Advanced Classical Mechanics

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Classical Mechanics III MIT 8.09 & 8.309

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Contents

1	A F	Review of Analytical Mechanics	1		
	1.1	Introduction	1		
	1.2	Lagrangian & Hamiltonian Mechanics	1		
	1.3	Symmetry and Conservation Laws	12		
	1.4	Constraints and Friction Forces	16		
	1.5	Calculus of Variations & Lagrange Multipliers	22		
2	Rig	id Body Dynamics	30		
	2.1	Coordinates of a Rigid Body	30		
	2.2	Time Evolution with Rotating Coordinates	34		
	2.3	Kinetic Energy, Angular Momentum, and			
		the Moment of Inertia Tensor for Rigid Bodies	38		
	2.4	Euler Equations	45		
	2.5	Symmetric Top with One Point Fixed	47		
3	Vibrations & Oscillations				
	3.1	Simultaneous Diagonalization of \hat{T} and \hat{V}	54		
	3.2	Vibrations and Oscillations with Normal Coordinates	56		
4	Canonical Transformations, Hamilton-Jacobi Equations, and Action-Angle				
	Var	iables	61		
	4.1	Generating Functions for Canonical Transformations	61		
	4.2	Poisson Brackets and the Symplectic Condition	67		
	4.3	Equations of Motion & Conservation Theorems	70		
	4.4	Hamilton-Jacobi Equation	74		
	4.5	Kepler Problem	80		
	4.6	Action-Angle Variables	84		
5	Per	Perturbation Theory			
	5.1	Time Dependent Perturbation Theory for the			
		Hamilton-Jacobi Equations	94		
	5.2	Periodic and Secular Perturbations to Finite Angle Pendulum	96		

CONTENTS

	5.3	Perihelion Precession from Perturbing a Kepler Orbit	99
6	Flui	id Mechanics 1	05
	6.1		05
	6.2	9	110
		•	110
		v ±	13
		1 10	114
	6.3		116
	6.4	· ·	122
	6.5		126
	6.6		129
	6.7	1	132
	6.8		136
	0.0		.00
7	Cha	U	43
	7.1	Introduction to Chaos	44
			44
		7.1.2 Evolution of Phase Space	146
		7.1.3 Fixed Points	147
			150
	7.2	Bifurcations	154
		7.2.1 Saddle-Node Bifurcation	155
		7.2.2 Transcritical Bifurcation	158
		7.2.3 Supercritical Pitchfork Bifurcation	160
		7.2.4 Subcritical pitchfork bifurcation	161
	7.3	Fixed Points in Two-Dimensional Systems	165
			165
		7.3.2 Systems with a conserved $E(\vec{x})$	172
	7.4	Limit Cycles and Bifurcations	174
		7.4.1 Poincaré-Bendixson Theorem	176
		7.4.2 Fixed Point Bifurcations Revisited and Hopf Bifurcations	177
	7.5		183
	7.6	Chaos in Differential Equations, Strange Attractors, and Fractals	190
			190
		-	92
		v I	195

Chapter 1

A Review of Analytical Mechanics

1.1 Introduction

These lecture notes cover the third course in Classical Mechanics, taught at MIT by Professor Stewart to advanced undergraduates (course 8.09) as well as to graduate students (course 8.309). In the prerequisite classical mechanics II course the students are taught both Lagrangian and Hamiltonian dynamics, including Kepler bound motion and central force scattering, and the basic ideas of canonical transformations. This course briefly reviews the needed concepts, but assumes some familiarity with these ideas. References used for this course include

- Goldstein, Poole & Safko, Classical Mechanics, 3rd edition.
- Landau and Lifshitz vol.6, *Fluid Mechanics*. Symon, *Mechanics* for reading material on non-viscous fluids.
- Strogatz, Nonlinear Dynamics and Chaos.
- Review: Landau & Lifshitz vol.1, *Mechanics*. (Typically used for the prerequisite Classical Mechanics II course and hence useful here for review)

1.2 Lagrangian & Hamiltonian Mechanics

Newtonian Mechanics

In Newtonian mechanics, the dynamics of a system of N particles are determined by solving for their coordinate trajectories as a function of time. This can be done through the usual vector spatial coordinates $\mathbf{r}_i(t)$ for $i \in \{1, ..., N\}$, or with generalized coordinates $q_i(t)$ for $i \in \{1, ..., 3N\}$ in 3-dimensional space. These generalized coordinates could be angles, distances between objects, etc.

Velocities will be represented through $\mathbf{v}_i \equiv \dot{\mathbf{r}}_i$ for spatial coordinates, or through \dot{q}_i for generalized coordinates. Note that dots above a symbol will always denote the total time derivative $\frac{d}{dt}$. Momenta are likewise either Newtonian momenta $\mathbf{p}_i = m_i \mathbf{v}_i$ or generalized momenta p_i .

For a fixed set of masses m_i Newton's 2nd law can be expressed in 2 equivalent ways:

- 1. It can be expressed as N second-order equations $\mathbf{F}_i = \frac{d}{dt} (m_i \dot{\mathbf{r}}_i)$ with 2N boundary conditions given in $\mathbf{r}_i(0)$ and $\dot{\mathbf{r}}_i(0)$. The problem then becomes one of determining the N vector variables $\mathbf{r}_i(t)$.
- 2. It can also be expressed as an equivalent set of 2N 1st order equations $\mathbf{F}_i = \dot{\mathbf{p}}_i$ & $\mathbf{p}_i/m_i = \dot{\mathbf{r}}_i$ with 2N boundary conditions given in $\mathbf{r}_i(0)$ and $\mathbf{p}_i(0)$. The problem then becomes one of determining the 2N vector variables $\mathbf{r}_i(t)$ and $\mathbf{p}_i(t)$.

Note that $\mathbf{F} = m\mathbf{a}$ holds in *inertial frames*. Inertial frames describe time and space homogeneously (invariant to displacements), isotropically (invariant to rotations), and in a time independent manner. These are frames where the motion of a particle not subject to any force is in a straight line with constant velocity. (The converse, that straight line motion with constant velocity implies there are no forces present, is not true. For example, there may be forces that are not visible for motion in a particular direction, like motion parallel to a constant magnetic field.) Noninertial frames also generically have "fictitious forces", such as the centrifugal and Coriolis effects. Inertial frames also play a key role in special relativity. In general relativity the concept of inertial frames is replaced by that of geodesic motion.

The existence of an inertial frame is a useful approximation for working out the dynamics of particles, and non-inertial terms can often be included as perturbative corrections. Examples of approximate inertial frames are that of a fixed Earth, or better yet, of fixed distant stars. We can still test for how noninertial we are by looking for fictitious forces that (a) may point back to an origin with no source for the force or (b) behave in a non-standard fashion in different frames (i.e. they transform in an unusual manner when going between different frames).

We will use primes to denote coordinate transformations. If \mathbf{r} is measured in an inertial frame S, and \mathbf{r}' is measured in frame S' whose relation to S is given by a transformation $\mathbf{r}' = \mathbf{f}(\mathbf{r},t)$, then S' is inertial iff $\ddot{\mathbf{r}} = 0 \leftrightarrow \ddot{\mathbf{r}}' = 0$. This is solved by the Galilean transformations,

$$\mathbf{r}' = \mathbf{r} + \mathbf{v}_0 t,$$

$$t' = t,$$

with constant \mathbf{v}_0 . This transformation preserves the inertiality of frames, with $\mathbf{F} = m\ddot{\mathbf{r}}$ equivalent to $\mathbf{F}' = m\ddot{\mathbf{r}}'$. Galilean transformations are the non-relativistic limit, $v \ll c$, of Lorentz transformations which preserve inertial frames in special relativity. A few examples related to the concepts of inertial frames are:

1. In a rotating frame, the transformation is given by

$$\begin{bmatrix} x' \\ y' \end{bmatrix} = \begin{bmatrix} \cos(\theta) & \sin(\theta) \\ -\sin(\theta) & \cos(\theta) \end{bmatrix} \begin{bmatrix} x \\ y \end{bmatrix}$$

If $\theta = \omega t$ for some constant ω , then $\ddot{\mathbf{r}} = 0$ does not imply $\{\ddot{x}', \ddot{y}'\} = 0$, so the primed frame is noninertial.

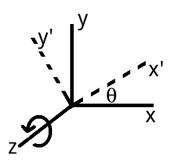


Figure 1.1: Frame rotated by an angle θ

2. In polar coordinates, $\mathbf{r} = r\hat{r}$, and the unit vectors are time dependent,

$$\frac{d\hat{r}}{dt} = \dot{\theta}\hat{\theta}, \qquad \frac{d\hat{\theta}}{dt} = -\dot{\theta}\hat{r} \tag{1.1}$$

and thus

$$\ddot{\mathbf{r}} = \ddot{r}\hat{r} + 2\dot{r}\dot{\theta}\hat{\theta} + r\left(\ddot{\theta}\hat{\theta} - \dot{\theta}^2\hat{r}\right). \tag{1.2}$$

Even if $\ddot{\mathbf{r}} = 0$ we can still have $\ddot{r} \neq 0$ and $\ddot{\theta} \neq 0$, and we can not in general form a simple Newtonian force law equation $m\ddot{q} = F_q$ for each of these coordinates. This is different than the first example, since here we are picking coordinates rather than changing the reference frame, so to remind ourselves about their behavior we will call these "non-inertial coordinates" (which we may for example decide to use in an inertial frame). In general, curvilinear coordinates are non-inertial.

Lagrangian Mechanics

In Lagrangian mechanics, the key function is the Lagrangian

$$L = L(q, \dot{q}, t). \tag{1.3}$$

Here, $q = (q_1, \ldots, q_N)$ and likewise $\dot{q} = (\dot{q}_1, \ldots, \dot{q}_N)$. We are now letting N denote the number of scalar (rather than vector) variables, and will often use the short form to denote

dependence on these variables, as in Eq. (1.3). Typically we can write L = T - V where T is the kinetic energy and V is the potential energy. In the simplest cases, $T = T(\dot{q})$ and V = V(q), but we also allow the more general possibility that $T = T(q, \dot{q}, t)$ and $V = V(q, \dot{q}, t)$. It turns out, as we will discuss later, that even this generalization does not describe all possible classical mechanics problems.

The solution to a given mechanical problem is obtained by solving a set of N second-order differential equations known as Euler-Lagrange equations of motion,

$$\frac{d}{dt} \left(\frac{\partial L}{\partial \dot{q}_i} \right) - \frac{\partial L}{\partial q_i} = 0. \tag{1.4}$$

These equations involve \ddot{q}_i , and reproduce the Newtonian equations $\mathbf{F} = m\mathbf{a}$. The principle of stationary action (Hamilton's principle),

$$\delta S = \delta \int_{t_1}^{t_2} L(q, \dot{q}, t) dt = 0, \tag{1.5}$$

is the starting point for deriving the Euler-Lagrange equations. Although you have covered the Calculus of Variations in an earlier course on Classical Mechanics, we will review the main ideas in Section 1.5.

There are several advantages to working with the Lagrangian formulation, including

- 1. It is easier to work with the scalars T and V rather than vectors like \mathbf{F} .
- 2. The same formula in equation (1.4) holds true regardless of the choice of coordinates. To demonstrate this, let us consider new coordinates

$$Q_i = Q_i(q_1, \dots, q_N, t). \tag{1.6}$$

This particular sort of transformation is called a *point transformation*. Defining the new Lagrangian by

$$L' = L'(Q, \dot{Q}, t) = L(q, \dot{q}, t),$$
 (1.7)

we claim that the equations of motion are simply

$$\frac{d}{dt} \left(\frac{\partial L'}{\partial \dot{Q}_i} \right) - \frac{\partial L'}{\partial Q_i} = 0. \tag{1.8}$$

Proof: (for N = 1, since the generalization is straightforward) Given $L'(Q, \dot{Q}, t) = L(q, \dot{q}, t)$ with Q = Q(q, t) then

$$\dot{Q} = \frac{d}{dt}Q(q,t) = \frac{\partial Q}{\partial q}\dot{q} + \frac{\partial Q}{\partial t}.$$
 (1.9)

Therefore

$$\frac{\partial \dot{Q}}{\partial \dot{q}} = \frac{\partial Q}{\partial q},\tag{1.10}$$

a result that we will use again in the future. Then

$$\frac{\partial L}{\partial q} = \frac{\partial L'}{\partial q} = \frac{\partial L'}{\partial Q} \frac{\partial Q}{\partial q} + \frac{\partial L'}{\partial \dot{Q}} \frac{\partial \dot{Q}}{\partial q},$$

$$\frac{\partial L}{\partial \dot{q}} = \frac{\partial L'}{\partial \dot{q}} = \frac{\partial L'}{\partial \dot{Q}} \frac{\partial \dot{Q}}{\partial \dot{q}} = \frac{\partial L'}{\partial \dot{Q}} \frac{\partial Q}{\partial q}.$$
(1.11)

Since $\frac{\partial Q}{\partial \dot{q}} = 0$ there is no term $\frac{\partial L'}{\partial Q} \frac{\partial Q}{\partial \dot{q}}$ in the last line.

Plugging these results into $0 = \frac{d}{dt} \left(\frac{\partial L}{\partial \dot{q}} \right) - \frac{\partial L}{\partial q}$ gives

$$0 = \left[\frac{d}{dt} \left(\frac{\partial L'}{\partial \dot{Q}} \right) \frac{\partial Q}{\partial q} + \frac{\partial L'}{\partial \dot{Q}} \frac{d}{dt} \left(\frac{\partial Q}{\partial q} \right) \right] - \left[\frac{\partial L'}{\partial Q} \frac{\partial Q}{\partial q} + \frac{\partial L'}{\partial \dot{Q}} \frac{\partial \dot{Q}}{\partial q} \right]$$
$$= \left[\frac{d}{dt} \left(\frac{\partial L'}{\partial \dot{Q}} \right) - \frac{\partial L'}{\partial Q} \right] \frac{\partial Q}{\partial q} . \tag{1.12}$$

Here we have used $\frac{d}{dt} \frac{\partial Q}{\partial q} = (\dot{q} \frac{\partial}{\partial q} + \frac{\partial}{\partial t}) \frac{\partial Q}{\partial q} = \frac{\partial}{\partial q} (\dot{q} \frac{\partial}{\partial q} + \frac{\partial}{\partial t}) Q = \frac{\partial \dot{Q}}{\partial q}$ so that the second and fourth terms in the first line of Eq. (1.12) cancel. Finally for non-trivial transformation where $\frac{\partial Q}{\partial q} \neq 0$ we have, as expected,

$$0 = \frac{d}{dt} \left(\frac{\partial L'}{\partial \dot{Q}} \right) - \frac{\partial L'}{\partial Q} \,. \tag{1.13}$$

Note two things:

- This implies we can freely use the Euler-Lagrange equations for noninertial coordinates.
- We can formulate L in whatever coordinates are easiest, and then change to convenient variables that better describe the symmetry of a system (for example, Cartesian to spherical).
- 3. Continuing our list of advantages for using L, we note that it is also easy to incorporate constraints. Examples include a mass constrained to a surface or a disk rolling without slipping. Often when using L we can avoid discussing forces of constraint (for example, the force normal to the surface).

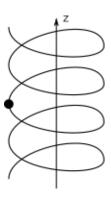


Figure 1.2: Bead on a helix

Lets discuss the last point in more detail (we will also continue to discuss it in the next section). The method for many problems with constraints is to simply make a good choice for the generalized coordinates to use for the Lagrangian, picking N-k independent variables q_i for a system with k constraints.

Example: For a bead on a helix as in Fig. 1.2 we only need one variable, $q_1 = z$.

Example: A mass m_2 attached by a massless pendulum to a horizontally sliding mass m_1 as in Fig. 1.3, can be described with two variables $q_1 = x$ and $q_2 = \theta$.

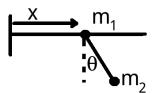


Figure 1.3: Pendulum of mass m_2 hanging on a rigid bar of length ℓ whose support m_1 is a frictionless horizontally sliding bead

Example: As an example using non-inertial coordinates consider a potential $V = V(r, \theta)$ in polar coordinates for a fixed mass m at position $\mathbf{r} = r\hat{r}$. Since $\dot{\mathbf{r}} = \dot{r}\hat{r} + r\dot{\theta}\hat{\theta}$ we have $T = \frac{m}{2}\dot{\mathbf{r}}^2 = \frac{m}{2}(\dot{r}^2 + r^2\dot{\theta}^2)$, giving

$$L = \frac{m}{2} \left(\dot{r}^2 + r^2 \dot{\theta}^2 \right) - V(r, \theta). \tag{1.14}$$

For r the Euler-Lagrange equation is

$$0 = \frac{d}{dt} \left(\frac{\partial L}{\partial \dot{r}} \right) - \frac{\partial L}{\partial r} = \frac{d}{dt} \left(m\dot{r} \right) - mr\dot{\theta}^2 + \frac{\partial V}{\partial r}. \tag{1.15}$$

This gives

$$m\ddot{r} - mr\dot{\theta}^2 = -\frac{\partial V}{\partial r} = F_r,\tag{1.16}$$

from which we see that $F_r \neq m\ddot{r}$. For θ the Euler-Lagrange equation is

$$0 = \frac{d}{dt} \left(\frac{\partial L}{\partial \dot{\theta}} \right) - \frac{\partial L}{\partial \theta} = \frac{d}{dt} \left(mr^2 \dot{\theta} \right) + \frac{\partial V}{\partial \theta}. \tag{1.17}$$

This gives

$$\frac{d}{dt}\left(mr^2\dot{\theta}\right) = -\frac{\partial V}{\partial \theta} = F_{\theta},\tag{1.18}$$

which is equivalent to the relation between angular momentum and torque perpendicular to the plane, $\dot{L}_z = F_\theta = \tau_z$. (Recall $\mathbf{L} = \mathbf{r} \times \mathbf{p}$ and $\tau = \mathbf{r} \times \mathbf{F}$.)

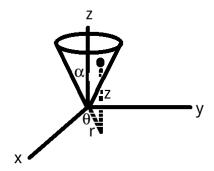


Figure 1.4: Particle on the inside of a cone

Example: Let us consider a particle rolling due to gravity in a frictionless cone, shown in Fig. 1.4, whose opening angle α defines an equation for points on the cone $\tan(\alpha) = \sqrt{x^2 + y^2}/z$. There are 4 steps which we can take to solve this problem (which are more general than this example):

- 1. Formulate T and V by N=3 generalized coordinates. Here it is most convenient to choose cylindrical coordinates denoted (r,θ,z) , so that $T=\frac{m}{2}\left(\dot{r}^2+r^2\dot{\theta}^2+\dot{z}^2\right)$ and V=mgz.
- 2. Reduce the problem to N-k=2 independent coordinates and determine the new Lagrangian L=T-V. In this case we eliminate $z=r\cot(\alpha)$ and $\dot{z}=\dot{r}\cot(\alpha)$, so

$$L = \frac{m}{2} \left[\left(1 + \cot^2 \alpha \right) \dot{r}^2 + r^2 \dot{\theta}^2 \right] - mgr \cot \alpha. \tag{1.19}$$

3. Find the Euler-Lagrange equations. For r, $0 = \frac{d}{dt} \left(\frac{\partial L}{\partial \dot{r}} \right) - \frac{\partial L}{\partial r}$, which here is

$$0 = \frac{d}{dt} \left[m \left(1 + \cot^2 \alpha \right) \dot{r} \right] - mr \dot{\theta}^2 + mg \cot \alpha$$
 (1.20)

giving

$$(1 + \cot^2 \alpha) \ddot{r} - r\dot{\theta}^2 + g \cot \alpha = 0. \tag{1.21}$$

For θ we have $0 = \frac{d}{dt} \left(\frac{\partial L}{\partial \dot{\theta}} \right) - \frac{\partial L}{\partial \theta}$, so

$$0 = \frac{d}{dt} \left(mr^2 \dot{\theta} \right) - 0, \tag{1.22}$$

giving

$$(2\dot{r}\dot{\theta} + r\ddot{\theta})r = 0. \tag{1.23}$$

4. Solve the system analytically or numerically, for example using Mathematica. Or we might be only interested in determining certain properties or characteristics of the motion without a full solution (we will discuss this last point in detail in Ch.7).

Hamiltonian Mechanics

In Hamiltonian mechanics, the canonical momenta $p_i \equiv \frac{\partial L}{\partial \dot{q}_i}$ are promoted to coordinates on equal footing with the generalized coordinates q_i . The coordinates (q, p) are canonical variables, and the space of canonical variables is known as phase space.

The Euler-Lagrange equations say $\dot{p}_i = \frac{\partial L}{\partial q_i}$. These need not equal the kinematic momenta $m_i \dot{q}_i$ if $V = V(q, \dot{q})$. Performing the Legendre transformation

$$H(q, p, t) = \dot{q}_i \, p_i - L(q, \dot{q}, t) \tag{1.24}$$

(where for this equation, and henceforth, repeated indices will imply a sum unless otherwise specified) yields the *Hamilton equations of motion*

$$\dot{q}_i = \frac{\partial H}{\partial p_i}$$

$$\dot{p}_i = -\frac{\partial H}{\partial q_i}$$
(1.25)

which are 2N 1st order equations. We also have the result that

$$\frac{\partial H}{\partial t} = -\frac{\partial L}{\partial t} \,. \tag{1.26}$$

Proof: (for N = 1) Consider how the left and right hand side of the Legendre transformation behave under differential changes to the variables,

$$dH = \frac{\partial H}{\partial q} dq + \frac{\partial H}{\partial p} dp + \frac{\partial H}{\partial t} dt$$

$$= p d\dot{q} + \dot{q} dp - \frac{\partial L}{\partial q} dq - \frac{\partial L}{\partial \dot{q}} d\dot{q} - \frac{\partial L}{\partial t} dt = \dot{q} dp - \dot{p} dq - \frac{\partial L}{\partial t} dt .$$
(1.27)

In the last line we used the definition $p = \frac{\partial L}{\partial \dot{q}}$ and the E-L equation, $\dot{p} = \frac{\partial L}{\partial q}$. Since we are free to independently vary dq, dp, and dt this implies (generalizing now to N variables):

$$-\frac{\partial H}{\partial q_i} = \dot{p}_i \,, \qquad \frac{\partial H}{\partial p_i} = \dot{q}_i \,, \qquad \frac{\partial H}{\partial t} = -\frac{\partial L}{\partial t} \,. \tag{1.28}$$

The first two results are Hamilton's equations. We can interpret these two Hamilton equations as follows:

- $\dot{q}_i = \frac{\partial H(q,p,t)}{\partial p_i} = \dot{q}_i(q,p,t)$ is an inversion of $p_i = \frac{\partial L(q,\dot{q},t)}{\partial \dot{q}_i} = p_i(q,\dot{q},t)$.
- $\dot{p}_i = -\frac{\partial H}{\partial q_i}$ provides the Newtonian dynamics.

However, these two equation have an have equal footing in Hamiltonian mechanics, since the coordinates and momenta are treated on a common ground. We can use $p_i = \frac{\partial L}{\partial \dot{q}_i}$ to construct H from L and then forget about L.

As an example of the manner in which we will usually consider transformations between Lagrangians and Hamiltonians, consider again the variables relevant for the particle on a cone from Fig. 1.4:

$$L(r,\theta,z,\dot{r},\dot{\theta},\dot{z}) \xrightarrow{z=r\cot\alpha} \text{new } L(r,\theta,\dot{r},\dot{\theta}) \longrightarrow \text{Euler-Lagrange Eqtns.} \quad (1.29)$$

$$\updownarrow \qquad \qquad \qquad \qquad \qquad \qquad \qquad \qquad \downarrow$$

$$H(r,\theta,z,p_r,p_\theta,p_z) \stackrel{\text{not here}}{\Longrightarrow} \qquad H(r,\theta,p_r,p_\theta) \longrightarrow \text{Hamilton Eqtns.}$$

Here we consider transforming between L and H either before or after removing the redundant coordinate z, but in this course we will only consider constraints imposed on Lagrangians and not in the Hamiltonian formalism (the step indicated by \Longrightarrow). For the curious, the topic of imposing constraints on Hamiltonians, including even more general constraints than those we will consider, is covered well in Dirac's little book "Lectures on Quantum Mechanics". Although Hamiltonian and Lagrangian mechanics provide equivalent formalisms, there is often an advantage to using one or the other. In the case of Hamiltonian mechanics potential advantages include the language of phase space with Liouville's Theorem, Poisson Brackets and the connection to quantum mechanics, as well as the Hamilton-Jacobi transformation theory (all to be covered later on).

Special case: Let us consider a special case that is sufficient to imply that the Hamiltonian is equal to the energy, $H = E \equiv T + V$. If we only have quadratic dependence on velocities in the kinetic energy, $T = \frac{1}{2}T_{jk}(q)\dot{q}_j\dot{q}_k$, and V = V(q) with L = T - V, then

$$\dot{q}_i p_i = \dot{q}_i \frac{\partial L}{\partial \dot{q}_i} = \frac{1}{2} \dot{q}_i T i k \dot{q}_k + \frac{1}{2} \dot{q}_j T_{ji} \dot{q}_i = 2T. \tag{1.30}$$

Hence,

$$H = \dot{q}_i p_i - L = T + V = E \tag{1.31}$$

which is just the energy.

Another Special case: Consider a class of Lagrangians given as

$$L(q, \dot{q}, t) = L_0 + a_j \dot{q}_j + \frac{1}{2} \dot{q}_j T_{jk} \dot{q}_k$$
(1.32)

where $L_0 = L_0(q, t)$, $a_j = a_j(q, t)$, and $T_{jk} = T_{kj} = T_{jk}(q, t)$. We can write this in shorthand as

$$L = L_0 + \vec{a} \cdot \dot{\vec{q}} + \frac{1}{2} \dot{\vec{q}} \cdot \hat{T} \cdot \dot{\vec{q}}. \tag{1.33}$$

Here the generalized coordinates, momenta, and coefficients have been collapsed into vectors, like \vec{q} (rather than the boldface that we reserve for Cartesian vectors), and dot products of vectors from the left imply transposition of that vector. Note that \vec{q} is an unusual vector, since its components can have different dimensions, eg. $\vec{q} = (x, \theta)$, but nevertheless this notation is useful. To find H,

$$p_j = \frac{\partial L}{\partial \dot{q}_j} = a_j + T_{jk} \dot{q}_k, \tag{1.34}$$

meaning $\vec{p} = \vec{a} + \hat{T} \cdot \dot{\vec{q}}$. Inverting this gives $\dot{\vec{q}} = \hat{T}^{-1}$. $(\vec{p} - \vec{a})$, where \hat{T}^{-1} will exist because of the positive-definite nature of kinetic energy, which implies that \hat{T} is a postive definite matrix. Thus, $H = \dot{\vec{q}} \cdot \vec{p} - L$ yields

$$H = \frac{1}{2}(\vec{p} - \vec{a}) \cdot \hat{T}^{-1} \cdot (\vec{p} - \vec{a}) - L_0(q, t)$$
(1.35)

as the Hamiltonian. So for any Lagrangian in the form of Eq. (1.32), we can find \hat{T}^{-1} and write down the Hamiltonian as in Eq. (1.35) immediately.

Example: let us consider $L = \frac{1}{2}m\mathbf{v}^2 - e\phi + e\mathbf{A} \cdot \mathbf{v}$, where e is the electric charge and SI units are used. In Eq. (1.32), because the coordinates are Cartesian, $\mathbf{a} = e\mathbf{A}$, $\hat{T} = m\mathbb{1}$, and $L_0 = -e\phi$, so

$$H = \frac{1}{2m} \left(\mathbf{p} - e\mathbf{A} \right)^2 + e\phi. \tag{1.36}$$

As you have presumably seen in an earlier course, this Hamiltonian does indeed reproduce the Lorentz force equation $e(\mathbf{E} + \mathbf{v} \times \mathbf{B}) = m\dot{\mathbf{v}}$.

