

Pattern Formation in a Simple Fluid System:

Taylor-vortex Flow

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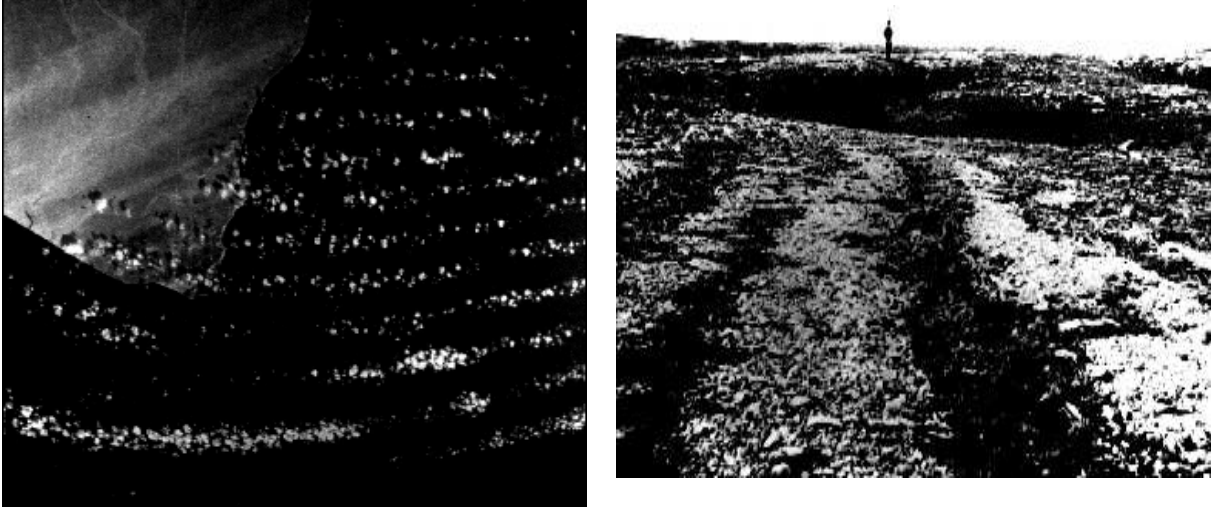


FIG. 1: Left: Cloud streets over the Arabian peninsular. Right: Permafrost stripes in Spitzbergen.

I. INTRODUCTION

In this experiment we are concerned with phenomena involving non-equilibrium non-linear systems. By “non-equilibrium” we mean systems that are driven away from the thermodynamic state, by the application of some stress parameter R such as a temperature difference, a shear rate, a concentration difference, or an electric voltage. Usually this leads to a current (for instance a heat, momentum, mass, or electrical current) passing through the system. This current produces dissipation. Thus, our systems are typically also dissipative. By that we mean that energy is converted at some rate, for instance by friction associated with the viscosity in fluid flow, into heat which is lost from the macroscopic dynamics of the system (such as the flow of a fluid). A particularly interesting aspect of nonlinear dissipative systems driven far from equilibrium is that they may undergo transitions from a temporally and/or spatially uniform state to a state of lower symmetry with characteristic temporal and/or spatial variations in the properties. Such transitions are called bifurcations.

In the first part of this experiment we shall study a bifurcation in a well defined laboratory system, namely the flow of a fluid between two concentric cylinders with the inner one rotating. This flow is known as Taylor-vortex flow (TVF). For careful laboratory study our system has the great virtue that it is relatively simple and that all factors influencing it can be carefully controlled. Consequently a complete theoretical analysis has been possible, allowing us to compare our results with quantitative theoretical predictions.



FIG. 2: Left: Striped pattern on the back of a tiger. Right: Zebra stripes.



FIG. 3: Left: Hexagonal permafrost patterns in Spitzbergen. Right: Hexagonal rock formation at Devil's Postpile in the Sierra Nevada.

Loosely, we refer to the spatial variation above the bifurcation as a “pattern”. Patterns can occur in one, two, or three dimensions. Here we shall study a one-dimensional pattern. The spatial variation of the pattern can be described by a superposition of plane-wave contributions of the form $A_q(\vec{r})\exp(i\vec{q}\cdot\vec{r})$ where \vec{r} is the position vector. A large part of the study of pattern formation involves the study of $\vec{q}(\vec{r}, t)$.

There are several reasons why non-equilibrium systems, and in particular pattern-forming systems, are of interest. One reason is that they are extremely common in the physical world which surrounds us. In nature, systems that can be well approximated by equilibrium thermodynamics are actually the exception, and most phenomena are associated with non-equilibrium states. In Figs. 1 to 3 several patterns that form in nature are illustrated.

Another reason why patterns are interesting to physicists is that the pattern-formation process is intrinsically nonlinear. By that we mean that the type of pattern which will form above a bifurcation under a given set of circumstances is determined by nonlinear terms in the equations of motion of the system. In the case of fluid systems you can see the non-linear term by going to Eq. 22 in the Appendix. There you find a contribution on the left-hand side that reads $(\vec{v} \cdot \vec{\nabla})\vec{v}$. This is a non-linear term because the fluid velocity \vec{v} occurs twice in it, i.e. the term is *quadratic* in the velocity. We know that, at least in principle, the linear problems of Physics are essentially soluble; but there is no general method by which nonlinear problems can be solved. In many cases the nonlinearities lead to qualitatively new physical phenomena, not present in linear systems, such as localized structures (solitary waves) and spatial or spatio-temporal chaos. Thus there is an important challenge here to physicists. An important way in which nonlinear non-equilibrium systems differ from equilibrium systems is that there is no extremum principle that enables us to predict the steady state. So there is no unique way to determine for instance the wave length of the striped patterns in Fig. 1 or 2 or the size of the hexagons in Fig. 3. In the second part of our experiment we shall see that we can produce Taylor vortices with wave lengths that span a continuous range even though the external conditions are fixed. This differs for instance from a crystalline sample in equilibrium; this sample usually would have a unique inter-atomic spacing. It is a challenging problem in non-linear physics to determine the permitted range of wave lengths, or to predict the unique wave length that might be chosen under very special circumstances.

