other hand, we need a cutoff function if $R \neq \infty$, since the random walk algorithm allows blobs to become arbitrarily close.

Hence if no cutoffs were used, the interactions would be unstable. We chose Chorin's cutoff function with $\delta = 2h$.

From (3.13), for a fixed number of blobs we expect that the error would be $O(R^{-\frac{1}{2}})$, at least for pure diffusion. Table 2 tabulates our results as we increased Reynolds number from 1250 to 80000. With 500 blobs, note that the errors for the random vortex method are larger than those obtained for random walks, but that the rates of convergence as $R \to \infty$ are similar. These results verify the estimates established in section 3. for the random walk simulation.

How does the number of vortex blobs affect the accuracy. With R=1250, figures 4a, 5 and 6 contain the relevent data. First note that comparison of the plots in figures 5 and 6 indicate that the random vortex method and the random walk algorithm produce similar results. In addition though, observe the way in which the error decreased with an increase in the number of particles. Figure 4a shows that for the random walk algorithm it is unlikely for the error to be greater than the standard deviation $\sigma(e_1(3T))$. We did not observe an error greater than 2 times this quantity. The results for the random vortex method are similar.

5. CONCLUSION.

We have shown that the choice of initial conditions are important in the random vortex method. With our initial vorticity a uniform grid initial distribution of blobs with the strength of the blobs obtained by averaging give the best results. The dependence on the initial conditions is also found in the vortex method for smooth initial data. Here though, the choice of cutoff function is also important (Beale and Madja [4] and Perlman [14]).

The time evolution of the errors obtained for the random vortex method is similar to the corresponding result for the random walk algorithm. This is a favourable outcome. We could not expect the fractional step algorithm to

produce less error than its individual parts, but it was conceivable that the errors would be much larger.

Finally, our results indicate that the error for the random vortex method is $O(N^{-\frac{1}{N}}R^{-\frac{1}{N}})$. Hence, for a fixed number of vortex blobs, the accuracy of the method increases as the Reynolds number increases.

6. ACKNOWLEDGEMENTS.

I would like to thank Prof. A. Chorin for suggesting this problem and for his many helpful suggestions during the writing of this paper. In addition I thank Prof. O. Hald for his many insightful comments.

REFERENCES

- [1] G. K. Batchelor, An Introduction to Fluid Mechanics, Cambridge University Press (1967).
- [2] J. T. Beale & A. Majda, The Design and Analysis of Vortex Methods, *Proc. Conf.* on Transonic, Shock, and Multi-Dimensional Flows, Madison, Wisc. (May 1981).
- [3] J. T. Beale & A. Majda, Rates of Convergence for Viscous Splitting of the Navier-Stokes Equations, *Math. Comp.* v. 37 (1981), pp. 243-259.
- [4] J. T. Beale & A. Majda, Vortex Methods. II: Higher Order Accuracy in Two and Three Dimensions, *Math. Comp.*, v. 39 (1982), pp. 29-52.
- [5] A. J. Chorin, Numerical study of slightly viscous flows, *J. Fluid Mech.*, v. 57 (1973), pp. 785-796.
- [6] A. J. Chorin, J. Comput. Phys., v. 26 (1978), pp. 453-454.
- [7] O. H. Hald, Convergence of Vortex Methods for Euler's Equation. II, Siam J. Numer. Anal., v. 16 (1979), pp. 726-755.
- [8] O. H. Hald, Convergence of Random Methods for a Reaction-Diffusion equation, Siam J. Sci. Stat. Comput., v. 2 (1981), pp. 85-94.
- [9] O. H. Hald & V. Mauceri Del Prete, Convergence of vortex methods for Euler's equations, *Math. Comp.*, v. 32 (1978), pp. 791-809.
- [10] J. Lamperti, *Probablity Theory*, Benjamin, New York (1966).
- [11] A. Leonard, Vortex Methods for Flow Simulation, J. Comput. Phys., v. 37 (1980), pp. 289-335.