A more detailed Example. Find L and H for the frictionless pendulum shown in Fig. 1.3. This system has two constraints, that m_1 is restricted to lie on the x-axis sliding without friction, and that the rod between m_1 and m_2 is rigid, giving

$$y_1 = 0,$$
 $(y_1 - y_2)^2 + (x_1 - x_2)^2 = \ell^2.$ (1.37)

Prior to imposing any constraints the Lagrangian is

$$L = T - V = \frac{m_1}{2}\dot{x}_1^2 + \frac{m_2}{2}(\dot{x}_2^2 + \dot{y}_2^2) - m_2gy_2 - m_1gy_1.$$
 (1.38)

Lets choose to use $x \equiv x_1$ and the angle θ as the independent coordinates after imposing the constraints in Eq. (1.37). This allows us to eliminate $y_1 = 0$, $x_2 = x + \ell \sin \theta$ and $y_2 = -\ell \cos \theta$, together with $\dot{x}_2 = \dot{x} + \ell \cos \theta \dot{\theta}$, $\dot{y}_2 = \ell \sin \theta \dot{\theta}$, $\dot{x}_1 = \dot{x}$. The Lagrangian with constraints imposed is

$$L = \frac{m_1}{2}\dot{x}^2 + \frac{m_2}{2}(\dot{x}^2 + 2\ell\cos\theta\,\dot{x}\dot{\theta} + \ell^2\cos^2\theta\,\dot{\theta}^2 + \ell^2\sin^2\theta\,\dot{\theta}^2) + m_2g\ell\cos\theta\,. \tag{1.39}$$

Next we determine the Hamiltonian. First we find

$$p_{x} = \frac{\partial L}{\partial \dot{x}} = m_{1}\dot{x} + m_{2}(\dot{x} + \ell\cos\theta\dot{\theta}) = (m_{1} + m_{2})\dot{x} + m_{2}\ell\cos\theta\dot{\theta}, \qquad (1.40)$$

$$p_{\theta} = \frac{\partial L}{\partial \dot{\theta}} = m_{2}\ell\cos\theta\dot{x} + m_{2}\ell^{2}\dot{\theta}.$$

Note that p_x is not simply proportional to \dot{x} here (actually p_x is the center-of-mass momentum). Writing $\begin{pmatrix} p_x \\ p_\theta \end{pmatrix} = \hat{T} \cdot \begin{pmatrix} \dot{x} \\ \dot{\theta} \end{pmatrix}$ gives

$$\hat{T} = \begin{pmatrix} m_1 + m_2 & m_2 \ell \cos \theta \\ m_2 \ell \cos \theta & m_2 \ell^2 \end{pmatrix}, \tag{1.41}$$

with $L = \frac{1}{2}(\dot{x}\,\dot{\theta}) \cdot \hat{T} \cdot \begin{pmatrix} \dot{x} \\ \dot{\theta} \end{pmatrix} + L_0$ where $L_0 = m_2 g \ell \cos \theta$. Computing

$$\hat{T}^{-1} = \frac{1}{m_1 m_2 \ell^2 + m_2 \ell^2 \sin^2 \theta} \begin{pmatrix} m_2 \ell^2 & -m_2 \ell \cos \theta \\ -m_2 \ell \cos \theta & m_1 + m_2 \end{pmatrix}, \tag{1.42}$$

we can simply apply Eq. (1.35) to find the corresponding Hamiltonian

$$H = \frac{1}{2} (p_x p_\theta) \cdot \hat{T}^{-1} \cdot \begin{pmatrix} p_x \\ p_\theta \end{pmatrix} - m_2 g \ell \cos \theta$$

$$= \frac{1}{2m_2 \ell^2 (m_1 + m_2 \sin^2 \theta)} \left[m_2 \ell^2 p_x^2 + (m_1 + m_2) p_\theta^2 - 2m_2 \ell \cos \theta p_x p_\theta \right] - m_2 g \ell \cos \theta .$$
(1.43)

Lets compute the Hamilton equations of motion for this system. First for (x, p_x) we find

$$\dot{x} = \frac{\partial H}{\partial p_x} = \frac{p_x}{m_1 + m_2 \sin^2 \theta} - \frac{\cos \theta \, p_\theta}{\ell (m_1 + m_2 \sin^2 \theta)}, \qquad (1.44)$$

$$\dot{p}_x = -\frac{\partial H}{\partial x} = 0.$$

As we might expect, the CM momentum is time independent. Next for (θ, p_{θ}) :

$$\dot{\theta} = \frac{\partial H}{\partial p_{\theta}} = \frac{1}{m_{2}\ell^{2}(m_{1} + m_{2}\sin^{2}\theta)} \Big[(m_{1} + m_{2})p_{\theta} - m_{2}\ell\cos\theta p_{x} \Big] ,$$

$$\dot{p}_{\theta} = -\frac{\partial H}{\partial \theta} = \frac{\sin\theta\cos\theta}{\ell^{2}(m_{1} + m_{2}\sin^{2}\theta)^{2}} \Big[m_{2}\ell^{2}p_{x}^{2} + (m_{1} + m_{2})p_{\theta}^{2} - 2m_{2}\ell\cos\theta p_{x}p_{\theta} \Big]$$

$$- m_{2}g\ell\sin\theta - \frac{\sin\theta p_{x}p_{\theta}}{\ell(m_{1} + m_{2}\sin\theta)} .$$
(1.45)

These non-linear coupled equations are quite complicated, but could be solved in mathematica or another numerical package. To test our results for these equations of motion analytically, we can take the small angle limit, approximating $\sin \theta \simeq \theta$, $\cos \theta \simeq 1$ to obtain

$$\dot{x} = \frac{p_x}{m_1} - \frac{p_\theta}{\ell m_1}, \qquad \dot{p}_x = 0, \qquad \dot{\theta} = \frac{1}{m_1 m_2 \ell^2} \Big[(m_1 + m_2) p_\theta - m_2 \ell p_x \Big],
\dot{p}_\theta = \frac{\theta}{\ell^2 m_1^2} \Big[m_2 \ell^2 p_x^2 + (m_1 + m_2) p_\theta^2 - 2m_2 \ell \cos \theta p_x p_\theta \Big] - \frac{\theta p_x p_\theta}{\ell m_1} - m_2 g \ell \theta.$$
(1.46)

To simplify it further we can work in the CM frame, thus setting $p_x = 0$, and linearize the equations by noting that $p_{\theta} \sim \dot{\theta}$ should be small for θ to remain small, and hence θp_{θ}^2 is a higher order term. For the non-trivial equations this leaves

$$\dot{x} = -\frac{p_{\theta}}{\ell m_1}, \qquad \dot{\theta} = \frac{p_{\theta}}{\mu \ell^2}, \qquad \dot{p}_{\theta} = -m_2 g \ell \theta, \qquad (1.47)$$

where $\mu = m_1 m_2/(m_1 + m_2)$ is the reduced mass for the two-body system. Thus $\ddot{\theta} = \dot{p}_{\theta}/(\mu \ell^2) = -\frac{m_2 g}{\mu \ell} \theta$ as expected for simple harmonic motion.

1.3 Symmetry and Conservation Laws

A cyclic coordinate is one which does not appear in the Lagrangian, or equivalently in the Hamiltonian. Because $H(q, p, t) = \dot{q}_i p_i - L(q, \dot{q}, t)$, if q_j is absent in L for some particular j, it will be absent in H as well. The absence of that q_j corresponds with a symmetry in the dynamics.

In this context, Noether's theorem means that a symmetry implies a cyclic coordinate, which in turn produces a conservation law. If q_j is a cyclic coordinate for some j, then we can change that coordinate without changing the dynamics given by the Lagrangian or Hamiltonian, and hence there is a symmetry. Furthermore the corresponding canonical momentum p_j is conserved, meaning it is a constant through time.

The proof is simple. If $\frac{\partial L}{\partial q_j} = 0$ then $\dot{p}_j = \frac{d}{dt} \frac{\partial L}{\partial \dot{q}_j} = \frac{\partial L}{\partial q_j} = 0$, or even more simply, $\frac{\partial H}{\partial q_j} = 0$ is equivalent to $\dot{p}_j = 0$, so p_j is a constant in time.

Special cases and examples of this abound. Lets consider a few important ones:

1. Consider a system of N particles where no external or internal force acts on the center of mass (CM) coordinate $\mathbf{R} = \frac{1}{M} m_i \mathbf{r}_i$, where the total mass $M = \sum_i m_i$. Then the CM momentum \mathbf{P} is conserved. This is because

$$\mathbf{F}_{\mathbf{R}} = -\nabla_{\mathbf{R}}V = 0 \tag{1.48}$$

so V is independent of **R**. Meanwhile, $T = \frac{1}{2} \sum_i m_i \dot{\mathbf{r}}_i^2$, which when using coordinates relative to the center of mass, $\mathbf{r}_i' \equiv \mathbf{r}_i - \mathbf{R}$, becomes

$$T = \frac{1}{2} \left(\sum_{i} m_{i} \right) \dot{\mathbf{R}}^{2} + \dot{\mathbf{R}} \cdot \frac{d}{dt} \left(\sum_{i} m_{i} \mathbf{r}'_{i} \right) + \frac{1}{2} \sum_{i} m_{i} \dot{\mathbf{r}}'^{2}_{i} = \frac{1}{2} M \dot{\mathbf{R}}^{2} + \frac{1}{2} \sum_{i} m_{i} \dot{\mathbf{r}}'^{2}_{i}.$$
 (1.49)

Note that $\sum_i m_i \mathbf{r}'_i = 0$ from the definitions of M, \mathbf{R} , and \mathbf{r}'_i , so T splits into two terms, one for the CM motion and one for relative motion. We also observe that T is independent of \mathbf{R} . This means that \mathbf{R} is cyclic for the full Lagrangian L, so $\mathbf{P} = M\dot{\mathbf{R}}$ is a conserved quantity. In our study of rigid bodies we will also need the forms of M and \mathbf{R} for a continuous body with mass distribution $\rho(\mathbf{r})$, which for a three dimensional body are $M = \int d^3r \, \rho(\mathbf{r})$ and $\mathbf{R} = \frac{1}{M} \int d^3r \, \rho(\mathbf{r}) \, \mathbf{r}$.

Note that $\dot{\mathbf{P}} = 0$ is satisfied by having no total external force, so $\mathbf{F}^{\text{ext}} = \sum_{i} \mathbf{F}_{i}^{\text{ext}} = 0$, and by the internal forces obeying Newton's 3^{rd} law $\mathbf{F}_{i \to j} = -\mathbf{F}_{j \to i}$. Hence,

$$M\ddot{\mathbf{R}} = \sum_{i} \mathbf{F}_{i}^{\text{ext}} + \sum_{i,j} \mathbf{F}_{i \to j} = 0.$$
 (1.50)

2. Let us consider a system that is invariant with respect to rotations of angle ϕ about a symmetry axis. This has a conserved angular momentum. If we pick ϕ as a generalized coordinate, then L = T - V is independent of ϕ , so $\dot{p}_{\phi} = \frac{\partial L}{\partial \phi} = 0$ meaning p_{ϕ} is constant. In particular, for a system where V is independent of the angular velocity $\dot{\phi}$ we have

$$p_{\phi} = \frac{\partial T}{\partial \dot{\varphi}} = \sum_{i} m_{i} \dot{\mathbf{r}}_{i} \cdot \frac{\partial \dot{\mathbf{r}}_{i}}{\partial \dot{\varphi}} = \sum_{i} m_{i} \mathbf{v}_{i} \cdot \frac{\partial \mathbf{r}_{i}}{\partial \varphi}.$$
 (1.51)

Simplifying further using the results in Fig. 2.2 yields

$$p_{\varphi} = \sum_{i} m_{i} \mathbf{v}_{i} \cdot (\hat{n} \times \mathbf{r}_{i}) = \hat{n} \cdot \sum_{i} \mathbf{r}_{i} \times m_{i} \mathbf{v}_{i} = \hat{n} \cdot \mathbf{L}_{\text{total}}.$$
 (1.52)

Note that **L** about the CM is conserved for systems with no external torque, $\boldsymbol{\tau}^{\text{ext}} = \sum_{i} \mathbf{r}_{i} \times \mathbf{F}_{i}^{\text{ext}} = 0$ and internal forces that are all central. Defining $\mathbf{r}_{ij} \equiv \mathbf{r}_{i} - \mathbf{r}_{j}$ and its magnitude appropriately, this means $V_{ij} = V_{ij}(r_{ij})$. This implies that $\mathbf{F}_{ji} = -\nabla_{i}V_{ij}$ (no sum on the repeated index) is parallel to \mathbf{r}_{ij} . Hence,

$$\frac{d\mathbf{L}}{dt} = \sum_{i} \mathbf{r}_{i} \times \dot{\mathbf{p}}_{i} = \sum_{i} \mathbf{r}_{i} \times \mathbf{F}_{i}^{\text{ext}} + \sum_{i,j} \mathbf{r}_{i} \times \mathbf{F}_{ji}.$$
 (1.53)

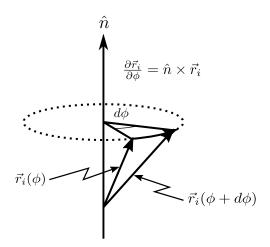


Figure 1.5: Rotation about a symmetry axis

However,
$$\sum_{i} \mathbf{r}_{i} \times \mathbf{F}_{i}^{\text{ext}} = 0$$
, so
$$\frac{d\mathbf{L}}{dt} = \sum_{i < j} \mathbf{r}_{ij} \times \mathbf{F}_{ji} = 0. \tag{1.54}$$

- 3. One can also consider a scaling transformation. Suppose that under the transformation $\mathbf{r}_i \to \lambda \mathbf{r}_i$ the potential is homogeneous and transforms as $V \to \lambda^k V$ for some constant k. Letting T be quadratic in $\dot{\mathbf{r}}_i$ and taking time to transform as $t \to \lambda^{1-k/2} t$ then gives $\dot{\mathbf{r}}_i \to \lambda^{k/2} \dot{\mathbf{r}}_i$. So by construction $T \to \lambda^k T$ also, and thus the full Lagrangian $L \to \lambda^k L$. This overall factor does not change the Euler-Lagrange equations, and hence the transformation is a symmetry of the dynamics, only changing the overall scale or units of the coordinate and time variables, but not their dynamical relationship. This can be applied for several well known potentials:
 - a) k=2 for a harmonic oscillator. In this case the scaling for time is given by 1-k/2=0, so it does not change with λ . The amplitude for the coordinates scales with λ . Thus, the frequency of the oscillator, which is a time variable, is independent of the amplitude.
 - b) k = -1 for the Coulomb potential. Here 1 k/2 = 3/2 so there is a more intricate relation between coordinates and time. This power is consistent with the behavior of bound state orbits, where the period of the orbit T obeys $T^2 \propto a^3$, for a the semi-major axis distance (Kepler's 3rd law).
 - c) k = 1 for a uniform gravitational field. Here 1 k/2 = 1/2 so for a freely falling object, the time of free fall goes as \sqrt{h} where h is the distance fallen.
- 4. Consider the Lagrangian for a charge in the presence of scalar (ϕ) and vector (\mathbf{A}) electromagnetic fields, $L = \frac{1}{2}m\dot{\mathbf{r}}^2 e\phi + e\mathbf{A} \cdot \dot{\mathbf{r}}$. As a concrete example, let us take ϕ and \mathbf{A} to be independent of the Cartesian coordinate x. The canonical momentum is

 $\mathbf{p} = \frac{\partial L}{\partial \dot{\mathbf{r}}} = m\dot{\mathbf{r}} + e\mathbf{A}$, which is notably different from the kinetic momentum. Then x being cyclic means the canonical momentum p_x is conserved.

5. Let us consider the conservation of energy and the relationship between energy and the Hamiltonian. Applying the time derivative gives $\dot{H} = \frac{\partial H}{\partial q} \dot{q} + \frac{\partial H}{\partial p} \dot{p} + \frac{\partial H}{\partial t}$. However, $\dot{q} = \frac{\partial H}{\partial p}$ and $\dot{p} = -\frac{\partial H}{\partial q}$. Thus

$$\dot{H} = \frac{\partial H}{\partial t} = -\frac{\partial L}{\partial t} \,. \tag{1.55}$$

There are two things to consider.

- If H (or L) has no explicit time dependence, then $H = \dot{q}_i p_i L$ is conserved.
- Energy is conserved if $\dot{E} = 0$, where energy is defined by E = T + V.

If H = E then the two points are equivalent, but otherwise either of the two could be true while the other is false.

Example: Let us consider a system which provides an example where H = E but energy is not conserved, and where $H \neq E$ but H is conserved. The two situations will be obtained from the same example by exploiting a coordinate choice. Consider a system consisting of a mass m attached by a spring, with spring constant k, to a cart moving at a constant speed v_0 in one dimension, as shown in Fig. 1.6. Let us call x

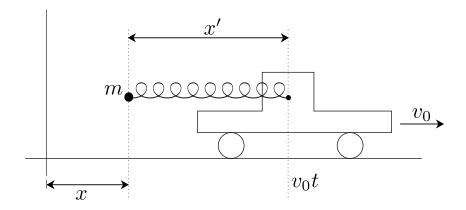


Figure 1.6: Mass attached by a spring to a moving cart

the displacement of m from the fixed wall and x' is its displacement from the center of the moving cart. Using x,

$$L(x, \dot{x}) = T - V = \frac{m}{2}\dot{x}^2 - \frac{k}{2}(x - v_0 t)^2,$$
(1.56)

where the kinetic term is quadratic in \dot{x} and the potential term is independent of \dot{x} . This means that H falls in the special case considered in Eq. (1.31) so

$$H = E = T + V = \frac{p^2}{2m} + \frac{k}{2} (x - v_0 t)^2, \qquad (1.57)$$

However $\frac{\partial H}{\partial t} \neq 0$ so the energy is not conserved. (Of course the full energy would be conserved, but we have not accounted for the energy needed to pull the cart at a constant velocity, treating that instead as external to our system. That is what led to the time dependent H.)

If we instead choose to use the coordinate $x' = x - v_0 t$, then

$$L'(x', \dot{x}') = \frac{m}{2}\dot{x}'^2 + mv_0x' + \frac{m}{2}v_0^2 - \frac{k}{2}x'^2.$$
 (1.58)

Note that $p' = m\dot{x}' + mv_0 = m\dot{x} = p$. This Lagrangian fits the general form in equation (1.32) with $a = mv_0$ and $L_0 = mv_0^2/2 - kx'^2/2$. So

$$H'(x',p') = \dot{x}'p' - L' = \frac{1}{2m} (p' - mv_0)^2 + \frac{k}{2}x'^2 - \frac{m}{2}v_0^2, \tag{1.59}$$

Here the last terms is a constant shift. The first and second terms in this expression for H' look kind of like the energy that we would calculate if we were sitting on the cart and did not know it was moving, which is not the same as the energy above. Hence, $H' \neq E$, but $\dot{H}' = 0$ because $\frac{\partial H'}{\partial t} = 0$, so H' is conserved.

1.4 Constraints and Friction Forces

So far, we've considered constraints to a surface or curve that are relationships between coordinates. These fall in the category of *holonomic constraints*. Such constraints take the form

$$f(q_1, \dots, q_N, t) = 0 (1.60)$$

where explicit time dependence is allowed as a possibility. An example of holonomic constrain is mass in a cone (Figure 1.4), where the constrain is $z - r \cot \alpha = 0$. Constraints that violate the form in Eq. (1.60) are non-holonomic constraints.

- An example of a non-holonomic constraint is a mass on the surface of a sphere. The constraint here is an *inequality* $r^2 a^2 \ge 0$ where r is the radial coordinate and a is the radius of the sphere.
- Another example of a non-holonomic constraint is an object rolling on a surface without slipping. The point of contact is stationary, so the constraint is actually on the velocities.

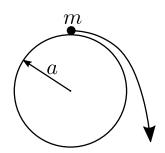


Figure 1.7: Mass on a sphere

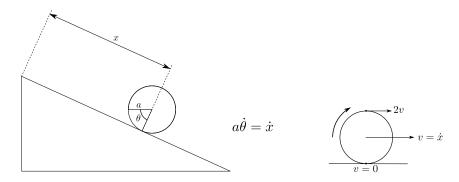


Figure 1.8: Disk rolling down an incline without slipping

A simple example is a disk of radius a rolling down an inclined plane without slipping, as shown in Fig. 1.8. Here the condition on velocities, $a\dot{\theta} = \dot{x}$ is simple enough that it can be integrated into a holonomic constraint.

As a more sophisticated example, consider a vertical disk of radius a rolling on a horizontal plane, as shown in Fig. 1.9. The coordinates are (x, y, θ, ϕ) , where (x, y) is the point of contact, ϕ is the rotation angle about its own axis, and θ is the angle of orientation along the xy-plane. We will assume that the flat edge of the disk always remains parallel to z, so the disk never tips over. The no-slip condition is $v = a\dot{\phi}$ where \mathbf{v} is the velocity of the center of the disk, and $v = |\mathbf{v}|$. This means $\dot{x} = v\sin(\theta) = a\sin(\theta)\dot{\phi}$ and $\dot{y} = -v\cos(\theta) = -a\cos(\theta)\dot{\phi}$, or in differential notation, $dx - a\sin(\theta)d\phi = 0$ and $dy + a\cos(\theta)d\phi = 0$.

In general, constraints of the form

$$\sum_{j} a_{j}(q)dq_{j} + a_{t}(q)dt = 0$$
(1.61)

are not holonomic. We will call this a semi-holonomic constraint.

Let us consider the special case of a holonomic constraint in differential form, $f(q_1, ..., q_{3N}, t) = 0$. This means

$$df = \sum_{j} \frac{\partial f}{\partial q_{j}} dq_{j} + \frac{\partial f}{\partial t} dt = 0, \qquad (1.62)$$

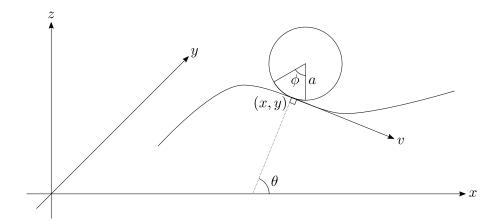


Figure 1.9: Vertical rolling disk on a two dimensional plane

so $a_j = \frac{\partial f}{\partial q_j}$ and $a_t = \frac{\partial f}{\partial t}$. The properties of mixed partial derivatives means

$$\frac{\partial a_j}{\partial q_i} = \frac{\partial a_i}{\partial q_j}, \qquad \frac{\partial a_t}{\partial q_i} = \frac{\partial a_i}{\partial t}. \qquad (1.63)$$

These conditions imply that a seemingly semi-holonomic constraint is in fact holonomic. (In math we would say that we have an exact differential form df for the holonomic case, but the differential form in Eq.(1.61) need not always be exact.)

Example: To demonstrate that not all semiholonomic constrants are secretly holonomic, consider the constraint in the example of the vertical disk. Here there is no function $h(x, y, \theta, \phi)$ that we can multiply the constraint df = 0 by to make it holonomic. For the vertical disk from before, we could try $(dx - a\sin(\theta) d\phi)h = 0$ with $a_x = h$, $a_{\phi} = -a\sin(\theta)h$, $a_{\theta} = 0$, and $a_y = 0$, all for some function h. As we must have $\frac{\partial a_{\phi}}{\partial \theta} = \frac{\partial a_{\theta}}{\partial \phi}$, then $0 = -a\cos(\theta) - a\sin(\theta)\frac{\partial h}{\partial \theta}$, so $h = \frac{k}{\sin(\theta)}$. That said, $\frac{\partial a_x}{\partial \theta} = \frac{\partial a_{\theta}}{\partial x}$ gives $\frac{\partial h}{\partial \theta} = 0$, yielding a contradiction for any non-trivial h with $k \neq 0$.

If the rolling is instead constrained to a line rather than a plane, then the constraint is holonomic. Take as an example $\theta = \frac{\pi}{2}$ for rolling along \hat{x} , then $\dot{x} = a\dot{\phi}$ and $\dot{y} = 0$. Integrating we have $x = a\varphi + x_0$, $y = y_0$, and $\theta = \frac{\pi}{2}$, which together form a set of holonomic constraints.

A useful concept for discussing constraints is that of the *virtual displacement* $\delta \mathbf{r}_i$ of particle *i*. There are a few properties to be noted of $\delta \mathbf{r}_i$.

- It is infinitesimal.
- It is consistent with the constraints.
- It is carried out at a fixed time (so time dependent constraints do not change its form).

Example: let us consider a bead constrained to a moving wire. The wire is oriented along

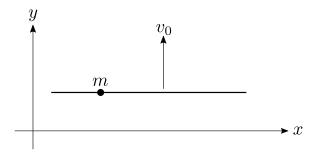


Figure 1.10: Bead on a moving wire

the x-axis and is moving with coordinate $y = v_0 t$. Here the virtual displacement of the bead $\delta \mathbf{r}$ is always parallel to \hat{x} (since it is determined at a fixed time), whereas the real displacement $d\mathbf{r}$ has a component along \hat{y} in a time interval dt.

For a large number of constraints, the constraint force \mathbf{Z}_i is perpendicular to $\delta \mathbf{r}_i$, meaning $\mathbf{Z}_i \cdot \delta \mathbf{r}_i = 0$, so the "virtual work" (in analogy to work $W = \int \mathbf{F} \cdot d\mathbf{r}$) of a constraint force vanishes. More generally, there is no *net* work from constraints, so $\sum_i \mathbf{Z}_i \cdot \delta \mathbf{r}_i = 0$ (which holds for the actions of surfaces, rolling constraints, and similar things). The Newtonian equation of motion is $\dot{\mathbf{p}}_i = \mathbf{F}_i + \mathbf{Z}_i$, where \mathbf{F}_i encapsulates other forces. Vanishing virtual work gives

$$\sum_{i} (\dot{\mathbf{p}}_i - \mathbf{F}_i) \cdot \delta \mathbf{r}_i = 0 \tag{1.64}$$

which is the D'Alembert principle. This could be taken as the starting principal for classical mechanics instead of the Hamilton principle of stationary action.

Of course Eq.(1.64) is not fully satisfactory since we are now used to the idea of working with generalized coordinates rather than the cartesian vector coordinates used there. So lets transform to generalized coordinates through $\mathbf{r}_i = \mathbf{r}_i(q,t)$, so $\delta \mathbf{r}_i = \frac{\partial \mathbf{r}_i}{\partial q_j} \delta q_j$, where again we sum over repeated indicies (like j here). This means

$$\mathbf{F}_{i} \cdot \delta \mathbf{r}_{i} = \mathbf{F}_{i} \cdot \frac{\partial \mathbf{r}_{i}}{\partial q_{i}} \delta q_{j} \equiv Q_{j} \delta q_{j} \tag{1.65}$$

where we have defined generalized forces

$$Q_j \equiv \mathbf{F}_i \cdot \frac{\partial \mathbf{r}_i}{\partial q_i} \,. \tag{1.66}$$

We can also transform the $\dot{\mathbf{p}}_i \cdot \delta \mathbf{r}_i$ term using our earlier point transformation results as well as the fact that $\frac{d}{dt} \left(\frac{\partial \mathbf{r}_i}{\partial q_j} \right) = \frac{\partial^2 \mathbf{r}_i}{\partial q_j \partial t} + \sum_k \frac{\partial^2 \mathbf{r}_i}{\partial q_j \partial q_k} \dot{q}_k = \frac{\partial \mathbf{v}_i}{\partial q_j}$. Writing out the index sums explicitly,

this gives

$$\sum_{i} \dot{\mathbf{p}}_{i} \cdot \delta \mathbf{r}_{i} = \sum_{i,j} m_{i} \ddot{\mathbf{r}}_{i} \cdot \frac{\partial \mathbf{r}_{i}}{\partial q_{j}} \delta q_{j}$$

$$= \sum_{i,j} \left(\frac{d}{dt} \left(m_{i} \dot{\mathbf{r}}_{i} \cdot \frac{\partial \mathbf{r}_{i}}{\partial q_{j}} \right) - m_{i} \dot{\mathbf{r}}_{i} \cdot \frac{d}{dt} \left(\frac{\partial \mathbf{r}_{i}}{\partial q_{j}} \right) \right) \delta q_{j}$$

$$= \sum_{i,j} \left(\frac{d}{dt} \left(m_{i} \mathbf{v}_{i} \cdot \frac{\partial \mathbf{v}_{i}}{\partial \dot{q}_{j}} \right) - m_{i} \mathbf{v}_{i} \cdot \frac{\partial \mathbf{v}_{i}}{\partial q_{j}} \right) \delta q_{j}$$

$$= \sum_{i} \left(\frac{d}{dt} \left(\frac{\partial T}{\partial \dot{q}_{i}} \right) - \frac{\partial T}{\partial q_{j}} \right) \delta q_{j} \tag{1.67}$$

for $T = \frac{1}{2} \sum_{i} m_{i} \mathbf{v}_{i}^{2}$. Together with the D'Alembert principle, we obtain the final result

$$\sum_{i} \left(\frac{d}{dt} \left(\frac{\partial T}{\partial \dot{q}_{j}} \right) - \frac{\partial T}{\partial q_{j}} - Q_{j} \right) \delta q_{j} = 0.$$
 (1.68)

We will see momentarily that this result is somewhat more general than the Euler-Lagrange equations, containing them as a special case.