While our system may seem relatively simple compared to the examples occurring in nature and shown in Figs. 1 to 3, it should always be kept in mind that the observed phenomena are of far broader applicability and pertain to a large class of nonlinear dissipative non-equilibrium systems. Other examples of such systems that have been studied in the laboratory include convection in a thin horizontal layer of fluid heated from below (Rayleigh-Bénard convection), dendritic crystal growth, flame fronts, certain chemical reac-

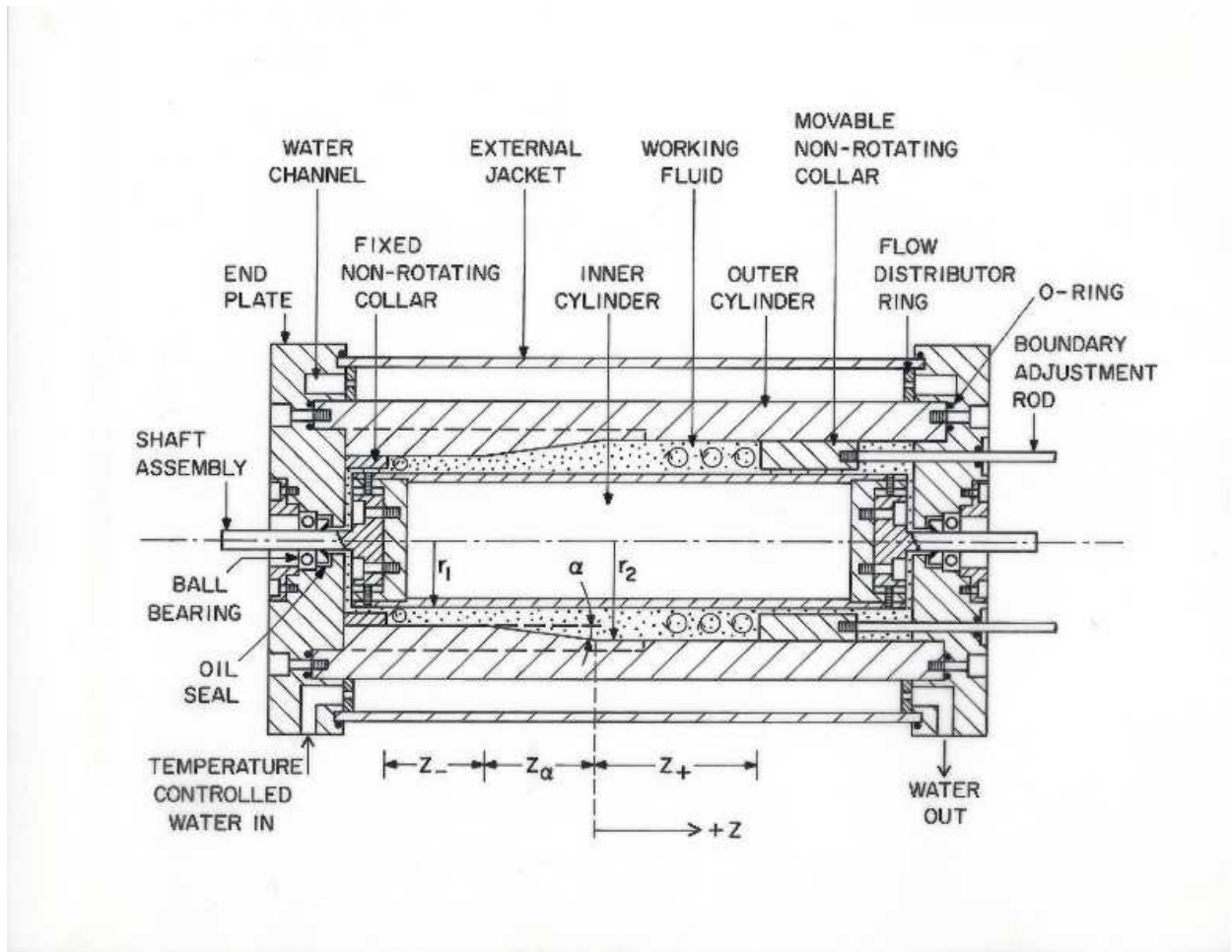


FIG. 4: Cross-sectional diagram of the apparatus. From Ref. [3].

tions, electric currents in semiconductors, convection in a nematic liquid crystal driven by an applied voltage, and even biological systems.

II. TAYLOR-VORTEX FLOW

In an Appendix you can find a derivation of the equations of fluid mechanics that describe our experiment. They follow very elegantly from the fundamental principles of mass, momentum, and energy conservation. If you are interested, and one hopes that you are, you should study this Appendix. But you do not absolutely have to do this in order to carry out this experiment.

A. The system

The apparatus used in our experiment is illustrated in Fig. 4. The fluid is contained between two concentric cylinders. In the illustration the left part of the outer cylinder has a varying radius. We will not investigate this case, and instead concentrate on the geometry where both cylinders have uniform radii. The apparatus also has an outer jacket that can be used to circulate temperature-controlled water. We will not use this feature either because our laboratory temperatures are sufficiently stable anyway.

A great richness of phenomena can be produced by rotating both the inner and the outer cylinder in various relative directions and at various rates.[1] We will consider only the case of a rotating inner cylinder, with the outer one stationary. In this case the angular velocity ω of that cylinder determines the distance from equilibrium. When ω is less than a critical value ω_c , the fluid flow will be purely azimuthal, and the velocity field $\vec{v}(r, \phi, x) = (v_r, v_\phi, v_x)$ has $v_x = v_r = 0$. This state is known as circular Couette flow and is described completely by $v_\phi(r)$. Above ω_c , however $\vec{v}(r, \phi, x)$ is periodic in the axial direction, but is unaltered by rotation around the axis. Thus, the pattern $\vec{q}(x)$ may be regarded as vortex rings stacked one next to the other. The direction of circulation within the vortices alternates between successive vortices. Thus one wave length $\tilde{\lambda}$ of the pattern includes two vortices (a vortex pair). The pattern is one-dimensional in the sense discussed earlier. For the axially infinite system, the theoretical analysis of the bifurcation from Couette to vortex flow was provided by G.I. Taylor. Taylor also performed extensive early experiments, which largely supported his theoretical findings.[2]

B. Dimensionless parameters that describe our apparatus and pattern

It is useful to describe the geometry of the apparatus and the cylinder speed in dimensionless form. That way, when these dimensionless parameters are the same for two experimental setups, we expect the same phenomena and flow states to occur. For instance, we might expect the same dimensionless wave lengths of the vortices.