- [12] C. Marchioro & M. Pulvirenti, Hydrodynamics in Two Dimensions and Vortex Theory, *Commun. Math. Phys.*, v. 84 (1982), pp. 483-503.
- [13] F. Milinazzo & P. G. Saffman, The calculation of large Reynolds number two-dimensional flow using discrete vortices with random walk, *J. Comput Phys.*, v. 23 (1977), pp. 380-392.
- [14] M. Perlman, private communication, 1983.

Table 1. Common cutoff functions

Chorin [5]

$$f_{\delta}(r) = \begin{cases} 1 & \text{if } r > \delta \\ r / \delta & \text{if } r < \delta \end{cases}$$

Milinazzo and Saffman [13]

$$f_{\delta}(\tau) = \begin{cases} 1 & \text{if } \tau > \delta \\ \tau^2 / \delta^2 & \text{if } \tau < \delta \end{cases}$$

Hald [7]

$$f_{\delta}(r) = \frac{r^2}{\delta^2} \left[14 - 105 \frac{r^2}{\delta^2} + 196 \frac{r^3}{\delta^3} - 140 \frac{r^4}{\delta^4} + 36 \frac{r^5}{\delta^5} \right] \quad \text{if } r < \delta, 1 \text{ if } r > \delta.$$

$$f_{\delta}(r) = \frac{r^2}{3\delta^2} \left[56 - 630 \frac{r^2}{\delta^2} + 1568 \frac{r^3}{\delta^3} - 1680 \frac{r^4}{\delta^4} + 864 \frac{r^5}{\delta^5} - 175 \frac{r^6}{\delta^6} \right] \quad \text{if } r < \delta, 1 \text{ if } r > \delta.$$

Beale and Majda [4]

$$f_{\delta}(r) = 1 - e^{-\frac{r^2}{\delta^2}}$$

$$f_{\delta}(r) = 1 - 2e^{-\frac{r^2}{\delta^2}} + e^{-\frac{r^2}{2\delta^2}}$$

$$f_{\delta}(r) = 1 - (\frac{r^2}{\delta^2} + 1)e^{-\frac{r^2}{\delta^2}}$$

Table 2. Comparison of error measurements (in %), for runs with 500 blobs and with $e_{startup}$ = .239%.

Reynolds number	Random vortex method			Random walk algorithm		
	e (T)	e ₁ (T)	$e_{MS}(T)$	e (T)	e 1(T)	$e_{MS}(T)$
80000 20000 5000 1250	.408 .196 .453 1.46	.310 .536 1.08 2.00	324. 39.1 22.5 18.2	.602 .516 .362 .434	.108 .193 .354 .634	478. 103. 18.0 5.39

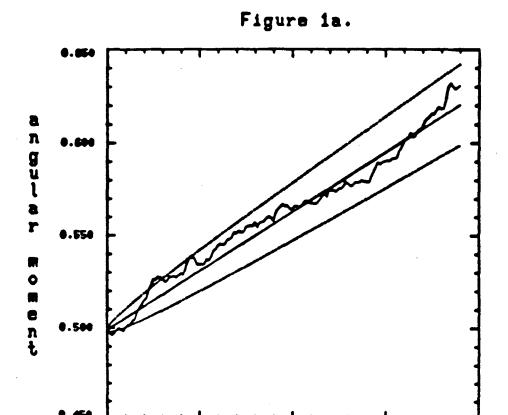
LIST OF FIGURE CAPTIONS.