Holonomic Constraints:

We will start by considering systems with only holonomic constraints, postponing other types of constraints to the next section. Here we can find the independent coordinates q_j with j = 1, ..., N-k that satisfy the k constraints. This implies that the generalized virtual displacements δq_j are independent, so that their coefficients in Eq. (1.68) must vanish,

$$\frac{d}{dt} \left(\frac{\partial T}{\partial \dot{q}_j} \right) - \frac{\partial T}{\partial q_j} - Q_j = 0, \quad \text{for } j = 1, \dots, N - k.$$
 (1.69)

There are several special cases of this result, which we derived from the d'Alembert principle.

1. For a conservative force $\mathbf{F}_i = -\nabla_i V$, then

$$Q_{j} = -(\nabla_{i}V) \cdot \frac{\partial \mathbf{r}_{i}}{\partial q_{j}} = -\frac{\partial V}{\partial q_{j}}$$
(1.70)

where we assume that the potential can be expressed in the generalized coordinates as V = V(q, t). Then using L = T - V, we see that Eq. (1.69) simply reproduces the Euler-Lagrange equations $\frac{d}{dt} \left(\frac{\partial L}{\partial \dot{q}_j} \right) - \frac{\partial L}{\partial q_j} = 0$.

2. If $Q_j = -\frac{\partial V}{\partial q_j} + \frac{d}{dt} \left(\frac{\partial V}{\partial \dot{q}_j} \right)$ for $V = V(q, \dot{q}, t)$, which is the case for velocity dependent forces derivable from a potential (like the electromagnetic Lorentz force), then the Euler-Lagrange equations $\frac{d}{dt} \left(\frac{\partial L}{\partial \dot{q}_j} \right) - \frac{\partial L}{\partial q_j} = 0$ are again reproduced.

3. If Q_j has forces obtainable from a potential as in case 2, as well as generalized forces R_j that cannot, then

$$\frac{d}{dt}\left(\frac{\partial L}{\partial \dot{q}_i}\right) - \frac{\partial L}{\partial q_i} = R_j \tag{1.71}$$

is the generalization of the Euler-Lagrange equations with non-conservative generalized forces.

We will continue to use the notation R_j for generalized forces that can not be obtained from a potential. An important example of such nonconservative forces R_j is given by friction.

- Static friction is $F_s \leq F_s^{\text{max}} = \mu_s F_N$ for a normal force F_N .
- Sliding friction is $\mathbf{F} = -\mu F_{\rm N} \frac{\mathbf{v}}{v}$, so this is a constant force that is always opposite the direction of motion (but vanishes when there is no motion).
- Rolling friction is $\mathbf{F} = -\mu_{\rm R} F_{\rm N} \frac{\mathbf{v}}{v}$.
- Fluid friction at a low velocity is $\mathbf{F} = -bv\frac{\mathbf{v}}{v} = -b\mathbf{v}$.

A general form for a friction force is $\mathbf{F}_i = -h_i(v_i) \frac{\mathbf{v}_i}{v_i}$ (where as a reminder there is no implicit sum on i here since we specified i on the left-hand-side). For this form

$$R_{j} = -\sum_{i} h_{i} \frac{\mathbf{v}_{i}}{v_{i}} \cdot \frac{\partial \mathbf{r}_{i}}{\partial q_{j}} = -\sum_{i} h_{i} \frac{\mathbf{v}_{i}}{v_{i}} \cdot \frac{\partial \mathbf{v}_{i}}{\partial \dot{q}_{j}}.$$
(1.72)

Simplifying further gives

$$R_{j} = -\sum_{i} \frac{h_{i}}{2v_{i}} \frac{\partial}{\partial \dot{q}_{j}} \left(\mathbf{v}_{i}^{2} \right) = -\sum_{i} h_{i} \frac{\partial v_{i}}{\partial \dot{q}_{j}} = -\sum_{i} \frac{\partial v_{i}}{\partial \dot{q}_{j}} \frac{\partial}{\partial v_{i}} \int_{0}^{v_{i}} dv'_{i} h_{i}(v'_{i}) = -\frac{\partial}{\partial \dot{q}_{j}} \sum_{i} \int_{0}^{v_{i}} dv'_{i} h_{i}(v'_{i}) = -\frac{\partial}{\partial \dot{q}_{i}} \sum_{i} \int_{0}^{v_{i}} dv'_{i} h_{i}(v'_{i}) dv'_{i} h_{i}(v'_{i}) = -\frac{\partial}{\partial \dot{q}_{i}} \sum_{i} \int_{0}^{v_{$$

where

$$\mathcal{F} = \sum_{i} \int_{0}^{v_i} dv_i' h_i(v_i') \tag{1.74}$$

is the "dissipation function". This is a scalar function like L so it is relatively easy to work with.

Example: Consider a sphere of radius a and mass m falling in a viscous fluid. Then $T = \frac{1}{2}m'\dot{y}^2$ where m' < m accounts for the mass of displaced fluid (recall Archimedes principle that the buoyant force on a body is equal to the weight of fluid the body displaces). Also

V=m'gy, and L=T-V. Here $h\propto\dot{y}$, so $\mathcal{F}=3\pi\eta a\dot{y}^2$, where by the constant of proportionality is determined by the constant η , which is the viscosity. From this, $\frac{d}{dt}\left(\frac{\partial L}{\partial\dot{y}}\right)-\frac{\partial L}{\partial y}=-\frac{\partial \mathcal{F}}{\partial\dot{y}}$ gives the equation of motion $m'\ddot{y}+m'g=-6\pi\eta a\dot{y}$. The friction force $6\pi\eta a\dot{y}$ is known as Stokes Law. (We will derive this equation for the friction force from first principles later on, in our discussion of fluids.) This differential equation can be solved by adding a particular solution $y_p(t)$ to a solution of the homogeneous equation $m'\ddot{y}_H+6\pi\eta a\dot{y}_H=0$. For the time derivatives the results are $\dot{y}_p=-m'g/(6\pi\eta a)$ and $\dot{y}_H=A\exp(-6\pi\eta at/m')$, where the constant A must be determined by an initial condition. The result $\dot{y}=\dot{y}_H+\dot{y}_p$ can be integrated in time once more to obtain the full solution y(t) for the motion.

Example: if we add sliding friction to the case of two masses on a plane connected by a spring (considered on problem set #1), then $h_i = \mu_f m_i g$ for some friction coefficient μ_f , and

$$\mathcal{F} = \mu_f g(m_1 v_1 + m_2 v_2) = \mu_f g\left(m_1 \sqrt{\dot{x}_1^2 + \dot{y}_1^2} + m_2 \sqrt{\dot{x}_2^2 + \dot{y}_2^2}\right). \tag{1.75}$$

If we switch to a suitable set of generalized coordinates q_j that simplify the equations of motion without friction, and then compute the generalized friction forces $R_j = -\frac{\partial \mathcal{F}}{\partial \dot{q}_j}$, we can get the equations of motion including friction. Further details of how this friction complicates the equations of motion were provided by an example done in lecture.

1.5 Calculus of Variations & Lagrange Multipliers

Calculus of Variations

In the calculus of variations, we wish to find a set of functions $y_i(s)$ between s_1 and s_2 that extremize the following functional (a function of functions),

$$J[y_i] = \int_{s_1}^{s_2} ds \, f(y_1(s), \dots, y_n(s), \dot{y}_1(s), \dots, \dot{y}_n(s), s) , \qquad (1.76)$$

where for this general discussion only we let $\dot{y}_i \equiv \frac{dy_i}{ds}$ rather than $\frac{d}{dt}$. To consider the action of the functional under a variation we consider $y_i'(s) = y_i(s) + \eta_i(s)$ where $\eta_i(s_1) = \eta_i(s_2) = 0$, meaning that while the two endpoints are fixed during the variation $\delta y_i = \eta_i$, the path in between is varied. Expanding the variation of the functional integral $\delta J = J[y_i'] - J[y_i] = 0$ to 1^{st} order in δy_i we have

$$0 = \delta J = \int_{s_1}^{s_2} ds \sum_{i} \left[\delta y_i \frac{\partial f}{\partial y_i} + \delta \dot{y}_i \frac{\partial f}{\partial \dot{y}_i} \right]. \tag{1.77}$$

Using integration by parts on the second term, and the vanishing of the variation at the endpoints to remove the surface term, δJ vanishes when

$$\int_{s_1}^{s_2} \sum_{i} \left[\frac{\partial f}{\partial y_i} - \frac{d}{ds} \left(\frac{\partial f}{\partial \dot{y}_i} \right) \right] \delta y_i(s) \, ds = 0.$$
 (1.78)

If we identify a set of independent variations δy_j (for example, after imposing holonomic constraints), this can only occur if

$$\frac{\partial f}{\partial y_j} - \frac{d}{ds} \left(\frac{\partial f}{\partial \dot{y}_j} \right) = 0. \tag{1.79}$$

The scope of this calculus of variation result for extremizing the integral over f is more general than its application to classical mechanics.

Example: Hamilton's principle states that motion $q_i(t)$ extremizes the action, so in this case s = t, $y_i = q_i$, f = L, and J = S. Demanding $\delta S = 0$ then yields the Euler-Lagrange equations of motion from Eq. (1.79).

Example: As an example outside of classical mechanics, consider showing that the shortest distance between points on a sphere of radius a are great circles. This can be seen by minimizing the distance $J = \int_{s_1}^{s_2} ds$ where for a spherical surface,

$$ds = \sqrt{(dx)^2 + (dy)^2 + (dz)^2} = \sqrt{a^2(d\theta)^2 + a^2 \sin^2(\theta)(d\phi)^2}$$
(1.80)

since dr = 0. Taking $s = \theta$ and $y = \phi$, then

$$ds = a\sqrt{1 + \sin^2(\theta) \left(\frac{d\phi}{d\theta}\right)^2} d\theta, \tag{1.81}$$

so $f = \sqrt{1 + \sin^2(\theta)\dot{\phi}^2}$. The solution for the minimal path is given by solving $\frac{d}{d\theta}\left(\frac{\partial f}{\partial \dot{\varphi}}\right) - \frac{\partial f}{\partial \varphi} = 0$. After some algebra these are indeed found to be great circles, described by $\sin(\phi - \alpha) = \beta \cot(\theta)$ where α , β are constants.

Example: Hamilton's principle can also be used to yield the Hamilton equations of motion, by considering the variation of a path in phase space. In this case

$$\delta J[q, p] = \delta \int_{t_1}^{t_2} dt \left[p_i \dot{q}_i - H(q, p, t) \right] = 0$$
 (1.82)

must be solved with fixed endpoints: $\delta q_i(t_1) = \delta q_i(t_2) = 0$ and $\delta p_i(t_1) = \delta p_i(t_2) = 0$. Here, the role of y_i is played by the 2N variables $(q_1, \ldots, q_N, p_1, \ldots, p_N)$. As $f = p_i \dot{q}_i - H$, then

$$\frac{d}{dt} \left(\frac{\partial f}{\partial \dot{q}_i} \right) - \frac{\partial f}{\partial q_i} = 0 \qquad \Longrightarrow \qquad \dot{p}_i = -\frac{\partial H}{\partial q_i}, \qquad (1.83)$$

$$\frac{d}{dt} \left(\frac{\partial f}{\partial \dot{p}_i} \right) - \frac{\partial f}{\partial p_i} = 0 \qquad \Longrightarrow \qquad \dot{q}_i = \frac{\partial H}{\partial p_i},$$

giving Hamilton's equations as expected. Note that because f is independent of \dot{p}_i , the term $(\partial f/\partial \dot{p}_i)\delta \dot{p}_i = 0$, and it would seem that we do not really need the condition that $\delta p_i(t_1) = \delta p_i(t_2) = 0$ to remove the surface term. However, these conditions on the variations δp_i are actually required in order to put q_i and p_i on the same footing, including making transformation that mix up these variables (we will exploit this in detail later on when discussing canonical transformations).

It is interesting and useful to note that D'Alembert's principle

$$\left(\frac{d}{dt}\left(\frac{\partial L}{\partial \dot{q}_i}\right) - \frac{\partial L}{\partial q_i} - R_j\right)\delta q_j = 0$$
(1.84)

is a "differential" version of the equations that encode the classical dynamics, while Hamilton's principle

$$\delta J = \int_{t_1}^{t_2} dt \, \left(\frac{\partial L}{\partial q_j} - \frac{d}{dt} \left(\frac{\partial L}{\partial \dot{q}_j} \right) \right) \delta q_j = 0 \tag{1.85}$$

(for $R_j = 0$ where all forces come from a potential) is an integrated version.

Method of Lagrange Multipliers

Next we will consider the method of Lagrange multipliers. For simplicity we will assume there are no generalized forces outside the potential, $R_j = 0$, until further notice. (It is straightforward to add back the R_j to the final equations derived below.) The method of Lagrange multipliers will be useful for two situations that we will encounter:

- 1. When we actually want to study the forces of constraint that are holonomic.
- 2. When we have semi-holonomic constraints.

Let us consider k constraints for n coordinates, with $\alpha \in \{1, \ldots, k\}$ being the index running over the constraints. These holonomic or semi-holonomic constraints take the form

$$g_{\alpha}(q,\dot{q},t) = a_{j\alpha}(q,t)\dot{q}_j + a_{t\alpha}(q,t) = 0$$
 (1.86)

where again repeated indices are summed. Thus, $g_{\alpha}dt = a_{j\alpha}dq_j + a_{t\alpha}dt = 0$. For a virtual displacement δq_j we have dt = 0, so

$$\sum_{j=1}^{n} a_{j\alpha} \delta q_j = 0, \qquad (1.87)$$

which gives us k equations constraining the virtual displacements. For each equation we can multiply by a function $\lambda_{\alpha}(t)$ known as Lagrange multipliers, and sum over α , and the combination will still be zero. Adding this zero to D'Alembert's principle (taking $R_j = 0$) yields

$$\left[\frac{d}{dt}\left(\frac{\partial L}{\partial \dot{q}_j}\right) - \frac{\partial L}{\partial q_j} - \lambda_{\alpha} a_{j\alpha}\right] \delta q_j = 0, \qquad (1.88)$$

where the sums implicitly run over both α and j. It is clear that the Lagrange multiplier term is zero if we sum over j first, but now we want to consider summing first over α for a fixed j. Our goal is make the term in square brackets zero. Only n-k of the virtual displacements δq_j are independent, so for these values of j the square brackets must vanish. For the remaining k values of j we can simply choose the k Lagrange multipliers λ_{α} to force the k square bracketed equations to be satisfied. This is known as the method of Lagrange multipliers. Thus all square bracketed terms are zero, and we have the generalization of the Euler-Lagrange equations which includes terms for the constraints:

$$\frac{d}{dt} \left(\frac{\partial L}{\partial \dot{q}_j} \right) - \frac{\partial L}{\partial q_j} = \lambda_{\alpha} a_{j\alpha} \,. \tag{1.89}$$

This is n equations, for the n possible values of j, and on the right-hand-side we sum over α for each one of these equations. The sum $\lambda_{\alpha}a_{j\alpha}$ can be interpreted as a generalized constraint force Q_j . The Lagrange multipliers λ_{α} and generalized coordinates q_j together form n + k parameters, and equation (1.89) in conjunction with $g_{\alpha} = 0$ for each α from (1.86) together form n + k equations to be solved.

There are two important cases to be considered.

1. In the holonomic case, $f_{\alpha}(q,t) = 0$. Here, $g_{\alpha} = \dot{f}_{\alpha} = \frac{\partial f_{\alpha}}{\partial q_{j}} \dot{q}_{j} + \frac{\partial f_{\alpha}}{\partial t}$, so $a_{j\alpha} = \frac{\partial f_{\alpha}}{\partial q_{j}}$. This gives

$$\frac{d}{dt} \left(\frac{\partial L}{\partial \dot{q}_j} \right) - \frac{\partial L}{\partial q_j} = \sum_{\alpha=1}^k \lambda_\alpha \frac{\partial f_\alpha}{\partial q_j} \tag{1.90}$$

for holonomic constraints. The same result can be derived from a generalized Hamilton's principle

$$J[q_j, \lambda_\alpha] = \int_{t_1}^{t_2} (L + \lambda_\alpha f_\alpha) dt$$
 (1.91)

by demanding that $\delta J = 0$. It is convenient to think of $-\lambda_{\alpha} f_{\alpha}$ as an extra potential energy that we add into L so that a particle does work if it leaves the surface defined by $f_{\alpha} = 0$. Recall that given this potential, the Force_q = $-\nabla_q(-\lambda_{\alpha} f_{\alpha}) = \lambda_{\alpha} \nabla_q f_{\alpha}$, where the derivative $\nabla_q f_{\alpha}$ gives a vector that is normal to the constraint surface of constant $f_{\alpha} = 0$. This agrees with the form of our generalized force above.

2. In the semi-holonomic case, we just have $g_{\alpha} = a_{j\alpha}(q,t)\dot{q}_j + a_{t\alpha}(q,t) = 0$, with $a_{j\alpha} = \frac{\partial g_{\alpha}}{\partial \dot{q}_j}$. This gives

$$\frac{d}{dt} \left(\frac{\partial L}{\partial \dot{q}_j} \right) - \frac{\partial L}{\partial q_j} = \sum_{\alpha=1}^k \lambda_\alpha \frac{\partial g_\alpha}{\partial \dot{q}_j} \tag{1.92}$$

for semi-holonomic constraints. This result cannot be derived from Hamilton's principle in general. This justifies the time we spent discussing d'Alembert's principle, which we have used to obtain (1.92). Recall that static friction imposes a no-slip constraint in the form of our equation $g_{\alpha} = 0$. For $g \propto \dot{q}$, the form, $\frac{\partial g}{\partial \dot{q}}$, is consistent with the form of generalized force we derived from our dissipation function, $\frac{\partial \mathcal{F}}{\partial \dot{q}}$ in our discussion of friction.

We end this chapter with several examples of the use of Lagrange multipliers.

Example: Consider a particle of mass m at rest on the top of a sphere of radius a, as shown above in Fig. 1.7. The particle is given an infinitesimal displacement $\theta = \theta_0$ so that it slides down. At what angle does it leave the sphere?

We use the coordinates (r, θ, ϕ) but set $\phi = 0$ by symmetry as it is not important. The constraint $r \geq a$ is non-holonomic, but while the particle is in contact with the sphere the constraint f = r - a = 0 is holonomic. To answer this question we will look for the point where the constraint force vanishes. Here $T = \frac{m}{2} \left(\dot{r}^2 + r^2 \dot{\theta}^2 \right)$ and $V = mgz = mgr \cos(\theta)$ so that L = T - V, then $\frac{d}{dt} \left(\frac{\partial L}{\partial \dot{r}} \right) - \frac{\partial L}{\partial r} = \lambda \frac{\partial f}{\partial r}$ gives

$$m\ddot{r} - mr\dot{\theta}^2 + mg\cos(\theta) = \lambda, \tag{1.93}$$

while $\frac{d}{dt} \left(\frac{\partial L}{\partial \dot{\theta}} \right) - \frac{\partial L}{\partial \theta} = \lambda \frac{\partial f}{\partial \theta} = 0$ gives

$$\frac{d}{dt}\left(mr^2\dot{\theta}\right) - mgr\sin(\theta) = 0. \tag{1.94}$$

This in conjunction with r = a gives 3 equations for the 3 variables (r, θ, λ) . At this point we can now consider solving the equations together. The constraint equation gives $\dot{r} = 0$ so $\ddot{r} = 0$. This means

$$ma^2\ddot{\theta} = mga\sin(\theta), \qquad -ma\dot{\theta}^2 + mg\cos(\theta) = \lambda.$$

Multiply the first of these by $\dot{\theta}$ and integrate over time, knowing that $\dot{\theta}=0$ when $\theta=0$, gives $\dot{\theta}^2=\frac{2g}{a}\left(1-\cos(\theta)\right)$. Plugging this into the second equation we find that

$$\lambda = mg(3\cos(\theta) - 2) \tag{1.95}$$

is the radial constraint force. The mass leaves the sphere when $\lambda = 0$ which is when $\cos(\theta) = \frac{2}{3}$ (so $\theta \approx 48^{\circ}$).

What if we instead imposed the constraint $f' = r^2 - a^2 = 0$? If we call its Lagrange multiplier λ' we would get $\lambda' \frac{\partial f'}{\partial r} = 2a\lambda'$ when r = a, so $2a\lambda' = \lambda$ is the constraint force from before. The meaning of λ' is different, and it has different units, but we still have the same constraint force.

What are the equations of motion for $\theta > \arccos(\frac{2}{3})$? Now we no longer have the constraint so

$$m\ddot{r} - mr\dot{\theta}^2 + mg\cos(\theta) = 0$$
 and $\frac{d}{dt}\left(mr^2\dot{\theta}\right) - mgr\sin(\theta) = 0.$

The initial conditions are $r_1 = a$, $\theta_1 = \arccos\left(\frac{2}{3}\right)$, $\dot{r}_1 = 0$, and $\dot{\theta}_1^2 = \frac{2g}{3a}$ from before. Simpler coordinates are $x = r\sin(\theta)$ and $z = r\cos(\theta)$, giving

$$L = \frac{m}{2} \left(\dot{x}^2 + \dot{z}^2 \right) - mgz, \tag{1.96}$$

so $\ddot{x}=0$ and $\ddot{z}=-g$ with initial conditions $z_1=\frac{2a}{3},\ x_1=\frac{\sqrt{5}a}{3}$, and the initial velocities $v_{z1}=\dot{z}(t_1)=-a\sin\theta_1\dot{\theta}_1=-\frac{\sqrt{5}a}{3}\left(\frac{2g}{3a}\right)^{1/2}$ and $v_{x1}=\dot{x}(t_1)=a\cos\theta_1\dot{\theta}_1=\frac{2a}{3}\left(\frac{2g}{3a}\right)^{1/2}$. Here t_1 is the time when the mass leaves the sphere. This means

$$x(t) = v_{x1}(t - t_1) + x_1, (1.97)$$

$$z(t) = -\frac{g}{2} (t - t_1)^2 + v_{z1}(t - t_1) + z_1.$$
(1.98)

The time t_1 can be found from

$$\dot{\theta}^2 = \frac{2g}{a} \left(1 - \cos(\theta) \right) = \frac{4g}{a} \sin^2\left(\frac{\theta}{2}\right),\tag{1.99}$$

so $t_1 = \sqrt{\frac{a}{4g}} \int_{\theta_0}^{\arccos(\frac{2}{3})} \frac{d\theta}{\sin(\frac{\theta}{2})}$ where θ_0 is the small initial angular displacement from the top of the sphere. Note that the calculation of t_1 is the only place where we see that it was important to have a small θ_0 different from zero.

Example: Consider a hoop of radius a and mass m rolling down an inclined plane of angle ϕ without slipping as shown in Fig. 1.11, where we define the \hat{x} direction as being parallel to the ramp as shown. What is the friction force of constraint, and how does the acceleration compare to the case where the hoop is sliding rather than rolling?

The no-slip constraint means $a\dot{\theta} = \dot{x}$, so $h = a\dot{\theta} - \dot{x} = 0$, which can be made holonomic but which we will treat as semi-holonomic. Here the kinetic energy has contributions from the center mass motion of the hoop and its rotation, $T = T_{\rm CM} + T_{\rm rotation} = \frac{1}{2}m\dot{x}^2 + \frac{1}{2}ma^2\dot{\theta}^2$, where the hoops moment of inertia is $I_{\rm hoop} = ma^2$. (We will discuss moments of inertia in detail in chapter 2.) Meanwhile, $V = mg(l-x)\sin(\phi)$ where we are free to pick the length l where V(x=l)=0, such as at the bottom of the ramp. This means

$$L = T - V = \frac{m}{2}\dot{x}^2 + \frac{ma^2}{2}\dot{\theta}^2 + mg(x - l)\sin(\phi). \tag{1.100}$$

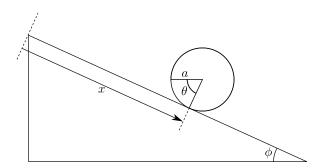


Figure 1.11: Hoop rolling on inclined plane

The equations of motion from $\frac{d}{dt} \left(\frac{\partial L}{\partial \dot{x}} \right) - \frac{\partial L}{\partial x} = \lambda \frac{\partial h}{\partial \dot{x}}$ and $\frac{d}{dt} \left(\frac{\partial L}{\partial \dot{\theta}} \right) - \frac{\partial L}{\partial \theta} = \lambda \frac{\partial h}{\partial \dot{\theta}}$ are

$$m\ddot{x} - mg\sin(\phi) = -\lambda \quad \text{and} \quad ma^2\ddot{\theta} = \lambda a,$$
 (1.101)

along with $\dot{x} = a\dot{\theta}$. Taking a time derivative of the constraint gives $\ddot{x} = a\ddot{\theta}$, so $m\ddot{x} = \lambda$, and $\ddot{x} = \frac{g}{2}\sin(\phi)$. This is one-half of the acceleration of a sliding mass. Plugging this back in we find that

$$\lambda = \frac{1}{2} mg \sin(\phi) \tag{1.102}$$

is the friction force in the $-\hat{x}$ direction for the no-sliding constraint, and also $\ddot{\theta} = \frac{g}{2a}\sin(\phi)$.

Example: Consider a wedge of mass m_2 and angle α resting on ice and moving without friction. Let us also consider a mass m_1 sliding without friction on the wedge and try to find the equations of motion and constraint forces. The constraints are that $y_2 = 0$ so the

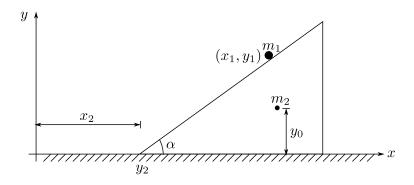


Figure 1.12: Wedge sliding on ice

wedge is always sitting on ice, and $\frac{y_1-y_2}{x_1-x_2} = \tan(\alpha)$ so the point mass is always sitting on the wedge. We will ignore the constraint force for no rotation of the wedge, and aim to find the constraint forces for these two constraints, and the equations for the motion.

The kinetic energy is simply $T = \frac{m_1}{2} (\dot{x}_1^2 + \dot{y}_1^2) + \frac{m_2}{2} (\dot{x}_2^2 + \dot{y}_2^2)$, while the potential energy is $V = m_1 g y_1 + m_2 g (y_2 + y_0)$, where y_0 is the height of the center of mass of the wedge. Then L = T - V, with the constraints $f_1 = (y_1 - y_2) - (x_1 - x_2) \tan(\alpha) = 0$ and $f_2 = y_2 = 0$. The equations of motion from the Euler-Lagrange equations with these holonomic constraints are

$$\frac{d}{dt}\frac{\partial L}{\partial \dot{x}_{1}} - \frac{\partial L}{\partial x_{1}} = \lambda_{1}\frac{\partial f_{1}}{\partial x_{1}} + \lambda_{2}\frac{\partial f_{2}}{\partial x_{1}} \qquad \Longrightarrow \qquad m_{1}\ddot{x}_{1} = -\lambda_{1}\tan(\alpha), \qquad (1.103)$$

$$\frac{d}{dt}\frac{\partial L}{\partial \dot{y}_{1}} - \frac{\partial L}{\partial y_{1}} = \lambda_{1}\frac{\partial f_{1}}{\partial y_{1}} + \lambda_{2}\frac{\partial f_{2}}{\partial y_{1}} \qquad \Longrightarrow \qquad m_{1}\ddot{y}_{1} + m_{1}g = \lambda_{1},$$

$$\frac{d}{dt}\frac{\partial L}{\partial \dot{x}_{2}} - \frac{\partial L}{\partial x_{2}} = \lambda_{1}\frac{\partial f_{1}}{\partial x_{2}} + \lambda_{2}\frac{\partial f_{2}}{\partial x_{2}} \qquad \Longrightarrow \qquad m_{2}\ddot{x}_{2} = \lambda_{1}\tan(\alpha),$$

$$\frac{d}{dt}\frac{\partial L}{\partial \dot{y}_{2}} - \frac{\partial L}{\partial y_{2}} = \lambda_{1}\frac{\partial f_{1}}{\partial y_{2}} + \lambda_{2}\frac{\partial f_{2}}{\partial y_{2}} \qquad \Longrightarrow \qquad m_{2}\ddot{y}_{2} + m_{2}g = -\lambda_{1} + \lambda_{2},$$

which in conjunction with $y_1 - y_2 = (x_1 - x_2) \tan(\alpha)$ and $y_2 = 0$ is six equations for six variables. We number them (1) to (6) and now consider putting the equations together. Equation (6) gives $\ddot{y}_2 = 0$ so (4) gives $m_2 g = \lambda_2 - \lambda_1$ where λ_2 is the force of the ice on the wedge and λ_1 is the *vertical* force (component) of the wedge on the point mass. Adding (1) and (3) gives $m_1\ddot{x}_1 + m_2\ddot{x}_2 = 0$ meaning that the CM of m_1 and m_2 has no overall force acting on it.