The geometry is easy to describe. For the physical system the ends are usually provided by two annular, non-rotating rigid inserts a distance ℓ apart. Thus we have an aspect ratio

$$L = \ell/d \tag{1}$$



FIG. 5: A photograph of a portion of the apparatus with Taylor vortices present..

that describes the length the system. Here $d = r_2 - r_1$ is the gap between the inner and outer cylinder which have radii $r_1 = 1.867$ cm and $r_2 = 2.499$ cm respectively. An additional geometric parameter is the radius ratio

$$\eta = r_1/r_2 \quad (2)$$

which for our apparatus is equal to 0.747.

For the inner-cylinder speed a somewhat more complicated dimensionless combination of parameters is used because it simplifies the equations of motion for the Taylor-Couette geometry. It is now known as the Taylor number and is given by

$$T = \frac{(1 - \eta)^5(1 + \eta)}{2\eta^2} \beta^2 \quad (3)$$

where

$$\beta = \frac{2r_1^2\omega}{\nu(1 - \eta^2)}. \quad (4)$$

Here ν is the kinematic viscosity (the kinematic viscosity is equal to the shear viscosity divided by the density and has the units of length²/time).

We have a choice of r_1 , r_2 , or $r_2 - r_1$ for scaling the wave length λ of the vortices that will form in our experiment. It is easy to see that $r_2 - r_1$ would not be a good choice because it vanishes as $\eta \rightarrow 1$, causing λ to diverge. Similarly r_1 is not so good because it vanishes as $\eta \rightarrow 0$. Thus, the choice we find in the literature is r_2 :

$$\lambda = \tilde{\lambda}/r_2 \quad (5)$$

where $\tilde{\lambda}$ is the wave length in the same units of length as those of r_2 (e.g. cm). Often it is more convenient to describe the vortex pattern in terms of the dimensionless wave numbers

$$q = 2\pi/\lambda . \quad (6)$$

C. The circular Couette flow

It is relatively easy to determine a solution for the velocity field of the fluid analytically if we ignore the influence of the ends (in the axial direction) on the flow. This is done in more detail in the Appendix. There we find that

$$v_\phi(r) = ar + b/r \quad (7)$$

where all quantities still have their natural physical units. The constants a and b are determined by the boundary conditions. For rigid BC's we have $v_\phi(r_1) = r_1\omega_1$ and $v_\phi(r_2) = r_2\omega_2$. In our case we have, for a stationary outer cylinder, $\omega_2 = 0$, and thus $v_\phi(r_2) = 0$. Dropping the subscript on ω_1 , this yields

$$a = -\omega r_1^2/(r_2^2 - r_1^2) \quad (8)$$

and

$$b = \omega r_1^2 r_2^2 / (r_2^2 - r_1^2) \quad (9)$$

where ω is now the angular speed, in radian/sec, of the inner cylinder.

Exercise 1: For Eqs. 7 to 9, show that $v_\phi = 0$ for $r = r_2$ and $v_\phi = \omega r_1$ for $r = r_1$.

Exercise 2: For $\omega = 1$ radian/sec and the cylinder radii of your apparatus, make a graph of $v_\phi(r)$ and put it in your notebook. Comment on what you find.

Your apparatus is a precision instrument for studying TVF. As mentioned above, it has a jacket through which temperature-controlled water can be circulated so as to keep the fluid temperature constant. This is useful both because the viscosity is temperature dependent, and because convection due to horizontal temperature gradients should be prevented, but not necessary for our work (but you should note the laboratory temperature while you do your experiment). The gap between the cylinders is uniform to within 0.003 cm, and the cylinder axis is straight to this accuracy. The cylinder speed is controlled extremely well by a function synthesizer and a stepper motor that has a rotation rate precisely synchronous with the synthesizer. Thus, the boundary conditions and the stress parameter are known with high accuracy in these experiments. To provide a better idea of how the apparatus performs, Fig. 5 shows a photograph of its top part. In that case, $\omega > \omega_c$ and the Taylor vortices are apparent.

We will want to study the vortices above onset, and thus it is useful to define a parameter that tells us how far above onset we are. A convenient parameter for this is $\epsilon \equiv \omega/\omega_c - 1$. Thus, when $\epsilon < 0$ we find the Couette state, and for $\epsilon > 0$ the vortices exist.

D. Stability of circular Couette flow

In Sect. II A we found the circular Couette solution Eqs. 7 to 9 of the flow between our cylinders. This solution exists for all ω . However, we do not know whether it is stable. If a small disturbance were to be applied to it, this disturbance might grow and a completely new flow state might evolve. It turns out that such disturbances always decay when ω is less than a critical value ω_c corresponding to a critical value T_c of the Taylor number, but that they grow for larger ω . So the Couette state is stable below ω_c , but at the bifurcation point $\omega = \omega_c$, an exchange of stability takes place between the circular Couette state and a new state that turns out to be the Taylor-vortex state shown in Fig. 5.

The precise value of the critical Taylor number depends on η . When it is reached, Taylor vortices of a unique wave length λ_c will first form. The corresponding critical wave numbers q_c and the values of T_c depend on η , and are given in Table I.

TABLE I: The critical wave number q_c and the critical Taylor number T_c as a function of the radius ratio η for Taylor-vortex flow in an infinitely long system.

η	q_c	T_c
0.975	3.127	1723.9
0.900	3.129	1823.4
0.800	3.133	1994.3
0.700	3.139	2230.1
0.600	3.148	2571.6
0.500	3.162	3099.6

III. THE EXPERIMENT

A. Experimental details

As you can see from Eq. 4, the Taylor number depends on the kinematic viscosity ν of the fluid. In case you don't remember, ν is related to the shear viscosity η (not to be confused with the radius ratio!) by $\nu = \eta/\rho$ where ρ is the density of the fluid. The unit of ν is the stokes or St, equal to 1 cm²/sec. The reference value that often is used for comparison is the viscosity of water at 20°C, which is 1.0038 cSt (centi-stokes).