Angular moment for 429 blobs evolved by random walk algorithm Figure 1a for R = 1250 together with its expected value and standard deviation. Figure 1b Angular moment for 468 blobs evolved by random walk algorithm for R = 1250 together with its expected value and standard deviation. Angular moment for 429 blobs evolved by random vortex method Figure 2a for R = 1250 together with its expected value and standard deviation. Figure 2b Angular moment for 468 blobs evolved by random vortex method for R = 1250 together with its expected value and standard deviation. Figure 3a Initial error, estartup, for uniformily distributed vortex blobs with averaged strength. Figure 3b Initial error, estartup, for uniformily distributed vortex blobs with unaveraged strength. Figure 3c Initial error, estartup, for randomly distributed vortex blobs with averaged strength. Figure 3d Initial error, estartup, for randomly distributed vortex blobs with unaveraged strength. Figure 4a Error e(3T) for uniformly distributed vortex blobs with averaged strength evolved by random walks with R = 1250. Error e(3T) for uniformly distributed vortex blobs with unaver-Figure 4b aged strength evolved by by random walks with R = 1250. Figure 4c Error e(3T) for randomly distributed vortex blobs with averaged strength evolved by by random walks with R = 1250. Figure 4d Error e(3T) for randomly distributed vortex blobs with unaveraged strength evolved by by random walks with R = 1250. Figure 5 Error e(3T) for uniformly distributed vortex blobs with averaged strength evolved by random walks with R = 1250.

Error e(3T) for uniformly distributed vortex blobs with averaged

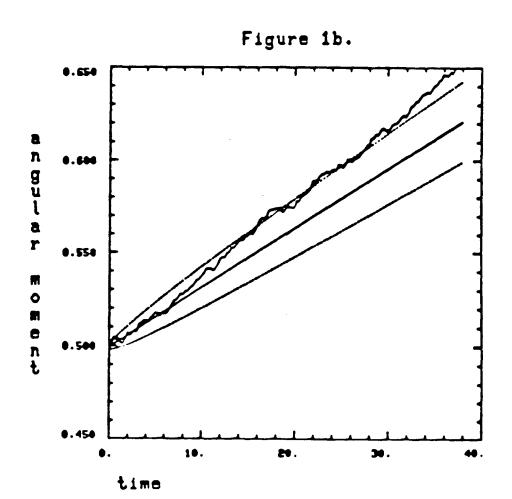
strength evolved by random vortex method with R = 1250.

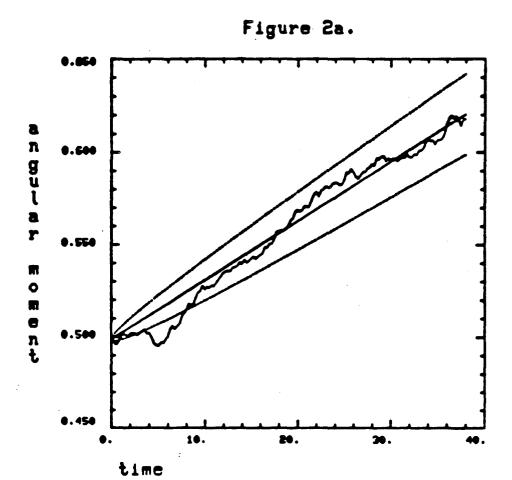
Figure 6



10.

time





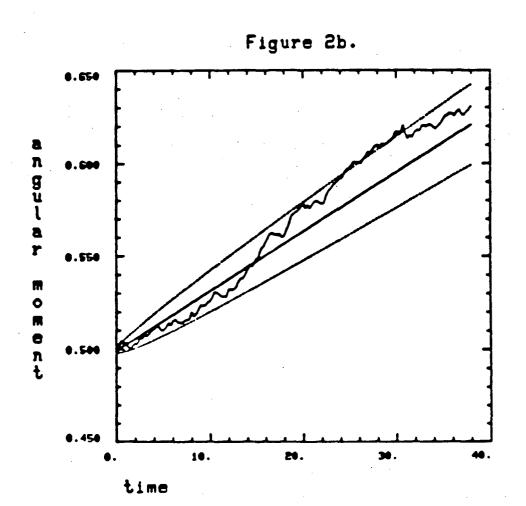


Figure 3a.