Additionally, as (5) implies $\ddot{y}_1 = (\ddot{x}_1 - \ddot{x}_2) \tan(\alpha)$, then using (1), (2), and (3) we find the constant force

$$\lambda_1 = \frac{g}{\frac{1}{m_1 \cos^2(\alpha)} + \frac{\tan^2(\alpha)}{m_2}}.$$
 (1.104)

With this result in hand we can use it in (1), (2), and (3) to solve for the trajectories. Since

$$\ddot{x}_2 = \frac{\tan(\alpha)}{m_2} \lambda_1,$$

$$\ddot{x}_1 = -\frac{\tan(\alpha)}{m_1} \lambda_1,$$

$$\ddot{y}_1 = \frac{\lambda_1}{m_1} - g,$$
(1.105)

the accelerations are constant. As a check on our results, if $m_2 \to \infty$, then $\ddot{x}_2 = 0$ so indeed the wedge is fixed; and for this case, $\ddot{x}_1 = -g\sin(\alpha)\cos(\alpha)$ and $\ddot{y}_1 = -g\sin^2(\alpha)$ which both vanish as $\alpha \to 0$ as expected (since in that limit the wedge disappears, flattening onto the icy floor below it).

Chapter 2

Rigid Body Dynamics

2.1 Coordinates of a Rigid Body

A set of N particles forms a *rigid body* if the distance between any 2 particles is fixed:

$$r_{ij} \equiv |\mathbf{r}_i - \mathbf{r}_j| = c_{ij} = \text{constant.}$$
 (2.1)

Given these constraints, how many generalized coordinates are there?

If we know 3 non-collinear points in the body, the remaing points are fully determined by triangulation. The first point has 3 coordinates for translation in 3 dimensions. The second point has 2 coordinates for spherical rotation about the first point, as r_{12} is fixed. The third point has one coordinate for circular rotation about the axis of \mathbf{r}_{12} , as r_{13} and r_{23} are fixed. Hence, there are 6 independent coordinates, as represented in Fig. 2.1. The result that there are six continuous coordinates is independent of N, so it also applies to a continuous body (in the limit of $N \to \infty$).

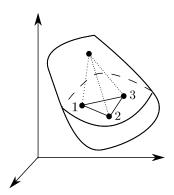


Figure 2.1: Three non-collinear points can be fully determined by using 6 continuous coordinates. Since the distances between any two other points are fixed in the rigid body, any other point of the body is fully determined by the distance to these 3 points (and a discrete orientation which is not a dynamical variable).

The translations of the body require three spatial coordinates. These translations can be taken from any fixed point in the body. Typically the fixed point is the center of mass (CM), defined as:

$$\mathbf{R} = \frac{1}{M} \sum_{i} m_i \mathbf{r}_i, \tag{2.2}$$

where m_i is the mass of the *i*-th particle and \mathbf{r}_i the position of that particle with respect to a fixed origin and set of axes (which will notationally be unprimed) as in Fig. 2.2. In the case of a continuous body, this definition generalizes as:

$$\mathbf{R} = \frac{1}{M} \int_{\mathcal{V}} \mathbf{r} \rho(\mathbf{r}) \, d\mathcal{V},\tag{2.3}$$

where $\rho(\mathbf{r})$ is the mass density at position \mathbf{r} and we integrate over the volume \mathcal{V} .

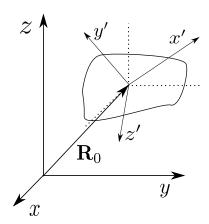


Figure 2.2: The three translational coordinates correspond to the position of the Center of Mass, and the three rotational coordinates correspond to the three angles necessary to define the orientation of the axis fixed with the body.

Rotations of the body are taken by fixing axes with respect to the body (we will denote these body fixed axes with primes) and describing their orientation with respect to the unprimed axes by 3 angles (ϕ, θ, ψ) .

A particularly useful choice of angles are called *Euler angles*. The angle ϕ is taken as a rotation about the z-axis, forming new \tilde{x} - and \tilde{y} -axes while leaving the z-axis unchanged, as shown in Fig. 2.3. The angle θ is then taken as a rotation about the \tilde{x} -axis, forming new \tilde{y}' - and z'-axes while leaving the \tilde{x} -axis unchanged, as shown in Fig. 2.4. Finally, the angle ψ is taken as a rotation about the z'-axis, forming new x'- and y'-axes while leaving the z'-axis unchanged, as shown in Fig. 2.5. (The \tilde{x} -axis is called the line of nodes, as it is the intersection of the xy- and x'y'-planes.)

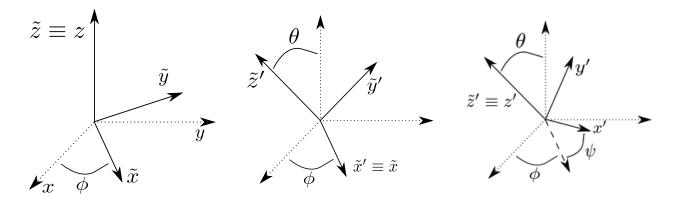


Figure 2.3: First rotation is by ϕ around the original z axis.

Figure 2.4: Second rotation is by θ around the intermediate \tilde{x} axis.

Figure 2.5: Final rotation is by ψ around the final z' axis.

Rotations can be described by 3×3 matrices U. This means each rotation step can be described as a matrix multiplication. Where $\mathbf{r} = (x, y, z)$, then

$$\tilde{\mathbf{r}} = U_{\phi}\mathbf{r} = \begin{bmatrix} \cos(\phi) & \sin(\phi) & 0 \\ -\sin(\phi) & \cos(\phi) & 0 \\ 0 & 0 & 1 \end{bmatrix} \begin{bmatrix} x \\ y \\ z \end{bmatrix}. \tag{2.4}$$

Similar transformations can be written for the other terms:

$$\tilde{\mathbf{r}}' = U_{\theta}\tilde{\mathbf{r}}$$
 , $\mathbf{r}' = U_{\psi}\tilde{\mathbf{r}}' = U_{\psi}U_{\theta}\tilde{\mathbf{r}} = U_{\psi}U_{\theta}U_{\phi}\mathbf{r}$.

Defining the total transformation as U, it can be written as:

$$U \equiv U_{\psi} U_{\theta} U_{\phi} \quad \Rightarrow \quad \mathbf{r}' = U \mathbf{r}.$$
 (2.5)

Care is required with the order of the terms since the matrices don't commute. Writing U out explicitly:

$$U = \begin{bmatrix} \cos(\psi) & \sin(\psi) & 0 \\ -\sin(\psi) & \cos(\psi) & 0 \\ 0 & 0 & 1 \end{bmatrix} \begin{bmatrix} 1 & 0 & 0 \\ 0 & \cos(\theta) & \sin(\theta) \\ 0 & -\sin(\theta) & \cos(\theta) \end{bmatrix} \begin{bmatrix} \cos(\phi) & \sin(\phi) & 0 \\ -\sin(\phi) & \cos(\phi) & 0 \\ 0 & 0 & 1 \end{bmatrix}.$$
(2.6)

All rotation matrices, including U_{ϕ} , U_{θ} , U_{ψ} , and U are orthogonal. Orthogonal matrices W satisfy

$$W^{\top}W = WW^{\top} = \mathbb{1} \quad \Leftrightarrow \quad W^{\top} = W^{-1}, \tag{2.7}$$

where $\mathbb{1}$ refers to the identity matrix and \top to the transpose. This ensures that the length of a vector is invariant under rotations:

$$\mathbf{r}^{\prime 2} = \mathbf{r}^{\top} (W^{\top} W) \mathbf{r} = \mathbf{r}^{2}. \tag{2.8}$$

Orthogonal matrices W have 9 entries but need to fulfill 6 conditions from orthogonality, leaving only 3 free parameters, corresponding to the 3 angles necessary to determine the rotation.

We can also view $\mathbf{r}' = U\mathbf{r}$ as a transformation from the vector \mathbf{r} to the vector \mathbf{r}' in the same coordinate system. This is an active transformation, as opposed to the previous perspective which was a passive transformation.

Finally, note that inversions like

$$U = \begin{bmatrix} -1 & 0 & 0 \\ 0 & -1 & 0 \\ 0 & 0 & -1 \end{bmatrix} \quad \text{or} \quad U = \begin{bmatrix} 1 & 0 & 0 \\ 0 & 1 & 0 \\ 0 & 0 & -1 \end{bmatrix}$$
 (2.9)

are not rotations. These have $\det(U) = -1$, so they can be forbidden by demanding that $\det(U) = 1$. All orthogonal matrices have $\det(W) = \pm 1$ because $\det(W^{\top}W) = (\det(W))^2 = 1$. In the language of group theory, the restriction to $\det(W) = 1$ gives the special orthogonal group SO(3) as opposed to simply O(3), the orthogonal group. We disregard the $\det(U) = -1$ subset of transformations because it is impossible for the system to undergo these transformations without the distance between the particles jumping abruptly in the process (they are not continuously connected to the identity transformation), so they are not allowed mechanical motions of our rigid body.

Intuitively, we could rotate the coordinates (x, y, z) directly into the coordinates (x', y', z') by picking the right axis of rotation. In fact, the *Euler theorem* states that a general displacement of a rigid body with one point fixed is equivalent to a single rotation about *some* axis that runs through the fixed point. This theorem will be true if a general rotation U leaves some axis fixed, which is satisfied by

$$U\mathbf{r} = \mathbf{r} \tag{2.10}$$

for any point \mathbf{r} on this axis. This is an eigenvalue equation for U with eigenvalue 1. To better understand this, we need to use the linear algebra concepts of eigenvalue equations, which we now review.

Although the notion of an eigenvalue equation generally holds for linear operators, for now the discussion will be restricted to orthogonal rotation matrices U. The eigenvalue equation is

$$U\boldsymbol{\xi} = \lambda \boldsymbol{\xi},\tag{2.11}$$

where ξ is an eigenvector and λ is the associated eigenvalue. Rewriting this as

$$(U - \lambda \mathbb{1})\boldsymbol{\xi} = 0 \tag{2.12}$$

requires that $\det(U - \lambda \mathbb{1}) = 0$, so that $U - \lambda \mathbb{1}$ is not invertible and the solution can be non-trivial, $\xi \neq 0$. $\det(U - \lambda \mathbb{1}) = 0$ is a cubic equation in λ , which has 3 solutions, which are the eigenvalues λ_{α} for $\alpha \in \{1, 2, 3\}$. The associated eigenvectors are $\xi^{(\alpha)}$ and satisfy

$$U\boldsymbol{\xi}^{(\alpha)} = \lambda_{\alpha}\boldsymbol{\xi}^{(\alpha)},\tag{2.13}$$

where no implicit sum over repeated indices is taken. Forming a matrix from the resulting eigenvectors as columns:

$$X = \begin{bmatrix} \uparrow & \uparrow & \uparrow \\ \boldsymbol{\xi}^{(1)} & \boldsymbol{\xi}^{(2)} & \boldsymbol{\xi}^{(3)} \\ \downarrow & \downarrow & \downarrow \end{bmatrix}$$
 (2.14)

then we can rewrite Eq.(2.13) as

$$UX = X \cdot \operatorname{diag}(\lambda_1, \lambda_2, \lambda_3) \Rightarrow X^{-1}UX = \operatorname{diag}(\lambda_1, \lambda_2, \lambda_3)$$
 (2.15)

This means X diagonalizes U. Since U is orthogonal, the matrix X is unitary ($X^{\dagger}X = XX^{\dagger} = 1$). Note that $^{\top}$ indicates transposition whereas † indicates Hermitian conjugation (i.e. complex conjugation * combined with transposition $^{\top}$).

Next we note that since $\det(U) = 1$, then $\lambda_1 \lambda_2 \lambda_3 = 1$. Additionally, $|\lambda_{\alpha}|^2 = 1$ for each α because:

$$U\boldsymbol{\xi} = \lambda \boldsymbol{\xi} \quad \Rightarrow \quad \boldsymbol{\xi}^{\dagger} U^{\top} = \lambda^{\star} \boldsymbol{\xi}^{\dagger} \quad \Rightarrow \quad \boldsymbol{\xi}^{\dagger} \boldsymbol{\xi} = \boldsymbol{\xi}^{\dagger} U^{\top} U \boldsymbol{\xi} = |\lambda|^{2} \boldsymbol{\xi}^{\dagger} \boldsymbol{\xi} \,. \tag{2.16}$$

Finally, if λ is an eigenvalue, then so is λ^* , since we can take the complex conjugate of the original equation:

$$U\boldsymbol{\xi} = \lambda \boldsymbol{\xi} \Rightarrow U\boldsymbol{\xi}^* = \lambda^* \boldsymbol{\xi}^* \tag{2.17}$$

where $\boldsymbol{\xi}^{\star}$ is still a column vector but with its elements undergoing complex conjugation with respect to $\boldsymbol{\xi}$. Without loss of generality, let us say for a rotation matrix U that $\lambda_2 = \lambda_3^{\star}$. Then $1 = \lambda_1 |\lambda_2|^2 = \lambda_1$, so one of the eigenvalues is 1, giving Eq.(2.10), and thus proving Euler's Theorem. The associated eigenvector $\boldsymbol{\xi}^{(1)}$ to the eigenvalue $\lambda_1 = 1$ is the rotation axis, and if $\lambda_2 = \lambda_3^{\star} = e^{i\Phi}$ then Φ is the rotation angle about that axis.

In fact, we can make good use of our analysis of Euler's theorem. Together the rotation axis and rotation angle can be used to define the instantaneous angular velocity $\omega(t)$ such that:

$$|\boldsymbol{\omega}| = \frac{d\Phi}{dt}$$
 and $\boldsymbol{\omega} \parallel \boldsymbol{\xi}^{(1)}$. (2.18)

The angular velocity will play an important role in our discussion of time dependence with rotating coordinates in the next section. If we consider several consecutive displacements of the rigid body, then each can have its own axis $\boldsymbol{\xi}^{(1)}$ and its own $\dot{\Phi}$, so $\boldsymbol{\omega}$ changes at each instance of time, and hence $\boldsymbol{\omega} = \boldsymbol{\omega}(t)$ (for the entire rigid body).

2.2 Time Evolution with Rotating Coordinates

Lets use unprimed axes (x, y, z) for the fixed (inertial) axes, with fixed basis vectors \mathbf{e}_i . We will also use primed axes (x', y', z') for the body axes with basis vectors \mathbf{e}'_i .

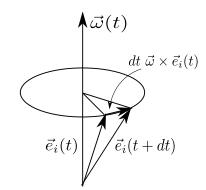
If we consider *any* vector then it can be decomposed with either set of basis vectors:

$$\mathbf{b} = \sum_{i} b_{i} \mathbf{e}_{i} = \sum_{i} b'_{i} \mathbf{e}'_{i}. \tag{2.19}$$

For fixed axes basis vectors by definition $\dot{\mathbf{e}}_i = 0$, while for those in the body frame,

$$\dot{\mathbf{e}}_i' = \boldsymbol{\omega}(t) \times \mathbf{e}_i' \tag{2.20}$$

meaning vectors of fixed length undergo a rotation at a time t. The derivation of this result is shown in the figure on the right, by considering the change to the vector after an infinitesimal time interval dt.



Summing over repeated indices, this means:

$$\dot{\mathbf{b}} = \dot{b}_i \mathbf{e}_i = \dot{b}_i' \mathbf{e}_i' + \boldsymbol{\omega}(t) \times (b_i' \mathbf{e}_i') = \dot{b}_i' \mathbf{e}_i' + \boldsymbol{\omega}(t) \times \mathbf{b}$$

Defining $\frac{d}{dt}$ as the time evolution in the fixed (F) frame and $\frac{d_{\rm R}}{dt}$ the time evolution in the rotating/body (R) frame, then vectors evolve in time according to

$$\frac{d\mathbf{b}}{dt} = \frac{d_{\mathbf{R}}\mathbf{b}}{dt} + \boldsymbol{\omega} \times \mathbf{b}.$$
 (2.21)

As a mnemonic we have the operation " $(d/dt) = d_R/dt + \omega \times$ " which can act on any vector. Let us apply this to the position **r** of a particle of mass m, which gives

$$\frac{d\mathbf{r}}{dt} = \frac{d_{R}\mathbf{r}}{dt} + \boldsymbol{\omega} \times \mathbf{r} \quad \Leftrightarrow \quad \mathbf{v}_{F} = \mathbf{v}_{R} + \boldsymbol{\omega} \times \mathbf{r}. \tag{2.22}$$

Taking another time derivative gives us the analog for acceleration:

$$\frac{\mathbf{F}}{m} = \frac{d\mathbf{v}_{F}}{dt} = \frac{d_{R}\mathbf{v}_{F}}{dt} + \boldsymbol{\omega} \times \mathbf{v}_{F}$$

$$= \frac{d_{R}\mathbf{v}_{R}}{dt} + \frac{d_{R}\boldsymbol{\omega}}{dt} \times \mathbf{r} + \boldsymbol{\omega} \times \frac{d_{R}\mathbf{r}}{dt} + \boldsymbol{\omega} \times \mathbf{v}_{R} + \boldsymbol{\omega} \times (\boldsymbol{\omega} \times \mathbf{r}).$$
(2.23)

As $\frac{d_{\rm R}\mathbf{r}}{dt} = \mathbf{v}_{\rm R}$ is the velocity within the body frame and $\frac{d_{\rm R}\mathbf{v}_{\rm R}}{dt} = \mathbf{a}_{\rm R}$ is the acceleration within the body frame, and $d_{\rm R}\boldsymbol{\omega}/dt = d\boldsymbol{\omega}/dt = \dot{\boldsymbol{\omega}}$ then

$$m\mathbf{a}_{\mathrm{R}} = \mathbf{F} - m\boldsymbol{\omega} \times (\boldsymbol{\omega} \times \mathbf{r}) - 2m\boldsymbol{\omega} \times \mathbf{v}_{\mathrm{R}} - m\dot{\boldsymbol{\omega}} \times \mathbf{r}$$
 (2.24)

gives the acceleration in the body frame with respect to the forces that seem to be present in that frame. The terms $-m\boldsymbol{\omega}\times(\boldsymbol{\omega}\times\mathbf{r})$ and $-2m\boldsymbol{\omega}\times\mathbf{v}_{\mathrm{R}}$ are, respectively, the centrifugal

and Coriolis ficticious forces respectively, while the last term $-m\frac{d_{\mathbf{R}}\boldsymbol{\omega}}{dt}\times\mathbf{r}$ is a ficticious force that arises from non-uniform rotational motion, so that there is angular acceleration within the body frame. The same result could also have been obtained with the Euler-Lagrange equations for L in the rotating coordinates:

$$L = \frac{m}{2}(\dot{\mathbf{r}} + \boldsymbol{\omega} \times \mathbf{r})^2 - V, \qquad (2.25)$$

and you will explore this on a problem set.

Note that the centrifugal term is radially outward and perpendicular to the rotation axis. To see this, decompose \mathbf{r} into components parallel and perpendicular to $\boldsymbol{\omega}$, $\mathbf{r} = \mathbf{r}_{\parallel} + \mathbf{r}_{\perp}$, then $\boldsymbol{\omega} \times \boldsymbol{r} = \boldsymbol{\omega} \times \boldsymbol{r}_{\perp}$, so $-\boldsymbol{\omega} \times (\boldsymbol{\omega} \times \boldsymbol{r}_{\perp}) = w^2 \boldsymbol{r}_{\perp}$. This term is present for any rotating body. On the other hand, the Coriolis force is nonzero when there is a nonzero velocity in the rotating/body frame: $\mathbf{v}_{\mathrm{R}} \neq 0$.

Example: Consider the impact of the Coriolis force on projectile motion on the rotating Earth, where the angular velocity is $\omega_{\text{Earth}} = \frac{2\pi}{24 \times 3600 \, \text{s}} \approx 7.3 \times 10^{-5} \, \text{s}^{-1}$. We work out the cross-product $-\boldsymbol{\omega} \times \boldsymbol{v}_r$ as shown in Fig. 2.6 for a particle in the northern hemisphere, where ω points to the north pole. Thus a projectile in the northern/southern hemisphere would be perturbed to the right/left relative to its velocity direction \boldsymbol{v}_r .

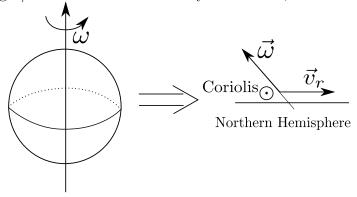
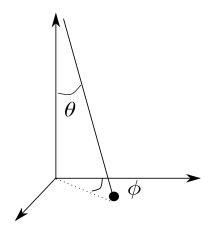
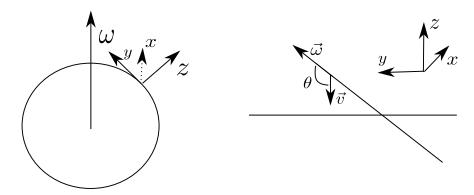


Figure 2.6: For a projectile, in the Northern Hemisphere, the Coriolis pushes it to its right, relative to its direction of motion.

Exercise: Consider a Foucault pendulum which hangs from a rigid rod, but is free to oscillate in two angular directions, as shown in Fig. 2.2. For $\theta \ll 1$ and working to first order in the small ω , the result derived from the Coriolis force gives $\dot{\phi} \approx \omega_{\text{Earth}} \sin(\lambda)$. Here λ is the latitude angle measured from equator. The precession is clockwise in the northern hemisphere, and is maximized at the north pole where $\lambda = 90^{\circ}$.



Example: Consider the Coriolis deflection of a freely falling body on Earth in the northern hemisphere. We use the coordinate system shown below, where z is perpendicular to the surface of the earth and y is parallel to the earth's surface and points towards the north pole.



Working to first order in the small ω gives us

$$m\mathbf{a}_{\mathrm{R}} = m\dot{\mathbf{v}}_{\mathrm{R}} = -mg\hat{z} - 2m\boldsymbol{\omega} \times \mathbf{v},$$
 (2.26)

where the centrifugal terms of order $O(\omega^2)$ are dropped. As an initial condition we take $\mathbf{v}(t=0) = v_0 \hat{z}$. The term $-\boldsymbol{\omega} \times \mathbf{v}$ points along \hat{x} , so:

$$\ddot{z} = -g + O(\omega^2) \quad \Rightarrow \quad v_z = v_0 - gt \tag{2.27}$$

Moreover implementing the boundary condition that $\dot{x}(t=0)=0$:

$$\ddot{x} = -2(\boldsymbol{\omega} \times \mathbf{v})_x = -2\omega \sin(\theta) v_z(t) \quad \Rightarrow \quad \dot{x} = -2\omega \sin(\theta) \left(v_0 t - \frac{g}{2} t^2 \right). \tag{2.28}$$

Taking also x(t=0)=0, and integrating one further time, the motion in the x direction is:

$$x(t) = -2\omega \sin(\theta) \left(\frac{v_0}{2}t^2 - \frac{g}{6}t^3\right). \tag{2.29}$$

Lets consider this effect for a couple simple cases. If the mass m is dropped from a height $z(t=0) = h_{\text{max}}$ with zero velocity, $v_0 = 0$, then:

$$z = h_{\text{max}} - \frac{g}{2}t^2 \tag{2.30}$$

and the mass reaches the floor at time

$$t_1 = \sqrt{\frac{2h_{\text{max}}}{g}}. (2.31)$$

From Eq.(2.28) we see that $\dot{x} > 0$ for all t, and that:

$$x(t = t_1) = \frac{2\sqrt{2}\,\omega\sin(\theta)h_{\text{max}}^{3/2}}{3\sqrt{g}} > 0.$$
 (2.32)

However, if the mass m is thrown up with an initial $\dot{z}(t=0)=v_0>0$ from the ground (z=0), then:

$$z(t) = v_0 t - \frac{g}{2} t^2 > 0. (2.33)$$

Here the particle rises to a maximum height $z = v_0^2/(2g)$ at time $t = v_0/g$, and then falls back to earth. Using Eq.(2.28) we see that $\dot{x} < 0$ for all t. If t_1 is the time it reaches the ground again $(t_1 = \frac{2v_0}{g})$, then:

$$x(t=t_1) = -\frac{4\omega \sin(\theta)v_0^3}{3g^2} < 0.$$
 (2.34)

2.3 Kinetic Energy, Angular Momentum, and the Moment of Inertia Tensor for Rigid Bodies

Returning to rigid bodies, consider one built out of N fixed particles The kinetic energy is best expressed using CM coordinates, where \mathbf{R} is the CM and we here take \mathbf{r}_i to be the *displacement* of particle i relative to the CM. Once again making sums over repeated subscripts as implicit, the kinetic energy (T) of the system is given by:

$$T = \frac{1}{2}M\dot{\mathbf{R}}^2 + \frac{1}{2}m_i\dot{\mathbf{r}}_i^2. \tag{2.35}$$

As the body is rigid, then points cannot translate relative to the body but can only rotate so that $\dot{\mathbf{r}}_i = \boldsymbol{\omega} \times \mathbf{r}_i$. The rotational kinetic energy is then

$$T_{\mathrm{R}} = \frac{1}{2} m_i \dot{\mathbf{r}}_i^2 = \frac{1}{2} m_i (\boldsymbol{\omega} \times \mathbf{r}_i)^2 = \frac{1}{2} m_i [\boldsymbol{\omega}^2 \mathbf{r}_i^2 - (\boldsymbol{\omega} \cdot \mathbf{r}_i)^2]. \tag{2.36}$$

Labeling Cartesian indices with a and b to reserve i and j for particle indices, then we can write out this result making the indices all explict as

$$T_{\rm R} = \frac{1}{2} \sum_{i,a,b} m_i (\delta_{ab} \mathbf{r}_i^2 - r_{ia} r_{ib}) \omega_a \omega_b.$$
 (2.37)

It is convenient to separate out the parts in this formula that depend on the shape and distributions of masses in the body by defining the moment of inertia tensor \hat{I} for the discrete body as

$$\hat{I}_{ab} \equiv \sum_{i} m_i (\delta_{ab} \mathbf{r}_i^2 - r_{ia} r_{ib}). \tag{2.38}$$

The analog for a continuous body of mass density $\rho(\mathbf{r})$ is:

$$\hat{I}_{ab} \equiv \int_{\mathcal{V}} (\mathbf{r}^2 \delta_{ab} - r_a r_b) \rho(\mathbf{r}) \, d\mathcal{V}. \tag{2.39}$$

In terms of the moment of inertia tensor, the kinetic energy from rotation can now be written as:

$$T_{\rm R} = \frac{1}{2} \sum_{ab} \hat{I}_{ab} \omega_a \omega_b = \frac{1}{2} \boldsymbol{\omega} \cdot \hat{I} \cdot \boldsymbol{\omega} , \qquad (2.40)$$

where in the last step we adopt a convenient matrix multiplication notation.

The moment of inertia tensor can be written with its components as a matrix in the form

$$\hat{I} = \sum_{i} m_{i} \begin{bmatrix} y_{i}^{2} + z_{i}^{2} & -x_{i}y_{i} & -x_{i}z_{i} \\ -x_{i}y_{i} & x_{i}^{2} + z_{i}^{2} & -y_{i}z_{i} \\ -x_{i}z_{i} & -y_{i}z_{i} & x_{i}^{2} + y_{i}^{2} \end{bmatrix},$$
(2.41)

where the diagonal terms are the "moments of inertia" and the off-diagonal terms are the "products of inertia". Note also that \hat{I} is symmetric in any basis, so $\hat{I}_{ab} = \hat{I}_{ba}$.

Special case: if the rotation happens about only one axis which can be defined as the z-axis for convenience so that $\boldsymbol{\omega} = (0,0,\omega)$, then $T_{\rm R} = \frac{1}{2}\hat{I}_{zz}\omega^2$ which reproduces the simpler and more familiar scalar form of the moment of inertia.