Exercise 3: Calculate the critical value $\tilde{\lambda}_c$, in cm, and the critical value ω_c , in radians per second, for your apparatus, assuming that you use pure water at 20°C as the fluid. You will have to use the data in Table I, Eqs. 3 and 4, and remember how everything is scaled.

The fluid is contained in an overflow volume mounted on a column next to the Taylor-vortex apparatus. You will have to take this volume to the sink and prepare your working fluid in it.

The inner cylinder is driven by a stepper motor in a so-called micro-stepping mode. This is done with a drive that has to be supplied with a square wave that goes from a low voltage (typically 0) to a voltage of at least 1.8 and no more than 5 volts. The motor will run at a speed of 1 Hz when the drive frequency is 10 kHz, i.e. at 1/10,000 the drive frequency.

B. Onset of Taylor vortices

Well, it looks like we are about ready to see whether all that theory is really correct. So let us prepare the apparatus. Drain whatever liquid may be in it into the overflow volume. You will have to take this volume to the sink and discard the old fluid. After rinsing it, fill the overflow volume with some de-ionized (DI) water and use this to flush out the apparatus a time or two. Who knows what was last in there, and we do not want our fluid to be contaminated! Now prepare one liter of DI water and 20 cc of Kalliroscope concentrate (i.e a 2% mixture) in a beaker. Pour it into the overflow volume, attach to the apparatus, and fill the apparatus. Again flush it once or twice with the mixture to be sure you have a uniform fluid in the apparatus.

You already calculated the angular speed, so you know about how fast to drive the motor to find the onset of Taylor vortices. Start at a speed well below that. At each speed, wait long enough to give potential vortices a chance to form. Make sure you record the lab temperature that prevails during your experiment; viscosities are temperature dependent!

At each speed, make a note in your notebook about whether vortices formed. Pay particular attention to where they first form. The theory pertains to an imagined system of infinite length without ends. Where would this be most likely to be a good approximation for your real apparatus?

Exercise 4: Why does it take so long for the vortices to form even after you have exceeded the critical angular speed?

Exercise 5: Can you explain the strong vortices that form first at the ends?

After you have determined the onset frequency and have a nice pattern of vortices in the apparatus, count how many there are. Also measure the distance between the ends of the liquid column (between the white teflon rings) and calculate the wave number of the vortices. How does this compare with the theory?

Now reverse the process and reduce the cylinder speed in small steps. You will have to be patient at each step to be sure whether stable vortices still exist. Does the onset you found going in this direction agree with what you found going up?

C. Eckhaus Boundary

Once you have established a nice vortex pattern above but close to onset, count the vortex pairs and then increase the speed suddenly to twice the onset speed, i.e. to $\epsilon \simeq 1$. Do the vortices survive this harsh treatment?

Now turn off the motor and let the vortices decay completely. Then go again suddenly to $\epsilon \simeq 1$. Watch the vortices grow and describe what you see. When everything settles down to a time independent pattern, count the vortices and calculate their wave number. Now reduce ϵ to about 0.2. Let the pattern equilibrate. Do you still have the same wave number? Continue to reduce ϵ in small steps and each time wait long enough to convince yourself that you have a stable pattern. What do you find?

In the above experiment you should have observed that the pattern with the wave number you generated by the sudden start experiment becomes unstable as you approach onset from above. A pair of vortices will be lost after one of your small steps.

Question: Where are the vortices lost? In the interior or at the ends?

This vortex-pair loss is due to the famous Eckhaus instability that occurs in many pattern-forming systems. Hopefully you did the above experiment slowly enough so that you lost only a single vortex pair. You will have to let the system equilibrate for some time to get a reasonably uniform wave number throughout; where you lost the pair the wave number initially will be too small. Then reduce the speed further to see whether you can find additional transitions. For each number of vortices in your system you can calculate a wave number at which the Eckhaus instability occurs. Make a graph of this in your notebook. How do your results compare with what you can find in the literature?

D. Secondary Bifurcations

When the speed of the inner cylinder gets large enough, the Taylor vortices become unstable to a time-periodic mode. Waves form on each vortex and travel azimuthally. A bifurcation to a time periodic state is called a Hopf bifurcation. Let us see whether we can find it.

Increase the speed to about two or three times critical and let the system equilibrate. If nothing happens, go up further in fairly large steps. Once you find this so-called wavy

mode, back off and see the minimum speed at which it survives. When it is gone, increase the speed in smaller steps and get a good value for ϵ_w at the onset.

Now that you have this wavy state, try to determine some of its properties. How many wave lengths are there around the circumference? What is the wavy-mode speed relative to the inner-cylinder speed?

E. Viscosity measurements

When you are done, drain the fluid from the apparatus into your overflow volume. Fill the viscometer with a small amount of the fluid and determine the viscosity. Did the 2% Kalliroscope that you added to the water alter the viscosity significantly? If the viscosity differs from what you used when you calculated ω_c , re-do the calculation and see whether it improved the agreement of your measurement with theory or whether it makes it worse.

F. Dependence on Viscosity

It would be nice to know whether the theory is right about the viscosity dependence of the onset. So let us change ν and do the above experiment again. For this, use a mixture of 50% water and 50% glycerol. Here too add about 20 cc of Kalliroscope concentrate to one liter of the mixture. Do make an estimate of the onset frequency *before* you start the experiment. You will need to know the viscosity of the mixture. So you better measure it. Discuss the agreement or disagreement of your results with the theory. Don't forget to see also how the wave number compares with what you found before and with theory.

IV. APPENDIX: THE EQUATIONS OF MOTION

A. Number of independent variables

Since we are going to use a fluid mechanical systems to investigate various bifurcation and pattern-formation problems, it seems appropriate to understand first the origin of the equations of motion for fluid flow. These equations are derivable from the conservation laws of Newtonian mechanics.

For a single-component system in thermodynamic equilibrium there are three extensive or intensive variables and one relation between them [e.g. the entropy, volume, and number of moles S, V, N , and the energy $U(S, V, N)$, or their conjugates and the appropriate Legendre transform of U]. Thus, there are only two independently specifiable variables needed to determine the state of the system. In fluid mechanics it is often convenient to use the pressure P and the density ρ .