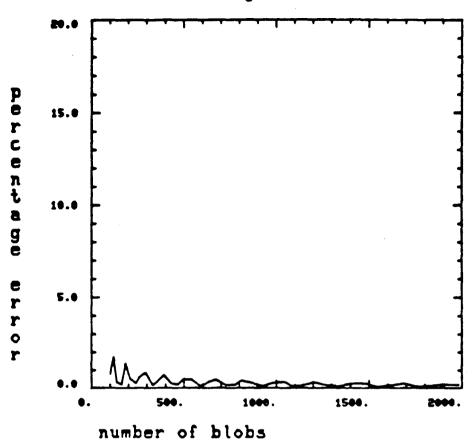
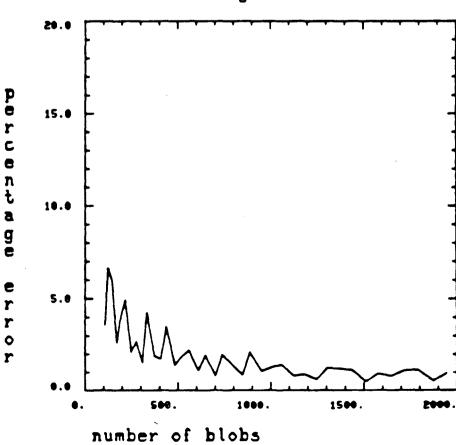
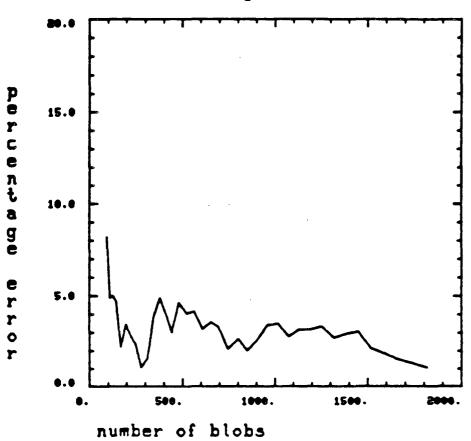


Figure 3b.







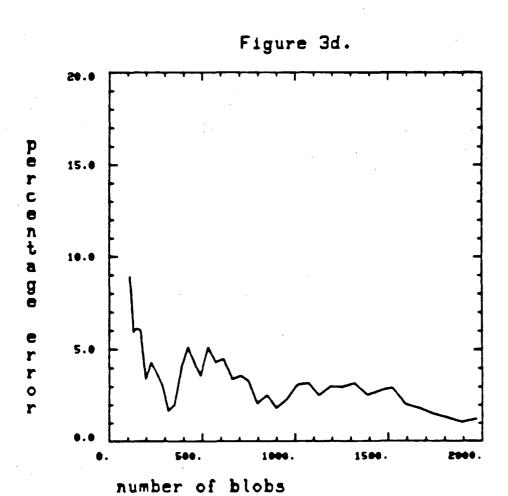


Figure 4a.

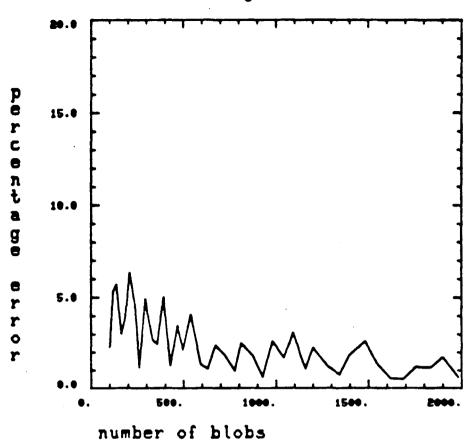
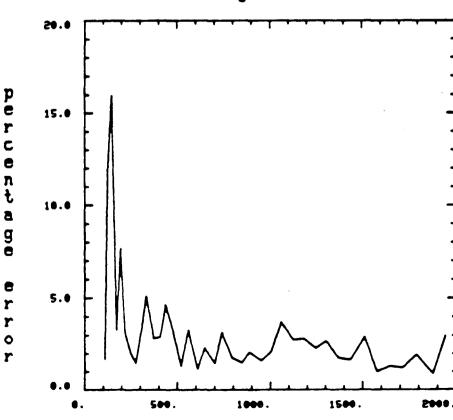
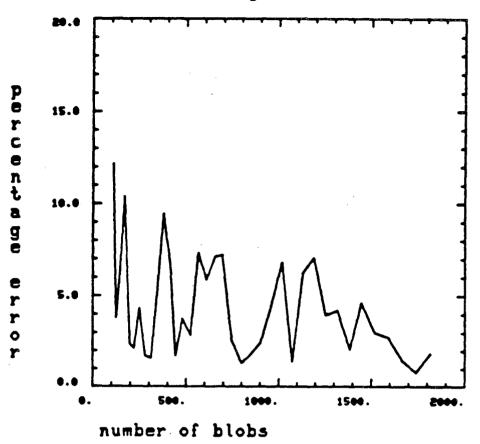