Lets now let \mathbf{r}_i be measured from a stationary point in the rigid body, which need not necessarily be the CM. The angular momentum can be calculated about this fixed point. Since $\mathbf{v}_i = \boldsymbol{\omega} \times \mathbf{r}_i$, we can write the angular momentum as:

$$\mathbf{L} = m_i \mathbf{r}_i \times \mathbf{v}_i = m_i \mathbf{r}_i \times (\boldsymbol{\omega} \times \mathbf{r}_i) = m_i [\mathbf{r}_i^2 \boldsymbol{\omega} - (\boldsymbol{\omega} \cdot \mathbf{r}_i) \mathbf{r}_i]. \tag{2.42}$$

Writing out the components

$$L_a = \sum_i m_i (\mathbf{r}_i^2 \omega_a - (\boldsymbol{\omega} \cdot \mathbf{r}_i) r_{ia}) = \sum_{i,b} \omega_b m_i (\delta_{ab} \mathbf{r}_i^2 - r_{ia} r_{ib}) = \sum_b \hat{I}_{ab} \omega_b, \tag{2.43}$$

which translates to the matrix equation:

$$\mathbf{L} = \hat{I} \cdot \boldsymbol{\omega} \,. \tag{2.44}$$

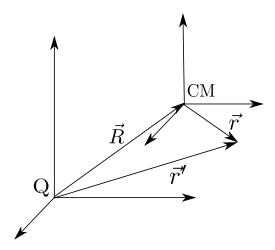
This allows us to write the corresponding rotational kinetic energy about the fixed point as:

$$T_{\rm R} = \frac{1}{2} \boldsymbol{\omega} \cdot \mathbf{L}. \tag{2.45}$$

Note that in general, **L** is *not* parallel to ω . We will see an explicit example of this below. Also note that the formula that we used for \hat{I} in this case is the same as we gave above. We use these formulas whether or not \mathbf{r}_i is taken with respect to the CM. (Do note that if the body is translating, such that no point in the body is fixed, then the CM point has a special role since the translational and rotational kinetic energy can be separated when using this point, and not in general other points.)

It is useful to pause to see what precisely the calculation of \hat{I} depends on. Since it involves components of the vectors r_i it depends on the choice of the origin for the rotation. Furthermore the entries of the matrix \hat{I}_{ab} obviously depend on the orientation of the axes used to define the components labeled by a and b. Given this, it is natural to ask whether given the result for \hat{I}_{ab} with one choice of axes and orientation, whether we can determine an $\hat{I}'_{a'b'}$ for a different origin and axes orientation. This is always possible with the help of a couple of theorems.

The parallel axis theorem: Given \hat{I}^{CM} about the CM, it is simple to find \hat{I}^{Q} about a different point Q with the *same* orientation for the axes. Referring to the figure below,



we define \mathbf{r}'_i as the coordinate of a particle i in the rigid body with respect to point Q and \mathbf{r}_i to be the coordinate of that particle with respect to the CM, so that:

$$\mathbf{r}_i' = \mathbf{R} + \mathbf{r}_i. \tag{2.46}$$

By definition of the CM:

$$\sum_{i} m_{i} \mathbf{r}_{i} = 0 \qquad \text{and we let} \qquad M = \sum_{i} m_{i}. \tag{2.47}$$

The tensor of inertia around the new axis is then:

$$\hat{I}_{ab}^{Q} = m_i (\delta_{ab} \mathbf{r}_i'^2 - r_{ia}' r_{ib}')$$

$$= m_i \left(\delta_{ab} (\mathbf{r}_i^2 + 2\mathbf{r}_i \cdot \mathbf{R} + \mathbf{R}^2) - \mathbf{r}_{ia} \mathbf{r}_{ib} - \mathbf{r}_{ia} \mathbf{R}_b - \mathbf{R}_a \mathbf{r}_{ib} - \mathbf{R}_a \mathbf{R}_b \right), \qquad (2.48)$$

where the cross terms involving a single \mathbf{r}_i or single component \mathbf{r}_{ia} sum up to zero by Eq.(2.47). The terms quadratic in \mathbf{r} are recognized as giving the moment of inertia tensor about the CM. This gives the *parallel axis theorem* for translating the origin:

$$\hat{I}_{ab}^{Q} = M(\delta_{ab}\mathbf{R}^2 - \mathbf{R}_a\mathbf{R}_b) + \hat{I}_{ab}^{CM}, \tag{2.49}$$

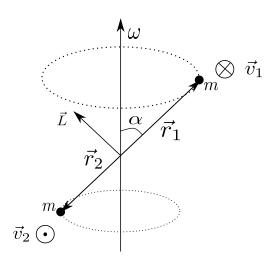
If we wish to carry out a translation between P and Q, neither of which is the CM, then we can simply use this formula twice. Another formula can be obtained by projecting the parallel axis onto a specific axis \hat{n} where $\hat{n}^2 = 1$ (giving a result that may be familiar from an earlier classical mechanics course):

$$\hat{n} \cdot \hat{I}^{Q} \cdot \hat{n} = M(\mathbf{R}^{2} - (\hat{n} \cdot \mathbf{R})^{2}) + \hat{n} \cdot \hat{I}^{CM} \cdot \hat{n} = M\mathbf{R}^{2}[1 - \cos^{2}(\theta)] + \hat{n} \cdot \hat{I}^{CM} \cdot \hat{n}$$

$$= M\mathbf{R}^{2} \sin^{2}(\theta) + \hat{n} \cdot \hat{I}^{CM} \cdot \hat{n}$$
(2.50)

where $\hat{n} \cdot \mathbf{R} \equiv R \cos(\theta)$.

Example: Lets consider an example of the calculation of \hat{I} for a situation where \boldsymbol{L} is not parallel to $\boldsymbol{\omega}$. Consider a dumbbell made of 2 identical point passes m attached by a massless rigid rod (but with different separations r_1 and r_2 from the axis of rotation), spinning so that $\boldsymbol{\omega} = \omega \hat{z}$ and so that the rod makes an angle α with the axis of rotation, as shown



We define body axes where the masses lie in the yz-plane. Here,

$$\mathbf{r}_1 = (0, r_1 \sin \alpha, r_1 \cos \alpha) \text{ and } \mathbf{r}_2 = (0, -r_2 \sin \alpha, -r_2 \cos \alpha). \tag{2.51}$$

Then using the definition of the moment inertia tensor:

$$I_{zz} = m(x_1^2 + y_1^2) + m(x_2^2 + y_2^2) = m(r_1^2 + r_2^2) \sin^2 \alpha$$

$$I_{xx} = m(y_1^2 + z_1^2) + m(y_2^2 + z_2^2) = m(r_1^2 + r_2^2)$$

$$I_{yy} = m(x_1^2 + z_1^2) + m(x_2^2 + z_2^2) = m(r_1^2 + r_2^2) \cos^2 \alpha$$

$$I_{xy} = I_{yx} = m(-x_1y_1 - x_2y_2) = 0$$

$$I_{xz} = I_{zx} = m(-x_1z_1 - x_2z_2) = 0$$

$$I_{yz} = I_{zy} = m(-y_1z_1 - y_2z_2) = -m(r_1^2 + r_2^2) \sin \alpha \cos \alpha$$
(2.52)

Plugging these into $\mathbf{L} = \hat{I} \cdot \boldsymbol{\omega}$, recalling that only ω_z is non-zero, this gives

$$\mathbf{L} = (0, I_{yz}\omega, I_{zz}\omega). \tag{2.53}$$

Thus in this example L is not parallel to ω .

Next, instead of translating the axes in a parallel manner, let us keep the origin fixed and rotate the axes according to an orthogonal rotation matrix U satisfying $U^{\top}U = UU^{\top} = 1$. Vectors are rotated as

$$\mathbf{L}' = U\mathbf{L}$$
 , $\boldsymbol{\omega}' = U\boldsymbol{\omega}$ and therefore $\boldsymbol{\omega} = U^{\top}\boldsymbol{\omega}'$. (2.54)

Putting these together

$$\mathbf{L}' = U\hat{I} \cdot \boldsymbol{\omega} = (U\hat{I}U^{\top}) \cdot \boldsymbol{\omega}' \quad \Rightarrow \quad \hat{I}' = U\hat{I}U^{\top}, \tag{2.55}$$

where \hat{I}' is the new moment of inertia tensor. (The fact that it transforms this way defines it as a tensor.) This allows us to calculate the new moment of inertia tensor after a rotation.

For a real symmetric tensor \hat{I} , there always exists a rotation from an orthogonal matrix U that diagonalizes \hat{I} giving a diagonal matrix \hat{I}' :

$$\hat{I}_D = \begin{pmatrix} I_1 & 0 & 0 \\ 0 & I_2 & 0 \\ 0 & 0 & I_3 \end{pmatrix} . \tag{2.56}$$

The entries of the diagonal moment of inertia tensor, I_{α} , are real and positive. This is just a special case of saying a Hermitian matrix can always be diagonalized by a unitary transformation (which is often derived in a Quantum Mechanics course as part of showing that a Hermitian matrix has real eigenvalues and orthogonal eigenvectors). The positivity

of diagonal matrix follows immediately from the definition of the moment of inertia tensor for the situation with zero off-diagonal terms.

The axes that make \hat{I} diagonal are called the *principal axes* and the components I_{α} are the *principal moments of inertia*. We find them by solving the eigenvalue problem

$$\hat{I} \cdot \boldsymbol{\xi} = \lambda \, \boldsymbol{\xi},\tag{2.57}$$

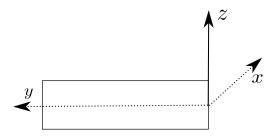
where the 3 eigenvalues λ give the principal moments of inertia I_{α} , and are obtained from solving $\det(\hat{I} - \lambda \mathbb{1}) = 0$. The corresponding 3 real *orthogonal* eigenvectors $\boldsymbol{\xi}^{(\alpha)}$ are the principal axes. Here $U^{\top} = \begin{bmatrix} \boldsymbol{\xi}^{(1)} & \boldsymbol{\xi}^{(2)} & \boldsymbol{\xi}^{(3)} \end{bmatrix}$, where the eigenvectors vectors fill out the columns. Then, without summing over repeated indices:

$$L_{\alpha} = I_{\alpha}\omega_{\alpha}$$
 and $T = \frac{1}{2}\sum_{\alpha}I_{\alpha}\omega_{\alpha}^{2}$, (2.58)

where L_{α} and ω_{α} are the components of **L** and $\boldsymbol{\omega}$, respectively, evaluated along the principal axes.

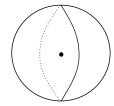
To summarize, for any choice of origin for any rigid body, there is a choice of axes that diagonalizes \hat{I} . For T to separate into translational and rotational parts, we must pick the origin to be the CM. Often, the principal axes can be identified by a symmetry of the body.

Example: for a thin rectangle lying in the yz-plane with one edge coinciding with the z-axis, and the origin chosen as shown below, then $I_{yz} = 0$ as the body is symmetric under $z \leftrightarrow -z$, while $I_{xz} = I_{xy} = 0$ as the body lies entirely within x = 0. Hence these are principal axes.

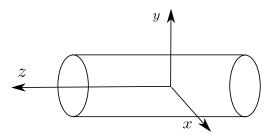


Sometimes, symmetry allows multiple choices for the principal axes.

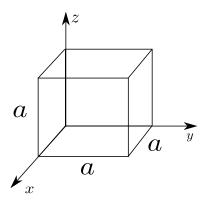
Example: for a sphere, any orthogonal axes through the origin are principal axes.



Example: for a cylinder whose central axis is aligned along the z-axis, because of rotational symmetry any choice of the x- and y-axes gives principal axes.



Example: Lets consider an example where the principal axes may not be apparent, which we can solve through the eigenvalue problem. Consider a uniform cube with sides of length a, mass m, and having the origin at one corner, as shown below.



By symmetry we have

$$I_{xx} = I_{yy} = I_{zz} = \frac{m}{a^3} \int_0^a \int_0^a \int_0^a (x^2 + y^2) \, dx \, dy \, dz = \frac{2}{3} m a^2,$$

$$I_{xy} = I_{yz} = I_{xz} = \frac{m}{a^3} \int_0^a \int_0^a \int_0^a -xz \, dx \, dy \, dz = -\frac{1}{4} m a^2.$$
(2.59)

Thus the matrix is

$$\hat{I} = ma^2 \begin{pmatrix} +\frac{2}{3} & -\frac{1}{4} & -\frac{1}{4} \\ -\frac{1}{4} & +\frac{2}{3} & -\frac{1}{4} \\ -\frac{1}{4} & -\frac{1}{4} & +\frac{2}{3} \end{pmatrix} . \tag{2.60}$$

The principal moments of inertia are found from

$$\det(\hat{I} - \lambda \mathbb{1}) = \left(\frac{11}{12}ma^2 - \lambda\right)^2 \left(\frac{1}{6}ma^2 - \lambda\right) = 0. \tag{2.61}$$

This gives $I_1 = \lambda_1 = \frac{1}{6}ma^2$. Solving

$$(\hat{I} - \lambda_1 \mathbb{1}) \boldsymbol{\xi}^{(1)} = 0$$
 we find $\boldsymbol{\xi}^{(1)} = (1, 1, 1)$. (2.62)

The remaining eigenvalues are degenerate:

$$I_2 = I_3 = \lambda_2 = \lambda_3 = \frac{11}{12} ma^2 \tag{2.63}$$

so there is some freedom in determining the corresponding principal axes from $(\hat{I} - \lambda_2 \mathbb{1}) \boldsymbol{\xi}^{(2,3)} = 0$, though they still should be orthogonal to each other (and $\boldsymbol{\xi}^{(1)}$). One example of a solution is:

$$\boldsymbol{\xi}^{(2)} = (1, -1, 0) \quad \text{and} \quad \boldsymbol{\xi}^{(3)} = (1, 1, -2)$$
 (2.64)

Using these principal axes and the same origin, the moment of inertia tensor becomes

$$\hat{I}_D = \frac{ma^2}{6} \begin{pmatrix} 1 & 0 & 0 \\ 0 & \frac{11}{2} & 0 \\ 0 & 0 & \frac{11}{2} \end{pmatrix} . \tag{2.65}$$

In contrast, if we had chosen the origin as the center of the cube, then one choice for the principal axes would have the same orientation, but with $\hat{I}_{\rm CM} = \frac{1}{6} m a^2 \mathbb{1}$. This result could be obtained from Eq. (2.65) using the parallel axis theorem. (Note that for a moment of inertia tensor $\propto \mathbb{1}$ that any set of orthogonal axes are principal axes due to the symmetry.)

2.4 Euler Equations

Consider the rotational motion of a rigid body about a fixed point (which could be the CM but could also be another point). We aim to describe the motion of this rigid body by exploiting properties of the body frame. To simplify things as much as possible, for this fixed point, we choose the principal axes fixed in the body frame indexed by $\alpha \in \{1, 2, 3\}$. Using the relation between time derivatives in the inertial and rotating frames, the torque is then given by:

$$\tau = \frac{d\mathbf{L}}{dt} = \frac{d_{\mathbf{R}}\mathbf{L}}{dt} + \boldsymbol{\omega} \times \mathbf{L}$$
 (2.66)

where $\boldsymbol{\omega} = \boldsymbol{\omega}(t)$. For example:

$$\tau_1 = \frac{d_{\rm R}L_1}{dt} + \omega_2 L_3 - \omega_3 L_2. \tag{2.67}$$

Not summing over repeated indices and using the formula for angular momentum along the principal axes gives $L_{\alpha} = I_{\alpha}\omega_{\alpha}$. Since we have fixed moments of inertia within the body we have $d_R I_{\alpha}/dt = 0$. Note that $d\omega/dt = d_R \omega/dt + \omega \times \omega = d_R \omega/dt$, so its rotating and inertial time derivatives are the same, and we can write $\dot{\omega}_{\alpha}$ without possible cause of confusion. Thus $d_R L_{\alpha}/dt = I_{\alpha} \dot{\omega}_{\alpha}$. This yields the Euler equations:

$$I_{1}\dot{\omega}_{1} - (I_{2} - I_{3})\omega_{2}\omega_{3} = \tau_{1}$$

$$I_{2}\dot{\omega}_{2} - (I_{3} - I_{1})\omega_{3}\omega_{1} = \tau_{2}$$

$$I_{3}\dot{\omega}_{3} - (I_{1} - I_{2})\omega_{1}\omega_{2} = \tau_{3}$$
(2.68)

where in all of these ω and τ are calculated in the rotating/body frame. This can also be written as

$$\tau_{\alpha} = I_{\alpha} \dot{\omega}_{\alpha} + \epsilon_{\alpha l k} \omega_{l} \omega_{k} I_{k}, \qquad (2.69)$$

with α fixed but a sum implied over the repeated l and k indicies. Here ϵ_{abc} is the fully antisymmetric Levi-Civita symbol.

Solving these equations gives $\omega_{\alpha}(t)$. Since the result is expressed in the body frame, rather than the inertial frame of the observer, this solution for $\omega(t)$ may not always make the physical motion transparent. To fix this we can connect our solution to the Euler angles using the relations

$$\omega_{1} = \omega_{x'} = \dot{\phi} \sin \theta \sin \psi + \dot{\theta} \cos \psi ,
\omega_{2} = \omega_{y'} = \dot{\phi} \sin \theta \cos \psi - \dot{\theta} \sin \psi ,
\omega_{3} = \omega_{z'} = \dot{\phi} \cos \theta + \dot{\psi} .$$
(2.70)

These results should be derived as exercise for the student.

Example: let us consider the stability of rigid-body free rotations ($\tau = 0$). Is a rotation $\omega = \omega_1 \mathbf{e}_1$ about the principal axis \mathbf{e}_1 stable?

Perturbations can be expressed by taking $\boldsymbol{\omega} = \omega_1 \mathbf{e}_1 + \kappa_2 \mathbf{e}_2 + \kappa_3 \mathbf{e}_3$, where κ_2 and κ_3 are small and can be treated to 1st order. The Euler equations are:

$$\dot{\omega}_1 = \frac{(I_2 - I_3)}{I_1} \,\kappa_2 \kappa_3 = O(\kappa^2) \approx 0, \tag{2.71}$$

so ω_1 is constant at this order, and

$$\dot{\kappa}_2 = \frac{(I_3 - I_1)}{I_2} \,\omega_1 \,\kappa_3 \quad \text{and} \quad \dot{\kappa}_3 = \frac{(I_1 - I_2)}{I_3} \,\omega_1 \,\kappa_2 \,.$$
(2.72)

Combining these two equations yields

$$\ddot{\kappa}_2 = \left[\frac{(I_3 - I_1)(I_1 - I_2)\omega_1^2}{I_2 I_3} \right] \kappa_2 \,, \tag{2.73}$$

and the same equation for κ_3 . The term in the square bracket is constant, and is either negative $= -w^2$ with an oscillating solution $\kappa_2 \propto \cos(wt + \phi)$, or is positive $= \alpha^2$ with exponential solutions $\kappa_2 \propto ae^{\alpha t} + be^{-\alpha t}$. If $I_1 < I_{2,3}$ or $I_{2,3} < I_1$ then the constant prefactor is negative, yielding stable oscillatory solutions. If instead $I_2 < I_1 < I_3$ or $I_3 < I_1 < I_2$ then the constant prefactor is positive, yielding an unstable exponentially growing component to their solution! This behavior can be demonstrated by spinning almost any object that has three distinct principal moments of inertia. A smart phone or tennis racket are good examples.

2.5 Symmetric Top with One Point Fixed

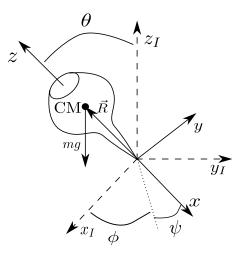
This section is devoted to a detailed analysis of a particular example that appears in many situations, the symmetric top with one point fixed, acted upon by a constant force.

Labeling the body axes as (x, y, z) and the fixed axes as (x_I, y_I, z_I) , as depicted in the right, symmetry implies that $I_1 = I_2$, and we will assume that $I_{1,2} \neq I_3$. The Euler angles are as usual (ϕ, θ, ψ) . From the figure we see that $\dot{\psi}$ is the rotation rate of the top about the (body) z-axis, $\dot{\phi}$ is the precession rate about the z_I fixed inertial axis, and $\dot{\theta}$ is the nutation rate by which the top may move away or towards the z_I axis. The Euler equations in this case are

$$I_{1}\dot{\omega}_{1} - (I_{2} - I_{3})\omega_{2}\omega_{3} = \tau_{1},$$

$$I_{1}\dot{\omega}_{2} - (I_{3} - I_{1})\omega_{3}\omega_{1} = \tau_{2},$$

$$I_{3}\dot{\omega}_{3} = 0 = \tau_{3}.$$
(2.74)



Since the CM coordinate **R** is aligned along the z-axis there is no torque along z, $\tau_3 = 0$, leading to a constant ω_3 .

There are two main cases that we will consider.

Case: $\tau = 0$ and $\dot{\theta} = 0$

The first case we will consider is when $\tau = 0$ (so there is no gravity) and $\dot{\theta} = 0$ (so there is no nutation). Then

$$\frac{d\mathbf{L}}{dt} = \boldsymbol{\tau} = 0 \quad \Rightarrow \quad \mathbf{L} = \text{constant} \tag{2.75}$$

Let us define the constant:

$$\Omega \equiv \frac{I_3 - I_1}{I_1} \omega_3 \,. \tag{2.76}$$

Then the Euler equations for this situation reduce to:

$$\dot{\omega}_1 + \Omega \omega_2 = 0 \quad \text{and} \quad \dot{\omega}_2 - \Omega \omega_1 = 0.$$
 (2.77)

The simplest solution correspond to $\omega_1(t) = \omega_2(t) = 0$, where we just have a rotation about the z-axis. Here:

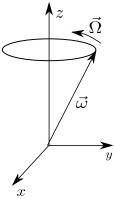
$$\mathbf{L} = L_3 \, \mathbf{e}_3$$
 where $L_3 = I_3 \omega_3$
 $\omega_1 = \omega_2 = 0 \quad \Rightarrow \quad \dot{\theta} = \dot{\phi} = 0 \text{ and } \dot{\psi} = \omega_3.$ (2.78)

In this case $\mathbf{L} \parallel \boldsymbol{\omega}$. A more general situation is when \mathbf{L} and $\boldsymbol{\omega}$ are not necessarily parallel, and ω_1 and ω_2 do not vanish. In this case Eq. (2.77) is solved by:

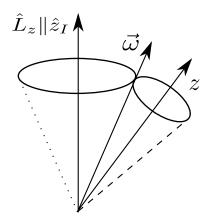
$$\omega_1 = C \sin(\Omega t + D)$$
 and $\omega_2 = -C \cos(\Omega t + D)$. (2.79)

The simple case corresponds to C=0, so now we take C>0 (since a sign can be accounted for by the constant phase D). This solution means ω precesses about the body z-axis at the rate Ω , as pictured on the right. Since $\omega_1^2 + \omega_2^2$ is constant, the full $\omega = |\omega|$ is constant, and is given by $\omega^2 = C^2 + \omega_3^2$.

The total energy here is just rotational kinetic energy $T_{\rm R} = \frac{1}{2} \boldsymbol{\omega} \cdot \mathbf{L}$ which is constant too. Thus $\boldsymbol{\omega}$ also precesses about \mathbf{L} .



We can picture this motion by thinking about a body cone that rolls around a cone in the fixed coordinate system, where in the case pictured with a larger cone about L we have $I_1 = I_2 > I_3$.



To obtain more explicit results for the motion we can relate Eq.(2.79) to Euler angles. Since $\dot{\theta} = 0$, we take $\theta = \theta_0$ to be constant. The other Euler angles come from:

$$\boldsymbol{\omega} = \begin{bmatrix} C \sin(\Omega t + D) \\ -C \cos(\Omega t + D) \\ \omega_3 \end{bmatrix} = \begin{bmatrix} \sin(\theta_0) \sin(\psi) \dot{\phi} \\ \sin(\theta_0) \cos(\psi) \dot{\phi} \\ \cos(\theta_0) \dot{\phi} + \dot{\psi} \end{bmatrix}. \tag{2.80}$$

Adding the squares of the 1st and 2nd components gives

$$C^2 = \sin^2(\theta_0) \ \dot{\phi}^2. \tag{2.81}$$

To be definite, take the positive square root of this equation to give

$$\dot{\phi} = \frac{C}{\sin(\theta_0)} \quad \Rightarrow \quad \phi = \frac{C}{\sin(\theta_0)} t + \phi_0.$$
 (2.82)

The first two equations in Eq. (2.80) are then fully solved by taking $\psi = \pi - \Omega t - D$, so we find that both ϕ and ψ have linear dependence on time. Finally the third equation gives a relation between various constants

$$\omega_3 = C \cot(\theta_0) - \Omega. \tag{2.83}$$

Thus, we see that the solution has $\dot{\phi}$ and $\dot{\psi}$ are constants with $\dot{\theta} = 0$. If we had picked the opposite sign when solving Eq. (2.81) then we would have found similar results:

$$\dot{\phi} = -\frac{C}{\sin(\theta_0)} \quad \Rightarrow \quad \psi = -\Omega t - D \quad \text{and} \quad \omega_3 = -C \cot(\theta_0) - \Omega.$$
 (2.84)

Case: $\tau \neq 0$ and $\dot{\theta} \neq 0$

Now we consider the general case where $\tau \neq 0$ and $\dot{\theta} \neq 0$. It is now more convenient to use the Lagrangian than the Euler equations directly. Since $I_1 = I_2$, using

$$T = \frac{1}{2} \left(I_1(\omega_1^2 + \omega_2^2) + I_3 \omega_3^2 \right) \quad \text{and} \quad \boldsymbol{\omega} = \begin{bmatrix} \sin(\theta) \sin(\psi) \dot{\phi} + \cos(\psi) \dot{\theta} \\ \sin(\theta) \cos(\psi) \dot{\phi} - \sin(\psi) \dot{\theta} \\ \cos(\theta) \dot{\phi} + \dot{\psi} \end{bmatrix} , \tag{2.85}$$

gives us the kinetic energy

$$T = \frac{I_1}{2} \left(\dot{\theta}^2 + \sin^2 \theta \, \dot{\phi}^2 \right) + \frac{I_3}{2} \left(\dot{\psi} + \cos \theta \, \dot{\phi} \right)^2. \tag{2.86}$$

Moreover, $V = mgR\cos(\theta)$, so in the Lagrangian L = T - V the variables ϕ and ψ are cyclic. This means that the momenta

$$p_{\phi} = \frac{\partial L}{\partial \dot{\phi}} = \left[I_1 \sin^2(\theta) + I_3 \cos^2(\theta) \right] \dot{\phi} + I_3 \cos(\theta) \dot{\psi}$$
 (2.87)

$$p_{\psi} = \frac{\partial L}{\partial \dot{\psi}} = I_3(\dot{\psi} + \cos(\theta)\dot{\phi}) = I_3\omega_3 \tag{2.88}$$

are conserved (constant). Here p_{ψ} is same as the angular momentum L_3 discussed in the case above. The torque is along the line of nodes, and p_{ϕ} and p_{ψ} correspond to two projections of **L** that are perpendicular to this torque (i.e. along $\hat{z}_{\rm I}$ and \hat{z}). Additionally, the energy is given by

$$E = T + V = \frac{I_1}{2} \left(\dot{\theta}^2 + \sin^2(\theta) \dot{\phi}^2 \right) + \frac{I_3}{2} \left(\dot{\psi} + \cos(\theta) \dot{\phi} \right)^2 + mgR\cos(\theta)$$
 (2.89)

and is also conserved. Solving the momentum equations, Eq. (2.87), for $\dot{\phi}$ and $\dot{\psi}$ gives

$$\dot{\phi} = \frac{p_{\phi} - p_{\psi} \cos(\theta)}{I_1 \sin^2(\theta)}$$

$$\dot{\psi} = \frac{p_{\psi}}{I_3} - \frac{(p_{\phi} - p_{\psi} \cos(\theta)) \cos(\theta)}{I_1 \sin^2(\theta)}$$
(2.90)

Note that once we have a solution for $\theta(t)$ that these two equations then allow us to immediately obtain solutions for $\phi(t)$ and $\psi(t)$ by integration. Eq. (2.90) can be plugged into the energy formula to give

$$E = \frac{I_1}{2}\dot{\theta}^2 + \frac{(p_\phi - p_\psi \cos(\theta))^2}{2I_1 \sin^2(\theta)} + \frac{p_\psi^2}{2I_3} + mgR\cos(\theta),$$
 (2.91)

which is a (nonlinear) differential equation for θ , since all other quantities that appear are simply constants. To simplify this result take $u = \cos(\theta)$ so that:

$$1 - u^2 = \sin^2(\theta), \qquad \dot{u} = -\sin(\theta)\dot{\theta}, \qquad \dot{\theta}^2 = \frac{\dot{u}^2}{1 - u^2}.$$
 (2.92)

Putting all this together gives:

$$\frac{\dot{u}^2}{2} = \left(\frac{2EI_3 - p_{\psi}^2}{2I_1I_3} - \frac{mgR}{I_1}u\right)\left(1 - u^2\right) - \frac{1}{2}\left(\frac{p_{\phi} - p_{\psi}u}{I_1}\right)^2 \equiv -V_{\text{eff}}(u), \qquad (2.93)$$

which is a cubic polynomial that we've defined to be the effective potential $V_{\text{eff}}(u)$. The solution to this from

$$dt = \pm \frac{du}{\sqrt{-2V_{\text{eff}}(u)}} \tag{2.94}$$

yields a complicated elliptic function, from which it is hard to get intuition for the motion. Instead, we can look at the form of $V_{\text{eff}}(u)$, because

$$\frac{1}{2}\dot{u}^2 + V_{\text{eff}}(u) = 0 \tag{2.95}$$

is the equation for the energy of a particle of unit mass m=1, kinetic energy $\dot{u}^2/2$, a potential $V_{\rm eff}(u)$, and with vanishing total energy. The cubic equation will have in general three roots where $V_{\rm eff}(u)=0$. Since the kinetic energy is always positive or zero, the potential energy must be negative or zero in the physical region, and hence the particle can not pass through any of the roots. The roots therefore serve as turning points. Furthermore, physical solutions are bounded by $-1 \le (u = \cos \theta) \le 1$. While the precise values for the roots will depend on the initial conditions or values of E, p_{ψ} , and p_{ϕ} , we can still describe the solutions in a fairly generic manner.