A flowing fluid is obviously not an equilibrium state. Nonetheless, it is considered to be an excellent approximation to assume that it is *locally* in equilibrium. However, the two thermodynamic variables are now functions of the position \vec{r} and of the time. If the fluid is in motion, it is necessary to specify also the velocity \vec{v} in addition to the two thermodynamic variables. The velocity generally has three components, which may also be functions of \vec{r} and t . Thus, $\vec{v} = \vec{v}(\vec{r}, t)$, or

$$\vec{v}(x, y, z, t) = u(x, y, z, t)\hat{i} + v(x, y, z, t)\hat{j} + w(x, y, z, t)\hat{k}$$

where \hat{i} , \hat{j} , and \hat{k} are the unit vectors in the x , y , and z directions. Specifying the five scalar variables P , ρ , u , v , and w , all as a function of x , y , z , and t , completely describes the state of the system. Thus, we need five scalar equations to completely describe the fluid flow.

B. Equations of Motion.

Since there are five independent variables, we need five equations which relate them to each other. One of them is the equation of continuity which corresponds to mass conservation. The next three can best be written as a single vector equation corresponding to Newton's Second Law. For a so-called ideal fluid, in which there is no dissipation due to viscosity or thermal conduction, this equation is known as Euler's equation. When dissipation is taken into consideration, as it must be for real fluids in most cases of interest to us, then this equation becomes the Navier-Stokes equation. The fifth equation is one which assures entropy or energy conservation. It will not be important to our particular case of Taylor-vortex flow because our experiments are done at constant temperature. But for completeness, we derive it below as well.

1. *The continuity equation.*

The velocity $\vec{v}(x, y, z, t)$ is the velocity of the fluid at a *particular point fixed in the laboratory reference frame*. Thus, \vec{v} does *not* represent the velocity of a particular small volume element of the fluid as it moves along. Rather it refers to the velocity of that fluid material which happens to be present at a particular location at a particular time, and this fluid material may be changing with time. Similarly, P and ρ refer to a particular point in space.

Let us consider some macroscopic volume V_0 which is fixed in space. The total mass in this volume is

$$M = \int \rho dV \quad (10)$$

where the integral is over the entire volume V_0 . The mass contained in V_0 may change with time because fluid is flowing into and out of V_0 . Through a small element da of the surface which bounds V_0 , there is a mass flux *out of* V_0 given by

$$dm = \rho \vec{v} \cdot d\vec{f} \quad (11)$$

where $d\vec{f} = \hat{n} da$, and where \hat{n} is the unit vector normal to the surface element da and pointing in the outward direction (see Fig. 6). This is easily seen as follows. The dot-product $\vec{v} \cdot d\hat{n}$ gives the velocity component in the direction of the normal, and multiplying it by da yields the volume flow-rate out of V_0 through da . Multiplying further by ρ leads to the corresponding mass flow rate dm . Note that dm is by convention taken to be positive when the flow is out of V_0 . The *total* mass flux *out of* V_0 is given by the integral of dm over the surface bounding V_0 , that is by $\int_S \rho \vec{v} \cdot d\vec{f}$. Using the divergence theorem, this can be written as the volume integral :

$$\int_S \rho \vec{v} \cdot d\vec{f} = \int \vec{\nabla} \cdot (\rho \vec{v}) dV . \quad (12)$$

But with Eq. 10, the total mass flux *out of* V_0 is also given by $\partial M / \partial t = -(\partial / \partial t) \int \rho dV$ with this last integral also over the entire volume V_0 . Thus we have

$$(\partial / \partial t) \int \rho dV = - \int \vec{\nabla} \cdot (\rho \vec{v}) dV . \quad (13)$$

Rearranging, one has

$$\int [\partial \rho / \partial t + \vec{\nabla} \cdot (\rho \vec{v})] dV = 0 . \quad (14)$$

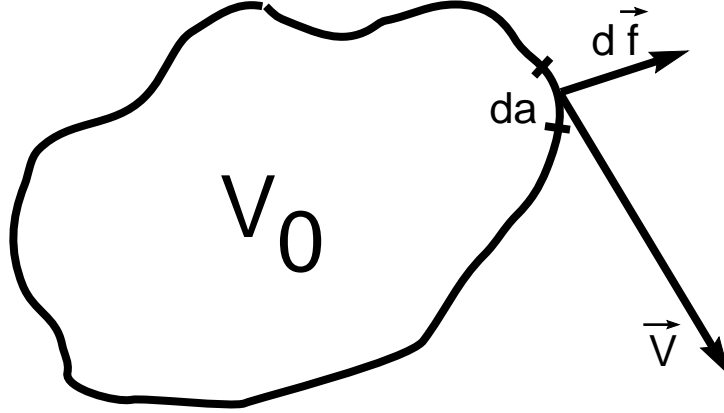


FIG. 6: Schematic illustration of the derivation of the continuity equation.

Since Eq. 14 must hold for any volume V_0 , it follows that

$$\partial\rho/\partial t + \vec{\nabla} \cdot (\rho\vec{v}) = 0 . \quad (15)$$

Equation 16 is the continuity equation in its most general form. It simplifies considerably when the density is constant throughout the fluid. In that case it is easy to see that Eq. 16 becomes

$$\vec{\nabla} \cdot \vec{v} = 0 . \quad (16)$$

2. Euler's equation and the Navier-Stokes equation.

In order to find the acceleration of a small fluid volume element, we consider the force $-Pd\vec{f}$ which acts *inwardly* (remember how $d\vec{f}$ was defined!) on the area da . Integrating over the surface bounding the volume and using the divergence theorem, we obtain

$$- \int_S Pd\vec{f} = - \int \vec{\nabla} P dV . \quad (17)$$

for the total force exerted by the surrounding fluid upon V_0 . Thus, the force acting on a small volume element dV containing a mass dM is $-(\vec{\nabla}P)dV$. Equating this with $(dM)\vec{a}$ where $\vec{a} = d\vec{v}/dt$ is the acceleration of the volume element, we have $(dM)(d\vec{v}/dt) = -(\vec{\nabla}P)dV$, or

$$\rho(d\vec{v}/dt) = -\vec{\nabla}P \quad (18)$$

where we have used $\rho = dM/dV$.