Figure 4b.



number of blobs





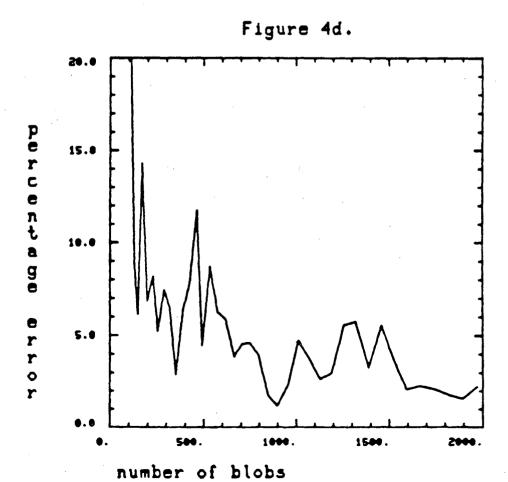


Figure 5.

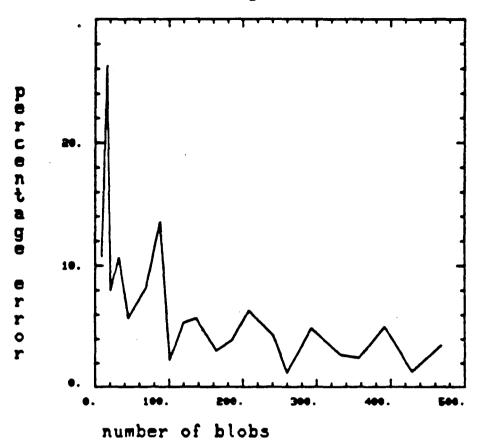
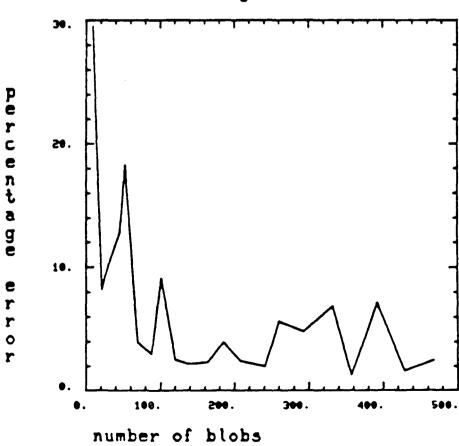


Figure 6.



This report was done with support from the Department of Energy. Any conclusions or opinions expressed in this report represent solely those of the author(s) and not necessarily those of The Regents of the University of California, the Lawrence Berkeley Laboratory or the Department of Energy.

Reference to a company or product name does not imply approval or recommendation of the product by the University of California or the U.S. Department of Energy to the exclusion of others that may be suitable.

TECHNICAL INFORMATION DEPARTMENT LAWRENCE BERKELEY LABORATORY UNIVERSITY OF CALIFORNIA BERKELEY, CALIFORNIA 94720