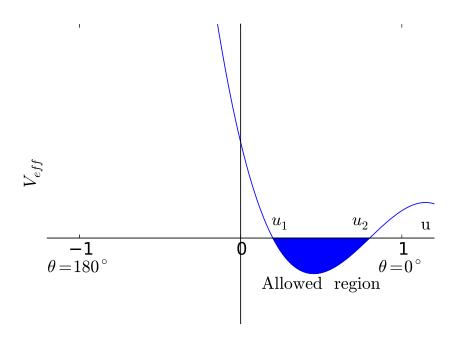


Figure 2.7: Allowed region for solutions for the top's nutation angle θ that solve Eq. (2.95).

Consider two roots u_1 and u_2 (corresponding respectively to some angles θ_1 and θ_2 as $u = \cos(\theta)$) satisfying $V_{\text{eff}}(u_1) = V_{\text{eff}}(u_2) = 0$, where $V_{\text{eff}}(u) < 0$ for $u_1 < u < u_2$; as shown in Fig. 2.7. We see that u_1 and u_2 correspond to the turning points of the motion. The region $u_1 < u < u_2$ corresponds to the region where the motion of our top lives and gives rise to a periodic nutation, where the solution bounces between the two turning points. Depending on the precise value of the various constants that appear in this V_{eff} this gives rise to different qualitative motions, with examples shown in Figs. 2.8–2.11. Recalling that $\dot{\phi} = (p_{\phi} - p_{\psi}u)/[I_1(1-u^2)]$, we see that the possible signs for $\dot{\phi}$ will depend on p_{ϕ} and p_{ψ} . In Fig. 2.8 the top nutates between θ_1 and θ_2 while always precessing in the same direction with $\dot{\phi} > 0$, whereas in Fig. 2.9 the precession is also in the backward direction, $\dot{\phi} < 0$, for part of the range of motion. In Fig. 2.10 the top has $\dot{\phi} = 0$ at θ_2 , before falling back down in the potential and gaining $\dot{\phi} > 0$ again. This figure also captures the case where we let go of a top at $\theta = \theta_2 \geq 0$ that initially has $\dot{\psi} > 0$ but $\dot{\phi} = 0$. Finally in Fig. 2.11 we have the situation where there is no nutation oscillation because the two angles coincide, $\theta_1 = \theta_2$.

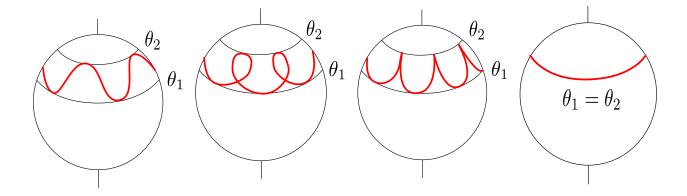


Figure 2.8: $\dot{\phi} > 0$ Figure 2.9: $\dot{\phi}$ has both Figure 2.10: at θ_2 we signs Figure 2.11: No nutation

Chapter 3

${f Vibrations} \ \& \ {f Oscillations}$

The topic of vibrations and oscillations is typically discussed in some detail in a course on waves (at MIT this is 8.03). Our goal for this chapter is to revisit aspects of oscillation phenomena using generalized coordinates. Many equations of motion we have encountered have been nonlinear. Here, we will expand about a minimum of the potential $V(q_1, \ldots, q_n)$, yielding linear equations.

Let us take $q_i = q_{0i} + \eta_i$, where \vec{q}_0 minimizes V(q), and expand in the η_i . Henceforth and until further notice, repeated indices will implicitly be summed over. Then

$$V(q_1, \dots, q_n) = V(q_{01}, \dots, q_{0n}) + \frac{\partial V}{\partial q_i} \bigg|_{0} \eta_i + \frac{1}{2} \frac{\partial^2 V}{\partial q_i \partial q_j} \bigg|_{0} \eta_i \eta_j + \dots,$$
(3.1)

where $|_0$ means "evaluate the quantity at $\vec{q_0}$ ". We already know that $\frac{\partial V}{\partial q_i}\Big|_0 = 0$ as by definition $\vec{q_0}$ minimizes V(q). As a matter of convention, we choose $V(q_0) = 0$, since this just corresponds to picking the convention for the zero of the Energy. Finally, we define the constants $V_{ij} \equiv \frac{\partial^2 V}{\partial q_i \partial q_j}\Big|_0$. There is no time dependence in the definition of our generalized coordinates, so the kinetic energy is

$$T = \frac{1}{2} m_{ij}(q_1, \dots, q_n) \dot{q}_i \dot{q}_j = \frac{1}{2} m_{ij}(q_{01}, \dots, q_{0n}) \dot{\eta}_i \dot{\eta}_j + \mathcal{O}(\eta \dot{\eta}^2), \tag{3.2}$$

where $m_{ij}(q_{01},\ldots,q_{0n}) \equiv T_{ij}$ are constants, and terms of $\mathcal{O}(\eta\dot{\eta}^2)$ and beyond are neglected. Thus, the Lagrangian to quadratic order in the η_i s is

$$L = \frac{1}{2} \left(T_{ij} \dot{\eta}_i \dot{\eta}_j - V_{ij} \eta_i \eta_j \right) . \tag{3.3}$$

From this, the equations of motion are

$$T_{ij}\ddot{\eta}_j + V_{ij}\eta_j = 0 (3.4)$$

to the same order. These are coupled linear equations of motion.

3.1 Simultaneous Diagonalization of \hat{T} and \hat{V}

To solve Eq. (3.4) lets try

$$\eta_i = a_i e^{-i\omega t} \tag{3.5}$$

where $a_i \in \mathbb{C}$ for all $i \in \{1, ..., n\}$, and the frequency ω is the same along all directions in the generalized coordinate space. Notationally, i and j will denote coordinate indices, while $i = +\sqrt{-1}$ is the imaginary unit. This gives

$$V_{ij}a_j = \omega^2 T_{ij}a_j \tag{3.6}$$

which can be rewritten in matrix form as

$$\hat{V} \cdot \vec{a} = \lambda \, \hat{T} \cdot \vec{a} \tag{3.7}$$

with $\lambda = \omega^2$. This looks like an eigenvalue equation except that when we act with the linear operator \hat{V} on \vec{a} we get back $\hat{T} \cdot \vec{a}$ instead of just the eigenvector \vec{a} . This can be rewritten as

$$(\hat{V} - \lambda \hat{T}) \cdot \vec{a} = 0 \tag{3.8}$$

where \hat{V} and \hat{T} are real and symmetric $n \times n$ matrices. In order to have a non-trivial solution of this equation we need

$$\det(\hat{V} - \lambda \hat{T}) = 0 \tag{3.9}$$

which is an n^{th} order polynomial equation with n solutions eigenvalues λ_{α} with $\alpha \in \{1, \ldots, n\}$. The solutions of $(\hat{V} - \lambda_{\alpha} \hat{T}) \cdot \vec{a}^{(\alpha)} = 0$ are the eigenvectors $\vec{a}^{(\alpha)}$. This means

$$\hat{V} \cdot \vec{a}^{(\alpha)} = \lambda_{\alpha} \hat{T} \cdot \vec{a}^{(\alpha)} \,, \tag{3.10}$$

and the solutions are much like a standard eigenvalue problem. Here and henceforth, there will be no implicit sum over repeated eigenvalue indices α (so any sums that are needed will be made explicit), but we will retain implicit sums over repeated coordinate indices i & j.

There are two cases we will consider.

1) Let us start by considering the case when \hat{T} is diagonal. In particular, let us consider the even easier case proportional to the unit matrix, where $T_{ij} = m\delta_{ij}$. This means

$$m\ddot{\eta}_i + V_{ij}\eta_j = 0. (3.11)$$

Here we have the standard eigenvalue equation

$$\hat{V} \cdot \vec{a}^{(\alpha)} = m\lambda_{\alpha}\vec{a}^{(\alpha)}. \tag{3.12}$$

The eigenvalues λ_{α} are real and nonnegative as $\lambda_{\alpha} = \omega_{\alpha}^2$; the quantities ω_{α} are the normal mode frequencies. The eigenvectors $\vec{a}^{(\alpha)}$ are orthogonal, and we can choose their normalization so that

$$m \, \vec{a}^{(\beta)\dagger} \cdot \vec{a}^{(\alpha)} = \delta_{\beta\alpha} \qquad (\text{or } \vec{a}^{(\beta)\dagger} \cdot \vec{a}^{(\alpha)} = \delta_{\beta\alpha}) \,.$$
 (3.13)

This implies that

$$\lambda_{\alpha} = \vec{a}^{(\alpha)\dagger} \cdot \hat{V} \cdot \vec{a}^{(\alpha)} \qquad \text{(or } m\lambda_{\alpha} = \vec{a}^{(\alpha)\dagger} \cdot \hat{V} \cdot \vec{a}^{(\alpha)} \text{)}. \tag{3.14}$$

The time-dependent eigenvectors are then

$$\vec{\eta}^{(\alpha)} = \vec{a}^{(\alpha)} e^{-i\omega_{\alpha}t}, \quad \text{or} \quad \eta_i^{(\alpha)} = a_i^{(\alpha)} e^{-i\omega_{\alpha}t}.$$
(3.15)

These are the normal mode solutions for the n coordinates labeled by i, and there are n such solutions labeled by α . The general solution of a linear equation is a superposition of the independent normal mode solutions:

$$\vec{\eta} = \sum_{\alpha} C_{\alpha} \, \vec{\eta}^{(\alpha)} \tag{3.16}$$

where $C_{\alpha} \in \mathbb{C}$ are fixed by initial conditions. To find real coordinate solutions, we take the real parts of these equations.

Lets prove the statements made above. Again, there will be no implicit sum over the eigenvalue index α . Dotting in $\vec{a}^{(\beta)\dagger}$ into Eq. (3.12) gives

$$\vec{a}^{(\beta)\dagger} \cdot \hat{V} \cdot \vec{a}^{(\alpha)} = m\lambda_{\alpha} \vec{a}^{(\beta)\dagger} \cdot \vec{a}^{(\alpha)}, \qquad (3.17)$$

taking the Hermitian conjugate of both sides, noting that $\hat{V}^{\dagger} = \hat{V}$, and then swapping $\alpha \leftrightarrow \beta$ gives $\vec{a}^{(\beta)\dagger} \cdot \hat{V} \cdot \vec{a}^{(\alpha)} = m \lambda_{\beta}^{\star} \vec{a}^{(\beta)\dagger} \cdot \vec{a}^{(\alpha)}$. Taking the difference of these results gives

$$(\lambda_{\alpha} - \lambda_{\beta}^{\star})\vec{a}^{(\beta)\dagger} \cdot \vec{a}^{(\alpha)} = 0, \tag{3.18}$$

and if $\alpha = \beta$ then $(\lambda_{\alpha} - \lambda_{\alpha}^{\star})\vec{a}^{(\alpha)\dagger} \cdot \vec{a}^{(\alpha)} = 0$ implies the eigenvalues are real $\lambda_{\alpha} \in \mathbb{R}$. For $\lambda_{\alpha} \neq \lambda_{\beta}$, Eq. (3.18) then implies $\vec{a}^{(\beta)\dagger} \cdot \vec{a}^{(\alpha)} = 0$ so the eigenvectors are orthogonal. If by chance $\lambda_{\alpha} = \lambda_{\beta}$ for some $\alpha \neq \beta$ then we can always simply choose the corresponding eigenvectors to be orthogonal. By convention, we then normalize the eigenvectors so that they satisfy Eq. (3.13). Finally, if $\alpha = \beta$ then Eq. (3.17) now gives $\lambda_{\alpha} = \vec{a}^{(\alpha)\dagger} \cdot \hat{V} \cdot \vec{a}^{(\alpha)}$. The statement that we are at a local minimum of the multivariable potential and not a saddle point or a maximum implies then that $\lambda_{\alpha} \geq 0$ (we have positive second derivatives in each eigenvector direction).

2) Let us now consider when \hat{T} is not diagonal and summarize which parts of the result are the same and where there are differences. Here we have $(\hat{V} - \lambda \hat{T}) \cdot \vec{a} = 0$. Again, the eigenvalues λ_{α} are real and nonnegative, with $\lambda_{\alpha} = \omega_{\alpha}^2$. Now, however,

$$\vec{a}^{(\beta)\dagger} \cdot \hat{T} \cdot \vec{a}^{(\alpha)} = 0, \qquad (3.19)$$

for $\alpha \neq \beta$, and we can replace the old normalization condition by a new one stating that

$$\vec{a}^{(\beta)\dagger} \cdot \hat{T} \cdot \vec{a}^{(\alpha)} = \delta_{\alpha\beta}, \tag{3.20}$$

which up to an overall prefactor reduces to the old orthonormality condition when $\hat{T} = m\hat{1}$. Here again,

$$\lambda_{\alpha} = \vec{a}^{(\alpha)\dagger} \cdot \hat{V} \cdot \vec{a}^{(\alpha)}, \qquad (3.21)$$

and the α^{th} normal mode solution is

$$\vec{\eta}^{(\alpha)} = \vec{a}^{(\alpha)} e^{-i\omega_{\alpha}t} \,. \tag{3.22}$$

The general solution is again the superposition

$$\vec{\eta} = \sum_{\alpha} C_{\alpha} \vec{\eta}^{(\alpha)} \,, \tag{3.23}$$

with the complex coefficients C_{α} fixed by the initial conditions (and a real part taken to get real coordinates).

Lets repeat the steps of our proof for this case. Dotting $\vec{a}^{(\beta)\dagger}$ into Eq. (3.10) gives

$$\vec{a}^{(\beta)\dagger} \cdot \hat{V} \cdot \vec{a}^{(\alpha)} = \lambda_{\alpha} \vec{a}^{(\beta)\dagger} \cdot \hat{T} \cdot \vec{a}^{(\alpha)}. \tag{3.24}$$

Taking the Hermitian conjugate and swapping α and β yields $\vec{a}^{(\beta)\dagger} \cdot \hat{V} \cdot \vec{a}^{(\alpha)} = \lambda_{\beta}^{\star} \vec{a}^{(\beta)\dagger} \cdot \hat{T} \cdot \vec{a}^{(\alpha)}$. Subtracting the two results this gives

$$(\lambda_{\alpha} - \lambda_{\beta}^{\star}) \, \vec{a}^{(\beta)\dagger} \cdot \hat{T} \cdot \vec{a}^{(\alpha)} = 0. \tag{3.25}$$

If we consider $\alpha = \beta$ then $(\lambda_{\alpha} - \lambda_{\alpha}^{\star})\vec{a}^{(\alpha)\dagger} \cdot \hat{T} \cdot \vec{a}^{(\alpha)} = 0$ implies $\lambda_{\alpha} \in \mathbb{R}$ since $\hat{T} = \hat{T}^{\star} = \hat{T}^{\top}$ and physically we know that the kinetic energy $T = \dot{\vec{\eta}} \cdot \hat{T} \cdot \dot{\vec{\eta}} > 0$ for any $\dot{\vec{\eta}} \neq 0$. For $\lambda_{\alpha} \neq \lambda_{\beta}$, then the condition instead implies $\vec{a}^{(\beta)\dagger} \cdot \hat{T} \cdot \vec{a}^{(\alpha)} = 0$ so the eigenvectors are orthogonal (with respect to an inner product defined with \hat{T}). If by chance $\lambda_{\alpha} = \lambda_{\beta}$ for some $\alpha \neq \beta$ then we can simply choose the corresponding eigenvectors to be orthogonal in this sense. By convention, we normalize the eigenvectors so that they will be orthonormal as in Eq. (3.20). Finally, if $\alpha = \beta$ then $\lambda_{\alpha} = \vec{a}^{(\alpha)\dagger} \cdot \hat{V} \cdot \vec{a}^{(\alpha)}$, which is positive, so $\lambda_{\alpha} > 0$ also. The statement that we are at a local minimum of the potential and not a saddle point or a maximum implies then that $\lambda_{\alpha} \geq 0$.

3.2 Vibrations and Oscillations with Normal Coordinates

Given these results, it is natural to ask whether a different set of generalized coordinates might be better for studying motion about the minimum?

We form the matrix A by placing the eigenvectors in columns

$$A = \begin{bmatrix} \vec{a}^{(1)} & \vec{a}^{(2)} & \dots & \vec{a}^{(n)} \end{bmatrix}$$
 (3.26)

and construct a diagonal eigenvalue matrix $\hat{\lambda} = \operatorname{diag}(\lambda_1, \lambda_2, \dots, \lambda_n)$. The matrix A can be ensured to be real because each $\vec{a}^{(\alpha)}$ only has at most an overall phase¹, and these can be removed by putting them into the coefficients C_{α} . The matrix A simultaneously diagonalizes \hat{T} and \hat{V} since

$$A^{\mathsf{T}}\hat{T}A = \mathbb{1} \quad \text{and} \quad A^{\mathsf{T}}\hat{V}A = \hat{\lambda}.$$
 (3.27)

We choose new normal coordinates $\vec{\xi}$ by letting

$$\vec{\eta} = A \vec{\xi} \quad \text{and} \quad \dot{\vec{\eta}} = A \dot{\vec{\xi}}$$
(3.28)

so that the Lagrangian

$$L = \frac{1}{2} \dot{\vec{\eta}} \cdot \hat{T} \cdot \dot{\vec{\eta}} - \frac{1}{2} \vec{\eta} \cdot \hat{V} \cdot \vec{\eta}$$

$$= \frac{1}{2} \dot{\vec{\xi}} \cdot (A^{\top} \hat{T} A) \cdot \dot{\vec{\xi}} - \frac{1}{2} \vec{\xi} \cdot (A^{\top} \hat{V} A) \cdot \vec{\xi}$$

$$= \frac{1}{2} \sum_{\alpha} \left(\dot{\xi}_{\alpha}^{2} - \omega_{\alpha}^{2} \xi_{\alpha}^{2} \right).$$
(3.29)

This gives the simple equations of motion for each α :

$$\ddot{\xi}_{\alpha} + \omega_{\alpha}^2 \, \xi_{\alpha} = 0 \,. \tag{3.30}$$

Thus, each normal coordinate describes the oscillations of the system with normal mode frequency ω_{α} .

Example: Let us consider the triatomic molecule CO_2 shown in Figure 3.1. We can picture it as a carbon atom of mass M in the middle of two oxygen atoms each of mass m. For the three particles there are 9 coordinates given by \mathbf{r}_1 , \mathbf{r}_2 , and \mathbf{r}_3 . Six of these coordinates correspond to translations and rotations of the mass system treated as a rigid body. This leaves 3 coordinates that correspond to internal motions of the system. To model the potential we connect each oxygen atom to the carbon atom with a spring of constant k and relaxed length k. This does not add any cost to relative motion of the atoms with fixed spring length, which we will address below by adding another potential term in order to favor the linear configuration.

It is straightforward to guess what the normal modes could be:

¹Why is it just an overall phase? The equation $\hat{V} \cdot \vec{a}^{(\alpha)} = \lambda_{\alpha} \hat{T} \cdot \vec{a}^{(\alpha)}$ alone does not fix the normalization of $\vec{a}^{(\alpha)}$. Let us say we pick $a_i^{(\alpha)} \in \mathbb{R}$ for some i. Then $V_{kj}a_j^{(\alpha)} = \lambda_{\alpha}K_{kj}a_j^{(\alpha)}$ is a set of equations with all real coefficients and one real term in the sums. Hence the solutions $a_j^{(\alpha)}/a_i^{(\alpha)} \in \mathbb{R}$ for all $j \in \{1, \ldots, n\}$, implying that at most there is an overall phase in $\vec{a}^{(\alpha)}$.

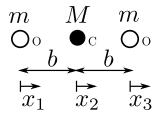


Figure 3.1: The CO_2 molecule.

- The oxygen atoms moving in the same direction along the line and the carbon atom moving in the opposite direction. This is a longitudinal oscillation.
- The oxygen atoms opposing each other along the line while the carbon atom remains at rest. This is a longitudinal oscillation.
- The oxygen atoms move in the same direction perpendicular to the line and the carbon atom moving in the opposite direction. This is a transverse oscillation.

These three normal modes are shown in Figure 3.2.

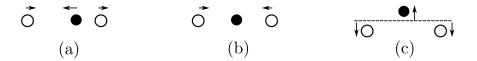


Figure 3.2: The three Normal Modes of the CO₂ molecule

We pick the body frame axes as follows:

- the three particles are in the xy-plane fixing 3 coordinates $z_i = 0$ for $i \in \{1, 2, 3\}$,
- the origin is the CM so $m(x_1 + x_3) + Mx_2 = m(y_1 + y_3) + My_2 = 0$, which fixes two more coordinates,
- the axes are oriented so that $y_1 = y_3$, which fixes one coordinate.

Defining the mass ratio as $\rho \equiv \frac{m}{M}$, then $x_2 = -\rho(x_1 + x_3)$ and $y_2 = -2\rho y_1$ can be eliminated. Altogether this fixes 6 coordinates, leaving the coordinates (x_1, x_3, y_1) . This setup is shown in Figure 3.3.

For the potential we take

$$V = \frac{k}{2} (s_1 - b)^2 + \frac{k}{2} (s_2 - b)^2 + \frac{\lambda b^2}{2} (\alpha_1^2 + \alpha_2^2).$$
 (3.31)

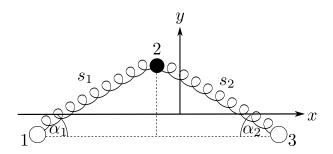


Figure 3.3: The orientation of the CO_2 molecule on xy plane

The first two terms are the springs discussed previously, and the last two provide a quadratic energy cost to the springs rotating away from the linear configuration, with strength given by λ . The spring lengths are

$$s_1 = \sqrt{(x_1 - x_2)^2 + (y_1 - y_2)^2} = \sqrt{[x_1 + \rho(x_1 + x_3)]^2 + (1 + 2\rho)^2 y_1^2}$$

$$s_2 = \sqrt{(x_2 - x_3)^2 + (y_2 - y_3)^2} = \sqrt{[x_3 + \rho(x_1 + x_3)]^2 + (1 + 2\rho)^2 y_1^2},$$
(3.32)

and the two angles are

$$\alpha_1 = \tan^{-1} \left(\frac{y_3 - y_2}{x_3 - x_2} \right) = \tan^{-1} \left[\frac{(1 + 2\rho)y_1}{(1 + \rho)x_1 + x_3} \right],$$

$$\alpha_2 = \tan^{-1} \left(\frac{y_2 - y_1}{x_2 - x_1} \right) = \tan^{-1} \left[\frac{(1 + 2\rho)y_1}{(1 + \rho)x_3 + x_1} \right].$$
(3.33)

These results give $V = V(x_1, y_1, x_3)$.

For the kinetic energy we have

$$T = \frac{m}{2} \left(\dot{x}_1^2 + \dot{x}_3^2 \right) + \frac{M}{2} \dot{x}_2^2 + \frac{m}{2} \left(\dot{y}_1^2 + \dot{y}_3^2 \right) + \frac{M}{2} \dot{y}_2^2, \tag{3.34}$$

which after eliminating coordinates becomes

$$T = \frac{m}{2} (1 + \rho) \left(\dot{x}_1^2 + \dot{x}_3^2 \right) + m\rho \dot{x}_1 \dot{x}_3 + m (1 + 2\rho) \dot{y}_1^2.$$
 (3.35)

Equilibrium comes from taking $y_1 = 0$, $x_3 = -x_1 = b$, which implies $\alpha_1 = \alpha_2 = 0$, $s_1 = s_2 = b$, and V = 0. We define coordinates for expanding about this potential minimum as $\eta_1 = x_1 + b$, $\eta_3 = x_3 - b$, and $\eta_2 = y_1$. Then as V(-b, 0, b) = 0 in equilibrium and $\frac{\partial V}{\partial \eta_i}\Big|_{0} = 0$ then $V = \frac{1}{2}V_{ij}\eta_i\eta_j + \ldots$ where

$$V_{ij} = \frac{\partial^2 V}{\partial \eta_i \partial \eta_j} \Big|_{0} = \begin{bmatrix} k(1 + 2\rho + 2\rho^2) & 0 & 2k\rho(1+\rho) \\ 0 & 2\lambda(1 + 2\rho)^2 & 0 \\ 2k\rho(1+\rho) & 0 & k(1 + 2\rho + 2\rho^2) \end{bmatrix}$$
(3.36)

for this system. Additionally,

$$T_{ij} = \begin{bmatrix} m(1+\rho) & 0 & m\rho \\ 0 & 2m(1+2\rho) & 0 \\ m\rho & 0 & m(1+\rho) \end{bmatrix}$$
(3.37)

for this system. Since there are no off-diagonal terms in the 2nd row or 2nd column in either \hat{V} or \hat{T} , the transverse and the longitudinal modes decouple. For the transverse mode, we are left with

$$\ddot{y_1} + \frac{2\lambda(1+2\rho)^2}{2m(1+2\rho)}y_1 = 0, \qquad (3.38)$$

which is a simple harmonic oscillator. For the longitudinal modes, we have $\vec{\eta} = (\eta_1, \eta_3)$. The frequencies come from

$$\det \begin{bmatrix} k(1+2\rho+2\rho^2) - \lambda m(1+\rho) & 2k\rho(1+\rho) - \lambda m\rho \\ 2k\rho(1+\rho) - \lambda m\rho & k(1+2\rho+2\rho^2) - \lambda m(1+\rho) \end{bmatrix} = 0$$
 (3.39)

The solutions give the normal mode frequencies

$$\lambda_1 = \omega_1^2 = \frac{k}{m}, \qquad \lambda_2 = \omega_2^2 = \frac{k}{m} (1 + 2\rho), \qquad (3.40)$$

with associated eigenvectors

$$\vec{a}^{(1)} = \frac{1}{\sqrt{2m}} \begin{pmatrix} 1 \\ -1 \end{pmatrix}, \qquad \vec{a}^{(2)} = \frac{1}{\sqrt{2m(1+2\rho)}} \begin{pmatrix} 1 \\ 1 \end{pmatrix}, \qquad (3.41)$$

which were chosen to satisfy $\vec{a}^{(\alpha)} \cdot \hat{T} \cdot \vec{a}^{(\beta)} = \delta_{\alpha\beta}$. Thus, the normal coordinates for the longitudinal modes are $\xi_1 \propto x_1 - x_3$ and $\xi_2 \propto x_1 + x_3$. Oscillations in these coordinates correspond to the normal mode motions in Fig. 3.2(b) and Fig. 3.2(a) respectively.