It is crucial to recognize that the derivative on the lhs of Eq. 18 is the *total* derivative of \vec{v} . Thus, it does *not* represent the change in \vec{v} with respect to t at a fixed location. Rather, in order to conform to Newton's Second Law, it represents the acceleration of a particular element of fluid material and thus corresponds to the change of \vec{v} in a frame *moving along* with this particular fluid element. Therefore, in Cartesian coordinates for instance,

$$d\vec{v} = (\partial\vec{v}/\partial t)dt + (\partial\vec{v}/\partial x)dx + (\partial\vec{v}/\partial y)dy + (\partial\vec{v}/\partial z)dz \quad (19)$$

or

$$d\vec{v} = (\partial\vec{v}/\partial t)dt + (d\vec{r} \cdot \vec{\nabla})\vec{v} . \quad (20)$$

With $d\vec{r}/dt = \vec{v}$ this gives

$$d\vec{v}/dt = \partial\vec{v}/\partial t + (\vec{v} \cdot \vec{\nabla})\vec{v} . \quad (21)$$

Substituting into Eq. 18, one has

$$\partial\vec{v}/\partial t + (\vec{v} \cdot \vec{\nabla})\vec{v} = -(1/\rho)\vec{\nabla}P . \quad (22)$$

Equation 22 is Euler's equation, and represents Newton's Second Law (or linear momentum conservation) for the ideal fluid in which there is no dissipation and no external force. If the fluid is subjected to the gravitational field, there is an external force equal to $\rho\vec{g}$ which must be taken into account, yielding

$$\partial\vec{v}/\partial t + (\vec{v} \cdot \vec{\nabla})\vec{v} = -(1/\rho)\vec{\nabla}P + \vec{g} . \quad (23)$$

The second term on the lhs of Eq. 22 or 23 is nonlinear in \vec{v} . It makes it impossible to solve the equations of motion under most circumstances in a straight forward manner, and is responsible for most of the phenomena which occur in fluid flow and which interest us here.

For real fluids, it is necessary to consider dissipation which was neglected in Euler's equation. We will not do that in detail here, but it can be accomplished by considering the momentum flux $\partial(\rho\vec{v})/\partial t$. As derived from Euler's equation, this flux is associated entirely with reversible processes. If there is irreversibility due to viscosity, the corresponding term can be subtracted. For an *incompressible* fluid in the gravitational field, this procedure leads to[5]

$$\partial\vec{v}/\partial t + (\vec{v} \cdot \vec{\nabla})\vec{v} = -(1/\rho)\vec{\nabla}P + (\eta/\rho)\Delta\vec{v} + \vec{g} . \quad (24)$$

Here Δ is the three-dimensional Laplace operator $[(\partial^2/\partial x^2 + \partial^2/\partial y^2 + \partial^2/\partial z^2)$ in Cartesian coordinates], and η is the shear viscosity (in the case of compressible fluids where the density may vary, there is an additional contribution involving the bulk viscosity ζ ; but we will have no need to consider this dissipative mechanism). Equation 24 is the Navier-Stokes equation for an incompressible fluid, and for us it will serve as the equation of motion appropriate for real fluids in the presence of gravity.

When we want to consider a fluid in the presence of temperature gradients, it is convenient to transform Eq. 24 as follows. We write the temperature $T(x, y, z, t)$ as

$$T(x, y, z, t) = T_0 + T_1(x, y, z, t) \quad (25)$$

where T_0 is a constant mean temperature. Similarly, we write the density as

$$\rho = \rho_0 + \rho_1 \quad (26)$$

and the pressure as

$$P = P_0 + P_1 . \quad (27)$$

For ρ_0 we take the constant density corresponding to T_0 and the equation of state. For small T_1/T_0 we then have

$$\rho_1 = -\rho_0\beta T_1 \quad (28)$$

where β is the isobaric thermal expansion coefficient $-(1/\rho)(\partial\rho/\partial T)_p$. For P_0 we take the pressure corresponding to mechanical equilibrium when $T = T_0$ and $\rho = \rho_0$. It is given by Eq. 23, and in the presence of gravity will vary with height according to

$$P_0 = \rho_0\vec{g} \cdot \vec{r} + constant . \quad (29)$$

To lowest order in ρ_1, T_1 , and P_1 the pressure term in Eq. 24 can now be written as

$$\vec{\nabla}P/\rho = \vec{\nabla}P_0/\rho_0 + \vec{\nabla}P_1/\rho_0 - (\vec{\nabla}P_0/\rho_0^2)\rho_1 \quad (30)$$

or

$$\vec{\nabla}P/\rho = \vec{g} + \vec{\nabla}P_1/\rho_0 + \vec{g}T_1\beta . \quad (31)$$

With this, the Navier-Stokes equation, Eq. 24, becomes

$$\partial\vec{v}/\partial t + (\vec{v} \cdot \vec{\nabla})\vec{v} = -(1/\rho_0)\vec{\nabla}P_1 + \nu\Delta\vec{v} - \beta T_1\vec{g} \quad (32)$$

where we have introduced the kinematic viscosity

$$\nu = \eta/\rho . \quad (33)$$

3. *The equation of continuity of the entropy or the energy.*

In an ideal fluid there cannot be any heat transport from one part to another, and the entropy per unit mass s of any fluid particle remains constant at all times; *i.e.* $ds/dt = 0$. By arguments similar to those used to derive the continuity equation Eq. 16, one has

$$\partial(\rho s)/\partial t + \vec{\nabla} \cdot (\rho s \vec{v}) = 0 . \quad (34)$$

If by virtue of the initial conditions s is uniform throughout the fluid, it will remain uniform at all times. Then the motion is isentropic and Eq. 34 becomes $s = \text{constant}$. Equations 16, 32, and 34 give the five scalar equations which determine the motion of the ideal fluid.