Chapter 4

Canonical Transformations, Hamilton-Jacobi Equations, and Action-Angle Variables

We've made good use of the Lagrangian formalism. Here we'll study dynamics with the Hamiltonian formalism. Problems can be greatly simplified by a good choice of generalized coordinates. How far can we push this?

Example: Let us imagine that we find coordinates q_i that are all cyclic. Then $\dot{p}_i = 0$, so $p_i = \alpha_i$ are all constant. If H is conserved, then:

$$H = H(\alpha_1, \dots, \alpha_n) \tag{4.1}$$

is also constant in time. In such a case the remaining equations of motion:

$$\dot{q}_i = \frac{\partial H}{\partial \alpha_i} = \omega_i(\alpha) \quad \Rightarrow \quad q_i = \omega_i t + \delta_i$$
 (4.2)

All coordinates are linear in time and the motion becomes very simple.

We might imagine searching for a variable transformation to make as many coordinates as possible cyclic. Before proceeding along this path, we must see what transformations are allowed.

4.1 Generating Functions for Canonical Transformations

Recall the Euler-Lagrange equations are invariant when:

- a point transformation occurs Q = Q(q, t) with L[q, t] = L'[Q, t];
- a total derivative is summed to the Lagrangian $L' = L + \frac{dF[q, t]}{dt}$.

For H we consider point transformations in phase space:

$$Q_i = Q_i(q, p, t) \quad \text{and} \quad P_i = P_i(q, p, t), \tag{4.3}$$

where the Hamilton's equations for the evolution of the canonical variables (q, p) are satisfied:

$$\dot{q}_i = \frac{\partial H}{\partial p_i}$$
 and $\dot{p}_i = -\frac{\partial H}{\partial q_i}$. (4.4)

Generally, not all transformations preserve the equations of motion. However, the transformation is *canonical* if there exists a new Hamiltonian:

$$K = K(Q, P, t), \tag{4.5}$$

where

$$\dot{Q}_i = \frac{\partial K}{\partial P_i}$$
 and $\dot{P}_i = -\frac{\partial K}{\partial Q_i}$. (4.6)

For notational purposes let repeated indices be summed over implicitly.

Hamilton's principle can be written as:

$$\delta \int_{t_1}^{t_2} (p_i \dot{q}_i - H(q, p, t)) dt = 0, \tag{4.7}$$

or in the new Hamiltonian as:

$$\delta \int_{t_1}^{t_2} \left(P_i \dot{Q}_i - K(Q, P, t) \right) dt = 0.$$
 (4.8)

For the Eq.(4.7) to imply Eq.(4.8), then we need:

$$\lambda(p_i \dot{q}_i - H) = P_i \dot{Q}_i - K + \dot{F}. \tag{4.9}$$

Since \dot{F} is a total derivative and the ends of the path are fixed:

$$\delta q|_{t_1}^{t_2} = 0$$
 and $\delta p|_{t_1}^{t_2} = 0$ \Rightarrow $\delta F|_{t_1}^{t_2} = 0$ (4.10)

There are a few things to be said about transformations and λ .

- If $\lambda = 1$ then the transformation is *canonical*, which is what we will study.
- If $\lambda \neq 1$ then the transformation is *extended canonical*, and the results from $\lambda = 1$ can be recovered by rescaling q and p appropriately.

• If $Q_i = Q_i(q, p)$ and $P_i = P_i(q, p)$ without explicit dependence on time, then the transformation is restricted canonical.

We will always take transformations $Q_i = Q_i(q, p, t)$ and $P_i = P_i(q, p, t)$ to be invertible in any of the canonical variables. If F depends on a mix of old and new phase space variables, it is called a *generating function* of the canonical transformation. There are four important cases of this.

1. Let us take

$$F = F_1(q, Q, t) \tag{4.11}$$

where the old coordinates q_i and the new coordinates Q_i are independent. Then:

$$p_i \dot{q}_i - H = P_i \dot{Q}_i - K + \dot{F}_1 = P_i \dot{Q}_i - K + \frac{\partial F_1}{\partial t} + \frac{\partial F_1}{\partial q_i} \dot{q}_i + \frac{\partial F_1}{\partial Q_i} \dot{Q}_i$$
(4.12)

from this we see that $P_i\dot{Q}_i$ cancels and equating the terms with a \dot{q}_i , a \dot{Q}_i and the remaining terms gives:

$$p_i = \frac{\partial F_1}{\partial q_i}$$
 , $P_i = -\frac{\partial F_1}{\partial Q_i}$ and $K = H + \frac{\partial F_1}{\partial t}$, (4.13)

which gives us formula for a transformation:

$$p_i = p_i(q, Q, t)$$
 and $P_i = P_i(q, Q, t)$ (4.14)

and connects K to an initial H.

Example: if

$$F_1 = -\frac{Q}{q},\tag{4.15}$$

then:

$$p = \frac{\partial F_1}{\partial q} = \frac{Q}{q^2}$$
 and $P = -\frac{\partial F_1}{\partial Q} = \frac{1}{q}$. (4.16)

Writing the new coordinates as function of the old ones yields

$$Q = pq^2 \quad \text{and} \quad P = \frac{1}{q} \tag{4.17}$$

Example: Given the transformations

$$Q = \ln\left(\frac{p}{q}\right)$$
 and $P = -\left(\frac{q^2}{2} + 1\right)\frac{p}{q}$, (4.18)

we can prove they are canonical by finding a corresponding generating function. We know:

$$\frac{\partial F_1}{\partial q} = p = qe^Q,\tag{4.19}$$

which gives us

$$F_1 = \int qe^Q dq + g(Q) = \frac{q^2}{2}e^Q + g(Q), \tag{4.20}$$

and

$$P = -\frac{\partial F_1}{\partial Q} = -\frac{q^2}{2}e^Q - \frac{dg}{dQ} = -\left(\frac{q^2}{2} + 1\right)\frac{p}{q} = -\left(\frac{q^2}{2} + 1\right)e^Q$$

$$\Rightarrow q(Q) = e^Q. \tag{4.21}$$

Thus F_1 is given by:

$$F_1 = \left(\frac{q^2}{2} + 1\right)e^Q. {(4.22)}$$

2. Let:

$$F = F_2(q, P, t) - Q_i P_i (4.23)$$

where we wish to treat the old coordinates q_i and new momenta P_i as independent variables. Then:

$$\dot{q}_i p_i - H = \dot{Q}_i P_i - K + \dot{F}_2 - \dot{Q}_i P_i - Q_i \dot{P}_i = -Q_i \dot{P}_i - K + \frac{\partial F_2}{\partial t} + \frac{\partial F_2}{\partial a_i} \dot{q}_i + \frac{\partial F_2}{\partial P_i} \dot{P}_i \quad (4.24)$$

This corresponds to

$$p_i = \frac{\partial F_2}{\partial q_i}$$
 ; $Q_i = \frac{\partial F_2}{\partial P_i}$ and $K = H + \frac{\partial F_2}{\partial t}$. (4.25)

3. We could also take

$$F = F_3(p, Q, t) + q_i p_i (4.26)$$

with the new coordinates Q_i and the old momenta p_i as independent variables.

4. Finally we could take

$$F = F_4(p, P, t) + q_i p_i - Q_i P_i (4.27)$$

with the old momenta p_i and new momenta P_i as independent variables.

This can be summarized in the table below.

CHAPTER 4. CANONICAL TRANSFORMATIONS, HAMILTON-JACOBI EQUATIONS, AND ACTION-ANGLE VARIABLES

Function	Transformations	Simplest case	
$F_1(q,Q,t)$	$p_i = \frac{\partial F_1}{\partial q_i}, P_i = -\frac{\partial F_1}{\partial Q_i}$	$F_1 = q_i Q_i$	$Q_i = p_i, P_i = -q_i$
$F_2(q, P, t)$	$p_i = \frac{\partial F_2}{\partial q_i}, Q_i = \frac{\partial F_2}{\partial P_i}$	$F_2 = q_i P_i$	$Q_i = q_i, P_i = p_i$
$F_3(p,Q,t)$	$q_i = -\frac{\partial F_3}{\partial p_i}, P_i = -\frac{\partial F_3}{\partial Q_i}$	$F_3 = p_i Q_i$	$Q_i = -q_i, P_i = -p_i$
$F_4(p, P, t)$	$q_i = -\frac{\partial F_4}{\partial p_i}, Q_i = \frac{\partial F_4}{\partial P_i}$	$F_4 = p_i P_i$	$Q_i = p_i, P_i = -q_i$

The simplest case of the 2^{nd} (F_2) transformation is just an identity transformation. For any of these F_i cases we also have:

$$K = H + \frac{\partial F_i}{\partial t}. (4.28)$$

If F_i is independent of time then this implies

$$K = H \tag{4.29}$$

Mixed cases may also occur when more than two old canonical coordinates are present. (In this chapter we will be using Einstein's repeated index notation for implicit summation, unless otherwise stated.)

Example: consider

$$F_2 = f_i(q, t)P_i \tag{4.30}$$

for some functions f_i where $i \in \{1, ..., n\}$. Then

$$Q_i = \frac{\partial F_2}{\partial P_i} = f_i(q, t) \tag{4.31}$$

is a coordinate point transformation. It is canonical with

$$p_i = \frac{\partial f_i}{\partial q_j} P_j, \tag{4.32}$$

which can be inverted to get $P_j = P_j(q, p, t)$.

Example: Consider the harmonic oscillator:

$$H = \frac{p^2}{2m} + \frac{kq^2}{2} \quad \text{where} \quad k = m\omega^2 \tag{4.33}$$

Let us try the transformation:

$$p = \alpha \sqrt{2mP} \cos(Q)$$

$$q = \frac{\alpha}{m\omega} \sqrt{2mP} \sin(Q)$$
(4.34)

for α constant. Then:

$$K = H = P\alpha^2 \left(\cos^2(Q) + \sin^2(Q)\right) = P\alpha^2,$$
 (4.35)

so the new momentum

$$P = \frac{E}{\alpha^2} \tag{4.36}$$

is just proportional to the energy, while Q is a cyclic variable.

Is this transformation canonical? We can find a generating function $F = F_1(q, Q)$ by dividing the old variables:

$$\frac{p}{q} = m\omega \cot(Q). \tag{4.37}$$

This gives us:

$$p = \frac{\partial F_1}{\partial q} \Rightarrow F_1 = \int p(q, Q) \, dq + g(Q) = \frac{1}{2} m \omega q^2 \cot(Q) + g(Q)$$

$$P = -\frac{\partial F_1}{\partial Q} = \frac{m \omega q^2}{2 \sin^2(Q)} - \frac{dg}{dQ}$$

$$(4.38)$$

Setting:

$$\frac{dg}{dQ} = 0 \Rightarrow q^2 = \frac{2P}{m\omega}\sin^2(Q),\tag{4.39}$$

which tells us the transformation is canonical if $\alpha = \sqrt{\omega}$. This means:

$$P = \frac{E}{\omega} \tag{4.40}$$

By Hamilton's equations Eq.(4.4):

$$\dot{Q} = \frac{\partial K}{\partial P} = \omega \Rightarrow Q = \omega t + \delta.$$
 (4.41)

Putting this altogether, this gives the familiar results:

$$q = \sqrt{\frac{2E}{m\omega^2}}\sin(\omega t + \delta)$$

$$p = \sqrt{2mE}\cos(\omega t + \delta).$$
(4.42)

Lets record for future use our final canonical transformation here:

$$q = \sqrt{\frac{2P}{m\omega}}\sin(Q)$$
, $p = \sqrt{2m\omega P}\cos(Q)$.

So far, a transformation Q = Q(q, p, t) and P = P(q, p, t) is canonical if we can find a generating function F. This involves integration, which could be complicated, so it would be nice to have a test that only involves differentiation. There is one!

4.2 Poisson Brackets and the Symplectic Condition

In Classical Mechanics II (8.223) the *Poisson bracket* of the quantities u and v was defined as

$$\{u, v\}_{q,p} = \sum_{i} \left(\frac{\partial u}{\partial q_i} \frac{\partial v}{\partial p_i} - \frac{\partial v}{\partial q_i} \frac{\partial u}{\partial p_i} \right)$$

$$(4.43)$$

It is easy to check that the following fundamental Poisson bracket relations are satisfied:

$$\{q_i, q_j\}_{q,p} = \{p_i, p_j\}_{q,p} = 0 \quad \text{and} \quad \{q_i, p_j\}_{q,p} = \delta_{ij}.$$
 (4.44)

There are a few other properties of note. These include:

$$\{u, u\} = 0, \tag{4.45}$$

$$\{u, v\} = -\{v, u\},\tag{4.46}$$

$$\{au + bv, w\} = a\{u, w\} + b\{v, w\}, \tag{4.47}$$

$$\{uv, w\} = u\{v, w\} + \{u, w\}v,$$
 (4.48)

$$\{u, \{v, w\}\} + \{v, \{w, u\}\} + \{w, \{u, v\}\} = 0, \tag{4.49}$$

for a, b constants. Eq.(4.49) is the Jacobi identity.

The above looks a lot like the commutators of operators in quantum mechanics, such as:

$$[\hat{x}, \hat{p}] = i\hbar \tag{4.50}$$

Indeed, quantizing a classical theory by replacing Poisson brackets with commutators through:

$$[u,v] = i\hbar\{u,v\} \tag{4.51}$$

is a popular approach (first studied by Dirac). It is also the root of the name "canonical quantization". (Note that Eq.(4.48) was written in a manner to match the analogous formula in quantum mechanics where the operator ordering is important, just in case its familiar. Here we can multiply functions in either order.)

Now we can state the desired criteria that only involves derivatives.

Theorem: A transformation $Q_i = Q_i(q, p, t)$ and $P_i = P_i(q, p, t)$ is canonical if and only if:

$${Q_i, Q_j}_{q,p} = {P_i, P_j}_{q,p} = 0$$
 and ${Q_i, P_j}_{q,p} = \delta_{ij}$. (4.52)

To prove it, we'll need some more notation. Let's get serious about treating q_i and p_i on an equal footing together, defining the following two quantities:

$$\vec{\eta} = \begin{pmatrix} q_1 \\ \vdots \\ q_n \\ p_1 \\ \vdots \\ p_n \end{pmatrix} \text{ and } J = \begin{bmatrix} 0_{n \times n} & \mathbb{1}_{n \times n} \\ -\mathbb{1}_{n \times n} & 0_{n \times n} \end{bmatrix}$$

$$(4.53)$$

where $0_{n\times n}$ is the $n\times n$ zero matrix, $1_{n\times n}$ is the $n\times n$ identity matrix. The following properties of J will be useful:

$$J^{\top} = -J$$
 , $J^{2} = -\mathbb{1}_{2n \times 2n}$ and $J^{\top}J = JJ^{\top} = \mathbb{1}_{2n \times 2n}$. (4.54)

We also note that det(J) = 1.

With this notation Hamilton's equations, Eq.(4.4), can be rewritten as:

$$\dot{\vec{\eta}} = J \frac{\partial H}{\partial \vec{\eta}} \quad \text{or} \quad \dot{\vec{\eta}} = J \nabla_{\vec{\eta}} H \,.$$
 (4.55)

The notation $\nabla_{\vec{\eta}} H$ better emphasizes that this quantity is a vector, but we will stick to using the first notation for this vector, $\partial H/\partial \vec{\eta}$, below.

Although the Theorem is true for time dependent transformations, lets carry out the proof for the simpler case of time independent transformations $Q_i = Q_i(q, p)$ and $P_i = P_i(q, p)$. This implies K = H. Let us define:

$$\vec{\xi} = \begin{pmatrix} Q_1 \\ \vdots \\ Q_n \\ P_1 \\ \vdots \\ P_n \end{pmatrix} \tag{4.56}$$

which is a function of the original coordinates, so we can write:

$$\vec{\xi} = \vec{\xi}(\vec{\eta}) \tag{4.57}$$

Now consider the time derivative of $\vec{\xi}$:

$$\dot{\xi}_i = \frac{\partial \xi_i}{\partial \eta_j} \dot{\eta}_j \quad \Leftrightarrow \quad \dot{\vec{\xi}} = M \dot{\vec{\eta}} \quad \text{where} \quad M_{ij} = \frac{\partial \xi_i}{\partial \eta_j}.$$
(4.58)

Here M corresponds to the Jacobian of the transformation.

From the Hamilton's equations, we know that

$$\dot{\vec{\eta}} = J \frac{\partial H}{\partial \vec{\eta}}.\tag{4.59}$$

We want to show that:

$$\dot{\vec{\xi}} = J \frac{\partial H}{\partial \vec{\xi}} \quad \text{for} \quad \vec{\xi} = \vec{\xi}(\vec{\eta}) \text{ a canonical transformation.}$$
(4.60)

Let us now consider:

$$\dot{\xi}_{i} = \frac{\partial \xi_{i}}{\partial \eta_{j}} \dot{\eta}_{j} = \frac{\partial \xi_{i}}{\partial \eta_{j}} J_{jk} \frac{\partial H}{\partial \eta_{k}} = \frac{\partial \xi_{i}}{\partial \eta_{j}} J_{jk} \frac{\partial \xi_{l}}{\partial \eta_{k}} \frac{\partial H}{\partial \xi_{l}} \quad \Leftrightarrow \quad \dot{\vec{\xi}} = M J M^{\top} \frac{\partial H}{\partial \vec{\xi}}$$
(4.61)

for any H. Then

$$\vec{\xi} = \vec{\xi}(\vec{\eta})$$
 is a canonical transformation iff $MJM^{\top} = J$ (4.62)

is satisfied. This is known as the "symplectic condition". Moreover, since

$$MJ = J(M^{\top})^{-1}$$
 and $J^2 = -1$, (4.63)

we can write:

$$J(MJ)J = -JM = J(JM^{T-1})J = -M^{T-1}J \Rightarrow JM = M^{T-1}J. \tag{4.64}$$

Thus we see the equivalence

$$MJM^{\top} = J \quad \Leftrightarrow \quad M^{\top}JM = J.$$
 (4.65)

Now consider Poisson brackets in this matrix notation:

$$\{u, v\}_{q,p} = \{u, v\}_{\vec{\eta}} = \left(\frac{\partial u}{\partial \vec{\eta}}\right)^{\perp} J \frac{\partial v}{\partial \vec{\eta}}$$

$$\tag{4.66}$$

and the fundamental Poisson brackets are:

$$\{\eta_i, \eta_j\}_{\vec{\eta}} = J_{ij} \tag{4.67}$$

Then we can calculate the Poisson brackets that appeared in the theorem we are aiming to prove as

$$\{\xi_i, \xi_j\}_{\vec{\eta}} = \left(\frac{\partial \xi_i}{\partial \vec{\eta}}\right)^{\top} J \frac{\partial \xi_j}{\partial \vec{\eta}} = (M^{\top} J M)_{ij}$$
(4.68)

This last equation is the same as Eq.(4.65). The new variables satisfy the Poisson bracket relationships Eq.(4.67):

$$\{\xi_i, \xi_j\}_{\vec{\eta}} = J_{ij} \tag{4.69}$$

if and only if

$$M^{\top}JM = J \tag{4.70}$$

which using Eqs. (4.62) and (4.65) is true if and only if $\vec{\xi} = \vec{\xi}(\vec{\eta})$ is canonical. This completes the proof.

There are two facts that arise from this.

• Poisson brackets are canonical invariants

$$\{u, v\}_{\vec{\eta}} = \{u, v\}_{\vec{\epsilon}} = \{u, v\}.$$
 (4.71)

This is true because:

$$\{u, v\}_{\vec{\eta}} = \left(\frac{\partial u}{\partial \vec{\eta}}\right)^{\top} J \frac{\partial v}{\partial \vec{\eta}} = \left(M^{\top} \frac{\partial u}{\partial \vec{\xi}}\right)^{\top} J \left(M^{\top} \frac{\partial v}{\partial \vec{\xi}}\right) \tag{4.72}$$

$$= \left(\frac{\partial u}{\partial \vec{\xi}}\right)^{\top} M J M^{\top} \frac{\partial v}{\partial \vec{\xi}} = \left(\frac{\partial u}{\partial \vec{\xi}}\right)^{\top} J \frac{\partial v}{\partial \vec{\xi}} = \{u, v\}_{\vec{\xi}}$$
(4.73)

• Phase space volume elements are preserved by canonical transformations, as discussed in 8.223. Phase space volume is given by:

$$V_{\vec{\xi}} = \prod_{i} dQ_{i} dP_{i} = |\det(M)| \prod_{j} dq_{j} dp_{j} = |\det(M)| V_{\vec{\eta}}. \tag{4.74}$$

Since we have

$$\det(M^{\top}JM) = \det(J) = (\det(M))^2 \det(J) \Rightarrow |\det(M)| = 1, \qquad (4.75)$$

and the two phase space volumes are the same.

4.3 Equations of Motion & Conservation Theorems

Let us consider a function:

$$u = u(q, p, t) \tag{4.76}$$

Then:

$$\dot{u} = \frac{\partial u}{\partial q_i} \dot{q}_i + \frac{\partial u}{\partial p_i} \dot{p}_i + \frac{\partial u}{\partial t} = \frac{\partial u}{\partial q_i} \frac{\partial H}{\partial p_i} - \frac{\partial u}{\partial p_i} \frac{\partial H}{\partial q_i} + \frac{\partial u}{\partial t}, \tag{4.77}$$

which can be written more concisely as

$$\dot{u} = \{u, H\} + \frac{\partial u}{\partial t} \tag{4.78}$$

for any canonical variables (q, p) and corresponding Hamiltonian H. Performing canonical quantization on this yields the Heisenberg equation of time evolution in quantum mechanics. There are a few easy cases to check.

• If $u = q_i$ then:

$$\dot{q}_i = \{q_i, H\} + \frac{\partial q_i}{\partial t} = \frac{\partial H}{\partial p_i}$$
 (4.79)

• If $u = p_i$ then:

$$\dot{p}_i = \{p_i, H\} + \frac{\partial p_i}{\partial t} = -\frac{\partial H}{\partial q_i} \tag{4.80}$$

Together the above two cases yield Hamilton's equations of motion.

• Also, if u = H then:

$$\dot{H} = \{H, H\} + \frac{\partial H}{\partial t} = \frac{\partial H}{\partial t}$$
 (4.81)

Next consider what we can say if a quantity u is conserved. Here:

$$\dot{u} = 0 = \{u, H\} + \frac{\partial u}{\partial t}.\tag{4.82}$$

As a corollary, if

$$\frac{\partial u}{\partial t} = 0, (4.83)$$

then

$${u, H} = 0 \Rightarrow u \text{ is conserved.}$$
 (4.84)

(In quantum mechanics this the analog of saying that u is conserved if u commutes with H.)

Another fact, is that if u and v are conserved then so is $\{u, v\}$. This could potentially provide a way to compute a new constant of motion. To prove it, first consider the special case where:

$$\frac{\partial u}{\partial t} = \frac{\partial v}{\partial t} = 0 \tag{4.85}$$

then using the Jacobi identity we have:

$$\{H, \{u, v\}\} = -\{u, \{v, H\}\} - \{v, \{H, u\}\} = -\{u, 0\} - \{v, 0\} = 0$$

$$(4.86)$$

For the most general case we proceed in a similar manner:

$$\{\{u, v\}, H\} = \{u, \{v, H\}\} + \{v, \{H, u\}\} = -\left\{u, \frac{\partial v}{\partial t}\right\} + \left\{v, \frac{\partial u}{\partial t}\right\}$$
$$= -\frac{\partial}{\partial t}\{u, v\} \quad \Rightarrow \quad \frac{d}{dt}\{u, v\} = 0 \tag{4.87}$$

<u>Infinitesimal Canonical Transformations</u>

Let us now consider the generating function:

$$F_2(q, P, t) = q_i P_i + \epsilon G(q, P, t), \qquad (4.88)$$

where $F_2 = q_i P_i$ is an identity transformation, and $|\epsilon| \ll 1$ is infinitesimal. The function G(q, P, t) is known as the generating function of an infinitesimal canonical transformation. Using the properties of an F_2 generating function we have:

$$p_{j} = \frac{\partial F_{2}}{\partial q_{i}} = P_{j} + \epsilon \frac{\partial G}{\partial q_{i}} \quad \Rightarrow \quad \delta p_{j} = P_{j} - p_{j} = -\epsilon \frac{\partial G}{\partial q_{i}}$$

$$(4.89)$$

giving the infinitesimal transformation in the momentum. Likewise:

$$Q_j = \frac{\partial F_2}{\partial P_j} = q_j + \epsilon \frac{\partial G}{\partial P_j} \tag{4.90}$$

Since $P_j = p_j + O(\epsilon)$ and ϵ is infinitesimal we can replace $\partial G(q, P, t)/\partial P_j = \partial G(q, p, t)/\partial p_j + O(\epsilon)$. Therefore we have:

$$Q_j = q_j + \epsilon \frac{\partial G}{\partial p_j} + O(\epsilon^2) \quad \Rightarrow \quad \delta q_j = Q_j - q_j = \epsilon \frac{\partial G}{\partial p_j}$$
(4.91)

where now we note that we can consider G = G(q, p, t), a function of q and p, to this order. Returning to the combined notation of $\vec{\eta}^{\top} = (q_1, \dots, q_n, p_1, \dots, p_n)$, Eq.(4.89) and Eq.(4.90) can be consistly written as the following Poisson bracket:

$$\delta \vec{\eta} = \epsilon \{ \vec{\eta}, G \} \tag{4.92}$$

Example: if $G = p_i$ then $\delta p_i = 0$ and $\delta q_j = \epsilon \delta_{ij}$, which is why momentum is the generator of spatial translations.

Example: if G is the z component of the angular momentum:

$$G = L_z = \sum_{i} (x_i p_{iy} - y_i p_{ix})$$
 and $\epsilon = \delta \theta$ (4.93)

then the infinitesimal change correponds to a rotation

$$\delta x_i = -y_i \delta \theta$$
 , $\delta y_i = x_i \delta \theta$, $\delta z_i = 0$ (4.94)

$$\delta p_{ix} = -p_{iy}\delta\theta$$
 , $\delta p_{iy} = p_{ix}\delta\theta$, $\delta p_{iz} = 0$ (4.95)

which is why angular momentum is the generator of rotations.

Important Example: if G = H and $\epsilon = dt$ then

$$\epsilon\{\vec{\eta},G\} = \{\vec{\eta},H\}dt = \dot{\vec{\eta}}dt = d\vec{\eta}$$

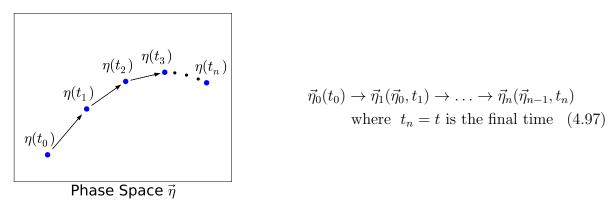
On the left hand side we have the change to the phase space coordinates due to our transformation. On the right hand side we have the physical increment to the phase space variables

that takes place in time dt. The fact that these are equivalent tells us that the Hamiltonian is the generator of time evolution. The infinitesimal transformation generated by the Hamiltonian gives the same function as solving for the physical motion.

Rather than trying to think of this as a passive coordinate change $(q, p) \to (Q, P)$, it is useful if we can take an *active view* of the infinitesimal canonical transformation generated by H. Let the time t be a parameter for the family of transformations with $\epsilon = dt$: the initial conditions are:

$$\vec{\eta}_0(t_0) = \vec{\eta}_0 \tag{4.96}$$

The result is a series of transformations of $\vec{\eta}$ that move us in a fixed set of phase space coordinates from one point to another:



All together, combining an infinite number of infinitesimal transformations allows us to make a finite transformation, resulting in:

$$\vec{\eta} = \vec{\eta}(\vec{\eta}_0, t) \quad \text{or} \quad \vec{\eta}_0 = \vec{\eta}_0(\vec{\eta}, t)$$
(4.98)

This is a canonical transformation that yields a solution for the motion!

How could we directly find this transformation, without resorting to stringing together infinitesimal transformations? We can simply look for a canonical transformation with new coordinates Q_i and new momenta P_i that are all constants, implying an equation of the type:

$$\vec{\eta}_0 = \vec{\eta}_0(\vec{\eta}, t) \tag{4.99}$$

Inverting this then gives the solution for the motion.