Alternatively to Eq. 34, we could have focused on an equation for the energy flux. We would again start with the ideal fluid, and consider some volume V_0 . We equate the time rate of change of the energy in this volume with the flux across the surface bounding the volume. The total energy in V_0 can be written as a sum of the kinetic and the internal energy, and its time rate of change is given by $(\partial/\partial t) \int (\rho v^2/2 + \rho \epsilon) dV$, where ϵ is the internal energy per unit mass. Equating with the flux across the surface, we get

$$(\partial/\partial t) \int (\rho v^2/2 + \rho \epsilon) dV = - \int_S \rho (v^2/2 + \epsilon + P/\rho) \vec{v} \cdot d\vec{f} . \quad (35)$$

On the rhs, the contribution P/ρ corresponds to the work done by pressure forces on the volume V_0 . Using the divergence theorem, this gives

$$(\partial/\partial t) \int (\rho v^2/2 + \rho \epsilon) dV = - \int \vec{\nabla} \cdot \rho (v^2/2 + \epsilon + P/\rho) \vec{v} dV . \quad (36)$$

Since Eq. 36 must hold for any V_0 , we have

$$(\partial/\partial t)(\rho v^2/2 + \rho \epsilon) = - \vec{\nabla} \cdot \rho (v^2/2 + \epsilon + P/\rho) \vec{v} . \quad (37)$$

Equations 34 and 37 should not be thought of as independent of each other. Indeed, one can be derived from the other because the energy and entropy are not independent thermodynamic variables. Thus there remain only five relationships between the five scalar fields of hydrodynamics.

Equations 34 and 37 contain only the reversible fluxes. When the fluid is not ideal, they must be modified by including dissipative terms. For our purposes, the most important contribution will come from the conduction of heat when the temperature is not uniform.

In that case, there is a heat-flux density \vec{q} which can be expanded in a power series of the temperature gradient $\vec{\nabla}T$. Since \vec{q} vanishes when $\vec{\nabla}T$ is zero, the lowest term in the expansion gives

$$\vec{q} = -\lambda\vec{\nabla}T . \quad (38)$$

The coefficient λ is called the thermal conductivity of the fluid. Since heat flows from high to low temperatures, the minus sign in the equation implies that $\lambda > 0$. We can now include the dissipative flux due to thermal conduction in the entropy- and energy-flux equations, Eqs. 34 and 37. We obtain

$$\rho T(\partial s/\partial t + \vec{v} \cdot \vec{\nabla}s) = \vec{\nabla} \cdot (\lambda\vec{\nabla}T) \quad (39)$$

and

$$(\partial/\partial t)((1/2)\rho v^2 + \rho\epsilon) = -\vec{\nabla} \cdot [\rho((1/2)v^2 + \epsilon + P/\rho)\vec{v} - \lambda\vec{\nabla}T] \quad (40)$$

respectively. In principle, viscous dissipative terms should be included in Eqs. 39 and 40 as well; but in practice for our applications the viscosity makes a major contribution only in the Navier-Stokes equation, Eq. 32, and we will neglect these terms in the energy or entropy equation.

It is often convenient to write the entropy-flux equation in terms of the temperature T . If the pressure variations in the fluid are so small that the corresponding changes in density can be neglected (which is the case when the velocity is much less than the velocity of sound), then $T\vec{\nabla}s = c_p\vec{\nabla}T$, and Eq. 39 yields

$$\rho c_p(\partial T/\partial t + \vec{v} \cdot \vec{\nabla}T) = \vec{\nabla} \cdot (\lambda\vec{\nabla}T) . \quad (41)$$

If we further assume that temperature variations are so small that λ may be regarded as constant, then

$$\partial T/\partial t + \vec{v} \cdot \vec{\nabla}T = \kappa\Delta T . \quad (42)$$

where

$$\kappa = \lambda/\rho c_p \quad (43)$$

is the thermal diffusivity. In real incompressible fluids with a nonuniform temperature, Eq. 42 together with the Navier-Stokes equation Eq. 32 and the continuity equation Eq. 16 provide an excellent approximation to the five relations which determine the fluid flow.

C. Boundary conditions.

In Sect. IV B we discussed the equations of motion of fluid flow. A particular fluid-flow problem is not defined, however, until boundary conditions (BC) have been specified. Usually there are solid walls which define the geometry of the fluid and which are impermeable. There may also be conditions imposed upon the temperature or upon the temperature gradient at the walls. For an ideal fluid, we generally have $v_n = 0$ at any stationary solid wall. Here v_n is the component of \vec{v} normal to the wall. This is so because there is no flow into or out of the wall. This BC by itself, without any constraint on the velocity components tangential to the wall, often is called a "free" BC because the fluid is free to move tangentially. In a real fluid, the components of \vec{v} tangential to a stationary wall also vanish because the viscosity assures that any fluid immediately adjacent to the wall is stationary. This BC often is referred to as a "rigid" BC. If the wall is moving, then for rigid BC's the fluid will move with it. Thus, for real fluids adjacent to a wall $\vec{v} = \vec{v}_w$, where \vec{v}_w is the wall velocity. For an incompressible fluid the condition that the velocity component tangential to the wall must vanish implies that the gradient of the velocity component perpendicular to the wall vanishes also. This is easy to see as follows. Let the z-component w of \vec{v} be normal to the wall, with the x and y components u and v tangential to the wall. Then $(\partial u/\partial x) = (\partial v/\partial y) = 0$ because u and v vanish for all x and y . In that case, by virtue of Eq. 16 $(\partial w/\partial z)$ must vanish also. That is, both the velocity and its gradient vanish at a rigid wall. Ideal-fluid (*i.e.* free) boundary conditions are often used because they greatly simplify the solution of many problems without losing the qualitative physics of the fluid flow. Although often the real fluid-flow problem differs from the ideal one only in quantitative aspects, there are cases where there is a qualitative difference.

D. Circular Couette flow.

As an example of a fluid-flow problem, let us consider an idealized version of the system described in Sect. II, and in particular by Fig. 4. It has a uniform temperature, *i.e.* $T_1 = 0$. Therefore Eq. 42 is trivially satisfied, and the last term on the rhs of Eq. 32 is zero. The system consists of two concentric cylinders rotating about their axis. We take this axis in the x-direction. Here we will take these cylinders to be infinitely long so as to avoid the

complications associated with the boundaries in the axial direction. The inner cylinder, of radius r_1 , rotates with angular speed ω_1 , and the outer one, of radius r_2 , with angular speed ω_2 . In most of the examples we will consider later on, $\omega_1 = \omega$ and $\omega_2 = 0$, *i.e.* the outer cylinder is stationary. It is convenient to make the variables dimensionless. Although there are several ways of doing this, a particularly good one is to measure length, time, and velocity in units of $d = r_2 - r_1$, d^2/ν , and $\omega_1 r$ respectively. Equation 32 then becomes

$$\partial \vec{v} / \partial t + R_i (\vec{v} \cdot \vec{\nabla}) \vec{v} = -\vec{\nabla} P + \Delta \vec{v} . \quad (44)$$

This equation contains only one independently variable parameter, namely the Reynolds number of the inner cylinder

$$R_i = \omega_1 r_1 d / \nu . \quad (45)$$

However, two other parameters are hidden away in the BC's. One of them is the outer cylinder speed, which can be written as a second Reynolds number

$$R_2 = \omega_2 r_2 d / \nu . \quad (46)$$

The third is the ratio

$$\eta = r_1 / r_2 \quad (47)$$

of the two cylinder radii. All Couette-Taylor systems with the same R_1 , R_2 , and η will behave in the same way, regardless of the detailed values of d , ν , ω_1 , and ω_2 . For the physical system there is of course one additional parameter, namely the axial length of the apparatus.

The boundary conditions are most easily imposed after re-writing Eq. 44 in cylindrical coordinates (r, ϕ, x) (which we will not do explicitly here). The symmetry of the boundaries suggests that there should be a time independent solution for which the components of $\vec{v}(r, \phi, x)$ are

$$v_x = v_r = 0 , \quad (48)$$

$$v_\phi = v_\phi(r) , \quad (49)$$

and

$$P = P(r) . \quad (50)$$

A more careful analysis confirms this, and one obtains the two ordinary differential equations

$$dP/dr = \rho v_\phi^2 / r \quad (51)$$

and

$$d^2v_\phi/dr^2 + (1/r)(dv_\phi/dr) - v_\phi/r^2 = 0 . \quad (52)$$

It is worth noting that the problem is so easy because the nonlinear term in the Navier-Stokes equation vanishes identically. This is so because, as implied by Eq. 26, \vec{v} has a non-vanishing component only in the ϕ -direction, whereas the gradient of \vec{v} is entirely in the r -direction. Thus the term $(\vec{v} \cdot \vec{\nabla})\vec{v}$ vanishes and we are left with a linear partial differential equation which trivially separates into ode's. This orthogonality between \vec{v} and its gradient is a general property of shear flows.

Equation 52 has a solution of the form

$$v_\phi = ar + b/r . \quad (53)$$

The constants a and b are determined by the boundary conditions. For rigid BC's we have $v_\phi = r_1\omega_1$ for $r = r_1$ and $v_\phi = r_2\omega_2$ at $r = r_2$. This yields

$$a = (\omega_2r_2^2 - \omega_1r_1^2)/(r_2^2 - r_1^2) \quad (54)$$

and

$$b = (\omega_1 - \omega_2)r_1^2r_2^2/(r_2^2 - r_1^2) . \quad (55)$$

The pressure can be found by integrating Eq. 51 after substituting the results for v_ϕ .

Although we have found a solution of the Navier-Stokes equation for the circular Couette boundary conditions, we have not yet established that this solution is stable. It turns out that Couette flow is stable for small ω_1 , but gives way to more complicated patterns such as Taylor-vortex flow for larger ω_1 .

E. Stability of fluid flow.

Let us assume that we have found a solution \vec{v}_0 to the equations of motion for a particular fluid-flow problem, such as the circular Couette state in the Couette-Taylor system given by Eqs. 7 to 9. We still have no assurance that these solutions are stable. By this we mean that there might be some disturbances which, if imposed upon the flow, would grow and would cause the system to choose another flow state. Here we only give a general outline of how to test for stability. To actually carry out the procedure usually involves numerical calculations [4] and its complexity is beyond the scope of our work.

We write the velocity, temperature, and pressure fields as

$$\vec{v} = \vec{v}_0 + \vec{v}_1, \tag{III.34a}$$

$$T = T_0 + T_1, \tag{III.34b}$$

and

$$P = P_0 + P_1. \tag{III.34c}$$

Here we regard $\vec{v}_1, T_1,$ and P_1 as a small perturbations of the solution $\vec{v}_0, T_0,$ and P_0 . Substituting into Eqs. 8, 23, and 14a, 14b, or 14c, the terms involving only $\vec{v}_0, T_0,$ and P_0 cancel because these functions satisfy the equations. There remains an equation for $\vec{v}_1, T_1,$ and P_1 which may contain the known functions $\vec{v}_0, T_0,$ and P_0 as coefficients. Since $\vec{v}_1, T_1,$ and P_1 are considered small perturbations, we can omit nonlinear terms such as $\vec{v}_1 \cdot \vec{\nabla} \vec{v}_1,$ and obtain, for instance from the Navier-Stokes and continuity equations,

$$\partial \vec{v}_1 / \partial t + (\vec{v}_0 \cdot \vec{\nabla}) \vec{v}_1 + (\vec{v}_1 \cdot \vec{\nabla}) \vec{v}_0 = -\vec{\nabla} P_1 / \rho + \nu \Delta \vec{v}_1 \tag{III.35}$$

and

$$\vec{\nabla} \cdot \vec{v}_1 = 0. \tag{III.36}$$

In the RB case, the disturbance of T must also be substituted in Eq. 31c. If we write the disturbance \vec{v}_1 as

$$\vec{v}_1 = A(t) \vec{f}(x, y, z), \tag{III.37}$$

and take A to have the form

$$A(t) = \text{constant} \times \exp(\gamma t) \exp(-i\omega t) \tag{III.38}$$

then we can determine whether the growth rate γ is positive or negative. The original solution \vec{v}_0 is stable only if the growth rates are negative for all $\vec{f}(x, y, z)$. Of course to determine the nature of the solution corresponding to a growing \vec{v}_1 requires the solution of nonlinear equations; and generally this is a difficult problem.

[1] A sizable literature now exists dealing with this system. A comprehensive review has been given by R.C. DiPrima and H.L. Swinney, in *Hydrodynamic Instabilities and Transitions to*

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