This logic can be used to extend our proof of the Theorem in Section 4.2 to fully account for time dependent transformations. (see Goldstein). Using $K = H + \epsilon \partial G / \partial t$, Goldstein also describes in some detail how the change to the Hamiltonian ΔH under an active infinitesimal canonical transformation satisfies:

$$\Delta H = -\epsilon \{G, H\} - \epsilon \frac{\partial G}{\partial t} = -\epsilon \dot{G}$$
 (4.100)

This says "the constants of motion are generating functions G of the infinitesimal canonical transformation that leave H invariant"; that is, $\dot{G} = 0$ if and only if $\Delta H = 0$ under the transformation. Thus a conservation law exists if and only if there is a symmetry present.

4.4 Hamilton-Jacobi Equation

Let us take the suggestion from the end of the previous section seriously and look for new canonical variables that are all cyclic, such that:

$$\dot{Q}_i = \dot{P}_i = 0 \quad \Rightarrow \quad (Q, P) \text{ are all constants}.$$
 (4.101)

If the new Hamiltonian K is independent of (Q, P) then:

$$\dot{Q}_i = \frac{\partial K}{\partial P_i} = 0$$
 and $\dot{P}_i = -\frac{\partial K}{\partial Q_i} = 0$. (4.102)

We could look for a constant K, but it it is simplest to simply look for K=0.

Using a generating function $F = F_2(q, P, t)$, then we need

$$K = H(q, p, t) + \frac{\partial F_2}{\partial t} = 0. \tag{4.103}$$

Because $p_i = \frac{\partial F_2}{\partial q_i}$, then we can rewrite this condition as

$$H\left(q_1, \dots, q_n, \frac{\partial F_2}{\partial q_1}, \dots, \frac{\partial F_2}{\partial q_n}, t\right) + \frac{\partial F_2}{\partial t} = 0, \tag{4.104}$$

which is the time dependent Hamilton-Jacobi equation (henceforth abbreviated as the H-J equation). This is a 1st order partial differential equation in n+1 variables (q_1, \ldots, q_n, t) for F_2 . The solution for F_2 has n+1 independent constants of integration, One of these constants is trivial $(F_2 \to F_2 + C$ for a pure constant C), so we'll ignore this one. Hence, suppose the solution is:

$$F_2 \equiv S = S(q_1, \dots, q_n, \alpha_1, \dots \alpha_n, t), \tag{4.105}$$

where S is called Hamilton's principal function and each α_i is an independent constant. We can pick our new momenta to be the constants of integration $P_i = \alpha_i$ for $i \in \{1, ..., n\}$ (so that $\dot{P}_i = 0$), thus specifying $F_2 = F_2(q, P, t)$ as desired. Then, using again the property of an F_2 generating function (and K = 0), we have that the new constant variables are:

$$P_i \equiv \alpha_i$$
 and $Q_i \equiv \beta_i = \frac{\partial S(q, \alpha, t)}{\partial \alpha_i}$. (4.106)

We introduce the notation β_i to emphasize that these are constants.

From these results we can obtain a solution for the motion as follows. From the invertibility of our transformations we have:

$$\beta_{i}(q,\alpha,t) = \frac{\partial S}{\partial \alpha_{i}} \quad \Rightarrow \quad q_{i} = q_{i}(\alpha,\beta,t),$$

$$p_{i}(q,\alpha,t) = \frac{\partial S}{\partial q_{i}} \quad \Rightarrow \quad p_{i} = p_{i}(q,\alpha,t) = p_{i}(q,\alpha,t),$$

$$p_{i}(q,\alpha,t) = \frac{\partial S}{\partial q_{i}} \quad \Rightarrow \quad p_{i} = p_{i}(q,\alpha,t) = p_{i}(q,\alpha,t),$$

$$(4.107)$$

(Note that function argument notation has been abused slightly here since $p_i(q, \alpha, t)$ and $p_i(\alpha, \beta, t)$ are technically different functions of their three arguments. Since we are always sticking explicit variables into the slots this should not cause confusion.) If desired, we can also swap our 2n constants α_i and β_i for 2n initial conditions q_{i0} and p_{i0} , to obtain a solution for the initial value problem. We obtain one set of constants in terms of the other set by solving the 2n equations obtained from the above results at $t = t_0$:

$$q_{i0} = q_i(\alpha, \beta, t_0), \qquad p_{i0} = p_i(\alpha, \beta, t_0).$$
 (4.108)

Thus we see that Hamilton's principal function S is the generator of canonical transformations of constant (Q, P), and provides a method of obtaining solutions to classical mechanics problems by way of finding a transformation.

There are a few comments to be made about this.

- 1. The choice of constants α_i is somewhat arbitrary, as any other independent choice $\gamma_i = \gamma_i(\alpha)$ is equally good. Thus, when solving the H-J equation, we introduce the constants α_i in whatever way is most convenient.
- 2. What is S? We know that:

$$\dot{S} = \frac{\partial S}{\partial q_i} \dot{q}_i + \frac{\partial S}{\partial P_i} \dot{P}_i + \frac{\partial S}{\partial t}, \tag{4.109}$$

but we also know that:

$$\frac{\partial S}{\partial q_i} = p_i$$
 , $\dot{P}_i = 0$ and $\frac{\partial S}{\partial t} = -H$ (4.110)

Putting Eq.(4.109) and Eq.(4.110) together we have:

$$\dot{S} = p_i \dot{q}_i - H = L \quad \Rightarrow \quad S = \int L \, dt \tag{4.111}$$

Thus S is the classical action which is an indefinite integral over time of the Lagrangian (so it is no coincidence that the same symbol is used).

3. The H-J equation is also the semiclassical limit of the quantum mechanical Schrödinger equation (0'th order term in the WKB approximation). To see this consider the Schrödinger equation

$$i\hbar \frac{\partial \psi}{\partial t} = H\psi = \left(-\frac{\hbar^2}{2m}\frac{\partial^2}{\partial q^2} + V(q)\right)\psi,$$
 (4.112)

with the wavefunction $\psi = \exp(iS/\hbar)$. At this point we are just making a change of variable, without loss of generality, and S(q,t) is complex. Plugging it in, and canceling an exponential after taking the derivative, we find

$$-\frac{\partial S}{\partial t} = -\frac{i\hbar}{2m} \frac{\partial^2 S}{\partial q^2} + \frac{1}{2m} \left(\frac{\partial S}{\partial q}\right)^2 + V(q). \tag{4.113}$$

This equation is just another way of writing the Schrödinger equation, to solve for a complex S instead of ψ . If we now take $\hbar \to 0$ then we find that the imaginary term goes away leaving

$$0 = \frac{\partial S}{\partial t} + \frac{1}{2m} \left(\frac{\partial S}{\partial q}\right)^2 + V(q) = \frac{\partial S}{\partial t} + H\left(q, \frac{\partial S}{\partial q}\right), \tag{4.114}$$

which is the Hamilton-Jacobi equation for S with a standard $p^2/2m$ kinetic term in H.

Having set things up, it is always good for us to test a new formalism on an example where we know the solution.

Example: let us consider the harmonic oscillator Eq.(4.33):

$$H = \frac{1}{2m} \left(p^2 + (m\omega q)^2 \right) = E \tag{4.115}$$

Here we will look for one constant $P = \alpha$ and one constant $Q = \beta$. The H-J equation says

$$\frac{1}{2m} \left(\left(\frac{\partial S}{\partial q} \right)^2 + (m\omega q)^2 \right) + \frac{\partial S}{\partial t} = 0. \tag{4.116}$$

In solving this, we note that the dependence of S on q and t is separable

$$S(q, \alpha, t) = W(q, \alpha) + g(\alpha, t), \tag{4.117}$$

which gives:

$$\frac{1}{2m} \left(\left(\frac{\partial W}{\partial q} \right)^2 + (m\omega q)^2 \right) = -\frac{\partial g}{\partial t} = \alpha.$$
 (4.118)

Since the left side is independent of t and the right hand side is independent of q, then the result must be equal to a separation constant α that is independent of q and t. We will choose our new $P = \alpha$. Now we have

$$\frac{\partial g}{\partial t} = -\alpha \Rightarrow g = -\alpha t \tag{4.119}$$

where we have avoided the addition of a further additive constant (since our convention was to always drop an additive constant when determining S). To identify what α is note that

$$H = -\frac{\partial S}{\partial t} = -\frac{\partial g}{\partial t} = \alpha, \tag{4.120}$$

which corresponds to the constant energy,

$$\alpha = E. \tag{4.121}$$

The other equation we have to solve is

$$\frac{1}{2m} \left(\left(\frac{\partial W}{\partial q} \right)^2 + (m\omega q)^2 \right) = \alpha, \tag{4.122}$$

so rearranging and integrating this gives the indefinite integral

$$W = \pm \int \sqrt{2m\alpha - (m\omega q)^2} \, dq \,, \tag{4.123}$$

which we will leave unintegrated until we must do so. The full solution is then given by:

$$S = -\alpha t \pm \int \sqrt{2m\alpha - (m\omega q)^2} \, dq. \tag{4.124}$$

With this result for Hamilton's Principal function in hand we can now solve for the equations of motion. The equations of motion come from (we now do the integral, after taking the partial derivative):

$$\beta = \frac{\partial S}{\partial \alpha} = -t \pm m \int \frac{dq}{\sqrt{2m\alpha - (m\omega q)^2}} \Rightarrow t + \beta = \pm \frac{1}{\omega} \arcsin\left(\sqrt{\frac{m\omega^2}{2\alpha}}q\right). \tag{4.125}$$

Inverting gives:

$$q = \pm \sqrt{\frac{2\alpha}{m\omega^2}} \sin(\omega(t+\beta)), \tag{4.126}$$

so β is related to the phase. Next we consider p and use this result to obtain:

$$p = \frac{\partial S}{\partial q} = \pm \sqrt{2m\alpha - (m\omega q)^2} = \pm \sqrt{2m\alpha}\cos(\omega(t+\beta))$$
 (4.127)

These results are as expected. We can trade (α, β) for the initial conditions (q_0, p_0) at t = 0. The choice of phase (from shifting β so that $\omega\beta \to \omega\beta + \pi$) allows taking the positive sign of each square root in the solutions above.

Separation of variables is the main technique to solve the H-J equation. In particular, for a time independent H where

$$\dot{H} = \frac{\partial H}{\partial t} = 0 \tag{4.128}$$

we can always separate time by taking:

$$S(q, \alpha, t) = W(q, \alpha) - \alpha_1 t, \tag{4.129}$$

where α_1 has been chosen as the separation constant, then plugging this into the time dependent H-J equation yields (just as in our Harmonic Oscillator example):

$$H\left(q_1,\ldots,q_n,\frac{\partial W}{\partial q_1},\ldots,\frac{\partial W}{\partial q_n}\right) = \alpha_1.$$
 (4.130)

This result is referred to as the *time independent Hamilton-Jacobi equation*. Since $\dot{H}=0$, H is conserved, and equal to a constant α_1 . If H=E then this constant is energy, $\alpha_1=E$. The function W is called Hamilton's characteristic function.

The idea is now to solve the time independent H-J equation for $W = W(q, \alpha)$ where $P = \alpha$ still. If we follow the setup from our time dependent solution above then the equations of motion are obtained from the following prescription for identifying variables:

$$p_{i} = \frac{\partial W}{\partial q_{i}} \text{ for } i \in \{1, \dots, n\},$$

$$Q_{1} = \beta_{1} = \frac{\partial S}{\partial \alpha_{1}} = \frac{\partial W}{\partial \alpha_{1}} - t,$$

$$Q_{j} = \beta_{j} = \frac{\partial W}{\partial \alpha_{j}} \text{ for } j \in \{2, \dots, n\} \text{ for } n > 1.$$

$$(4.131)$$

Here all the Q_i are constants.

There is an alternative to the above setup, which allows us to not refer to the time dependent solution. The alternative is to consider $W = F_2(q, P)$ as the generating function, instead of S and only demand that all the new momenta P_i are constants with $P_1 = \alpha_1 = H$ for a time independent Hamiltonian H. At the start of chapter 4 we saw that this less restrictive scenario would lead to Q_S that could have a linear time dependence, which is still pretty simple.

This is almost identical to the above setup but we rename and reshuffle a few things. The following three equations are the same as before:

$$p_i = \frac{\partial W}{\partial q_i}$$
 , $P_i = \alpha_i$ and $H\left(q, \frac{\partial W}{\partial q}\right) = \alpha_1$ (4.132)

However, now we have a non-zero K and different equation for Q_1 :

$$K = H = \alpha_1$$
 and $Q_i = \frac{\partial W}{\partial \alpha_i}$ for all $i \in \{1, \dots, n\}$. (4.133)

This means:

$$\dot{Q}_1 = \frac{\partial K}{\partial \alpha_1} = 1 \quad \Rightarrow \quad Q_1 = t + \beta_1 = \frac{\partial W}{\partial \alpha_1}$$
 (4.134)

which is Eq. (4.131) but rearranged from the perspective of Q_1 . For j > 1, the equations are the same as before Eq.(4.131):

$$\dot{Q}_j = \frac{\partial K}{\partial \alpha_j} = 0 \quad \Rightarrow \quad Q_j = \beta_j = \frac{\partial W}{\partial \alpha_j}$$
 (4.135)

In this language we do not need to mention S only W. There are a few comments to be made:

1. Again, the choice of α is arbitrary, and $\alpha_i = \alpha_i(\gamma)$ is fine. If we do replace $\alpha_1 = \alpha_1(\gamma)$ then $\dot{Q}_i = \frac{\partial K}{\partial \gamma_i} = v_i$ is a constant so that (potentially) all of the Q_i become linear in time:

$$Q_i = v_i t + \beta_i \text{ for all } i \in \{1, \dots, n\}$$
 (4.136)

2. What is W? We know that:

$$\dot{W} = \frac{\partial W}{\partial q_i} \dot{q}_i = p_i \dot{q}_i \Rightarrow W = \int p_i \dot{q}_i \, dt = \int p_i \, dq_i, \tag{4.137}$$

which is a different sort of "action".

3. The time independent H-J equation has some similarity to the time-independent Schrödinger energy eigenvalue equation (both involve H and and constant E, but the former is a non-linear equation for W, while the latter is a linear equation for the wavefunction ψ).

The H-J method is most useful when there is a separation of variables in H.

Example: if

$$H = h_1(q_1, q_2, p_1, p_2) + h_2(q_1, q_2, p_1, p_2) f(q_3, p_3) = \alpha_1, \tag{4.138}$$

so that q_3 is separable, then

$$f(q_3, p_3) = \frac{\alpha_1 - h_1}{h_2} \tag{4.139}$$

is a constant because the right hand side is independent of q_3 and p_3 . Thus we assign

$$f(q_3, p_3) = \alpha_2 (4.140)$$

for convenience. We can then write:

$$W = W'(q_1, q_2, \alpha) + W_3(q_3, \alpha) \quad \Rightarrow \quad f\left(q_3, \frac{\partial W_3}{\partial q_3}\right) = \alpha_2 \quad \text{and}$$
 (4.141)

$$h_1\left(q_1, q_2, \frac{\partial W'}{\partial q_1}, \frac{\partial W'}{\partial q_2}\right) + \alpha_2 h_2\left(q_1, q_2, \frac{\partial W'}{\partial q_1}, \frac{\partial W'}{\partial q_2}\right) = \alpha_1 \tag{4.142}$$

Here, q_1 and q_2 may or may not be separable.

If all variables are separable then we use the solution:

$$W = \sum_{i} W_i(q_i, \alpha) \tag{4.143}$$

We can simply try a solution of this form to test for separability.

Note that cyclic coordinates are always separable.

Proof: let us say that q_1 is cyclic. Then

$$p_1 \equiv \gamma$$
 and $H\left(q_2, \dots, q_n, \frac{\partial W}{\partial q_1}, \frac{\partial W}{\partial q_2}, \dots, \frac{\partial W}{\partial q_n}\right) = \alpha_1,$ (4.144)

where γ is constant. Let us now write

$$W(q,\alpha) = W_1(q_1,\alpha) + W'(q_2,\dots,q_n,\alpha). \tag{4.145}$$

This gives us:

$$p_1 = \frac{\partial W_1}{\partial q_1} = \gamma \Rightarrow W_1 = \gamma q_1. \tag{4.146}$$

Which gives us:

$$W(q,\alpha) = \gamma q_1 + W'(q_2, \dots, q_n, \alpha)$$
(4.147)

This procedure can be repeated for any remaining cyclic variables.

Note that the choice of variables is often important in finding a result that separates. A problem with spherical symmetry may separate in spherical coordinates but not Cartesian coordinates.

4.5 Kepler Problem

As an extended example, let us consider the Kepler problem of two masses m_1 and m_2 in a central potential (with the CM coordinate $\mathbf{R} = 0$). The Lagrangian is:

$$L = \frac{1}{2}m\dot{\mathbf{r}}^2 - V(r)$$
 where $\frac{1}{m} \equiv \frac{1}{m_1} + \frac{1}{m_2}$, (4.148)

and here m is the reduced mass. Any V(r) conserves $\mathbf{L} = \mathbf{r} \times \mathbf{p}$, so the motion of \mathbf{r} and \mathbf{p} is in a plane perpendicular to \mathbf{L} . The coordinates in the plane can be taken as (r, ψ) , so:

$$L = \frac{m}{2} \left(\dot{r}^2 + r^2 \dot{\psi}^2 \right) - V(r), \tag{4.149}$$

with ψ being cyclic, which implies:

$$p_{\psi} = mr^2 \dot{\psi}$$
 is a constant. (4.150)

In fact $p_{\psi} = |\mathbf{L}| \equiv \ell$. Notationally, we use ℓ for the magnitude of the angular momentum \mathbf{L} to distinguish it from the Lagrangian L.

The energy is then:

$$E = \frac{m}{2}\dot{r}^2 + \frac{\ell^2}{2mr^2} + V(r), \tag{4.151}$$

which is constant, and this can be rewritten as:

$$E = \frac{m}{2}\dot{r}^2 + V_{\text{eff}}(r) \quad \text{where} \quad V_{\text{eff}}(r) = V(r) + \frac{\ell^2}{2mr^2},$$
 (4.152)

where V_{eff} is the effective potential, as plotted below for the gravitational potential.

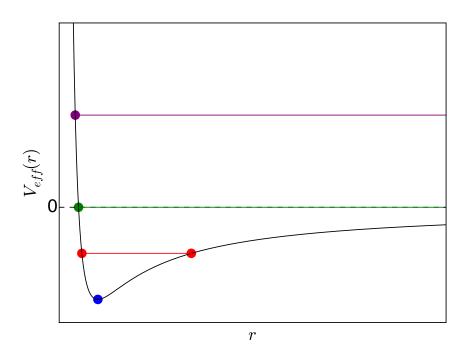


Figure 4.1: Plot of the effective potential V_{eff} along with the different qualitative orbits allowed in a gravity-like potential. The points correspond to turning points of the orbit.

Writing the E-L equation for $\dot{r} = dr/dt = \dots$ and then solving for it as $dt = dr/(\dots)$, and integrating yields

$$t = t(r) = \int_{r_0}^{r} \frac{dr'}{\sqrt{\frac{2}{m} \left(E - V(r') - \frac{\ell^2}{2mr'^2}\right)}}$$
(4.153)

as an implicit solution to the radial motion.

The orbital motion comes as $r = r(\psi)$ or $\psi = \psi(r)$ by using Eq.(4.150) and substituting, in Eq.(4.153). We have $\dot{\psi} = d\psi/dt = \ell/(mr^2)$, so we can use this to replace dt by $d\psi$ in dt = dr/(...) to get an equation of the form $d\psi = dr/(...)$. The result is given by

$$\psi - \psi_0 = \ell \int_{r_0}^r \frac{dr'}{r'^2 \sqrt{2m \left(E - V(r') - \frac{\ell^2}{2mr'^2}\right)}}$$
(4.154)

In the particular case of $V(r) = -\frac{k}{r}$, the solution of the orbital equation is:

$$\frac{1}{r(\psi)} = \frac{mk}{\ell^2} \left(1 + \varepsilon \cos(\psi - \psi') \right) \tag{4.155}$$

where the eccentricity ε is given by:

$$\varepsilon \equiv \sqrt{1 + \frac{2E\ell^2}{mk^2}} \tag{4.156}$$

Below are plotted the different qualitative orbits for the gravitic potential, with different ε or E (circular, elliptical, parabolic, and hyperbolic respectively).

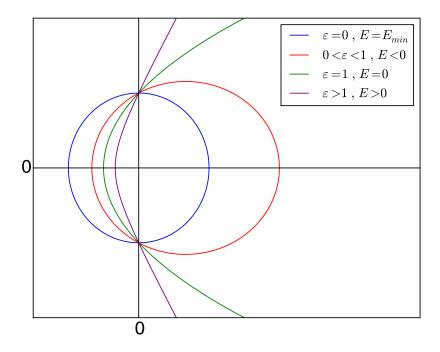


Figure 4.2: Different Orbits for the gravity-like potential. The orbits' colors match those of Fig.(4.1). The unbounded orbits occur for $E \geq 0$. The different curves correspond to the different possible conic sections.

Consider solving this problem instead by the H-J method. Lets start by considering as the variables (r, ψ) so that we assume that the motion of the orbit is in a plane. Here

$$H = \frac{1}{2m} \left(p_r^2 + \frac{p_\psi^2}{r^2} \right) + V(r) = \alpha_1 = E.$$
 (4.157)

As ψ is cyclic, then $p_{\psi} \equiv \alpha_{\psi}$ is constant. Using:

$$W = W_1(r) + \alpha_{\psi}\psi, \tag{4.158}$$

then the time independent H-J equation is:

$$\frac{1}{2m} \left(\left(\frac{\partial W_1}{\partial r} \right)^2 + \frac{\alpha_{\psi}^2}{r^2} \right) + V(r) = \alpha_1. \tag{4.159}$$

This is simplified to

$$\frac{\partial W_1}{\partial r} = \sqrt{2m(\alpha_1 - V(r)) - \frac{\alpha_\psi^2}{r^2}} \tag{4.160}$$

and solved by

$$W = \alpha_{\psi}\psi + \int \sqrt{2m(\alpha_1 - V(r)) - \frac{\alpha_{\psi}^2}{r^2}} dr.$$
 (4.161)

The transformation equations are:

$$t + \beta_1 = \frac{\partial W}{\partial \alpha_1} = m \int \frac{dr}{\sqrt{2m(\alpha_1 - V(r)) - \frac{\alpha_\psi^2}{r^2}}},$$

$$\beta_2 = \frac{\partial W}{\partial \alpha_\psi} = \psi - \alpha_\psi \int \frac{dr}{r^2 \sqrt{2m(\alpha_1 - V(r)) - \frac{\alpha_\psi^2}{r^2}}}.$$
(4.162)

Thus we immediately get the radial equation t = t(r) and orbital equation $\psi = \psi(r)$ from this, with $\alpha_{\psi} = \ell$ and $\alpha_1 = E$, showing that the constants are physically relevant parameters.

Let's solve this problem again, but suppose the motion is in 3 dimensions (as if we did not know the plane of the orbit). Using spherical coordinates (r, θ, φ) this corresponds to

$$H = \frac{1}{2m} \left(p_r^2 + \frac{p_\theta^2}{r^2} + \frac{p_\varphi^2}{r^2 \sin^2 \theta} \right) + V(r) = \alpha_1.$$
 (4.163)

Lets try a separable solution

$$W = W_r(r) + W_{\theta}(\theta) + W_{\varphi}(\varphi). \tag{4.164}$$

Since φ is cyclic we know it is separable and that:

$$W_{\varphi}(\varphi) = \alpha_{\varphi} \, \varphi \,. \tag{4.165}$$

Together, this leaves;

$$\left(\frac{\partial W_r}{\partial r}\right)^2 + \frac{1}{r^2} \left(\left(\frac{\partial W_\theta}{\partial \theta}\right)^2 + \frac{\alpha_\varphi^2}{\sin^2(\theta)}\right) + 2mV(r) = 2m\alpha_1.$$
(4.166)

Because the term:

$$\left(\frac{\partial W_{\theta}}{\partial \theta}\right)^2 + \frac{\alpha_{\varphi}^2}{\sin^2(\theta)} \tag{4.167}$$

only depends on θ while the rest of the equation depends on r, it must be a constant so we can say:

$$\left(\frac{\partial W_{\theta}}{\partial \theta}\right)^2 + \frac{\alpha_{\varphi}^2}{\sin^2(\theta)} \equiv \alpha_{\theta}^2 \tag{4.168}$$

and the separation works. This then gives:

$$\left(\frac{\partial W_r}{\partial r}\right)^2 + \frac{\alpha_\theta^2}{r^2} = 2m(\alpha_1 - V(r)), \tag{4.169}$$

which is the same equation we considered before when assuming the motion was in a plane, with $\alpha_1 = E$ and $\alpha_\theta = \ell$. Eq. (4.168) says that

$$p_{\theta}^{2} + \frac{p_{\varphi}^{2}}{\sin^{2}(\theta)} = \ell^{2}. \tag{4.170}$$

Here p_{φ} is the constant angular momentum about the \hat{z} axis.

4.6 Action-Angle Variables

For many problems, we may not be able to solve analytically for the exact motion or for orbital equations, but we can still characterize the motion. For *periodic* systems we can find the frequency by exploiting *action-angle variables*.

The simplest case is for a single dimension of canonical coordinates (q, p). If $H(q, p) = \alpha_1$ then $p = p(q, \alpha_1)$. There are two types of periodic motion to consider.

1. Libration (oscillation) is characterized by a closed phase space orbit, so that q and p evolve periodically with the same frequency.

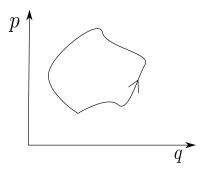


Figure 4.3: Phase space orbit of a libration (oscilation). The trajectory closes on itself, the state returns to the same position after some time τ .

2. Rotation is characterized by an open phase space path, so p is periodic while q evolves without bound.

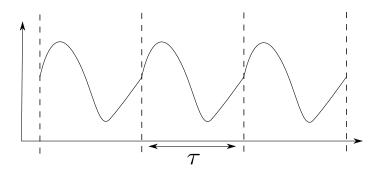


Figure 4.4: Phase space orbit of a rotation. Although the orbit is not closed, each period the evolution of the system is the same, leading to a orbit that repeats itself with a translation.

Example: a pendulum of length a may be characterized by canonical coordinates (θ, p_{θ}) , where:

$$E = H = \frac{p_{\theta}^2}{2ma^2} - mga\cos\theta \tag{4.171}$$

This means:

$$p_{\theta} = \pm \sqrt{2ma^2(E + mga\cos\theta)} \tag{4.172}$$

must be real. A rotation occurs when E > mga, and oscillations occur when E < mga. The critical point in between (when the pendulum just makes it to the top) is when E = mga exactly, and is depicted by a dashed line in the figure below.

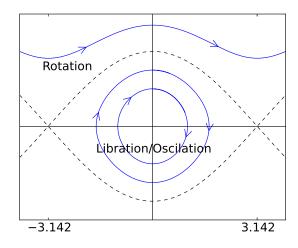


Figure 4.5: The pendulum exhibits both librations and rotations depending on the initial conditions.

For either type of periodic motion, it is useful to replace $P = \alpha_1$ by a different (constant) choice called the *action variable*

$$J = \oint p \, dq,\tag{4.173}$$

where \oint refers to a definite integral over one period in phase space. To see that J is constant, recall that $p = p(q, \alpha_1)$, so plugging this into the definite integral we are left with $J = J(\alpha_1)$. Also, we have the inverse, $\alpha_1 = H = H(J)$, and can rewrite Hamilton's characteristic function in terms of J by $W = W(q, \alpha_1) = W(q, H(J)) = W(q, J)$ (where again the argument notation is abused slightly).

The coordinate conjugate to J is the angle variable

$$\omega = \frac{\partial W}{\partial J} \tag{4.174}$$

(where ω is not meant to imply an angular velocity). This means

$$\dot{\omega} = \frac{\partial H(J)}{\partial J} = \nu(J) \text{ is a constant.}$$
 (4.175)

As a result the angle variable has linear time dependence,

$$\omega = \nu t + \beta \,, \tag{4.176}$$

for some initial condition β . Dimensionally, J has units of angular momentum, while ω has no dimensions (like an angle or a phase).

To see why it is useful to use the canonical variables (ω, J) , let us consider the change in ω when q goes through a complete cycle.

$$\Delta\omega = \oint \frac{\partial\omega}{\partial q} \, dq = \oint \frac{\partial^2 W}{\partial q \partial J} \, dq = \frac{\partial}{\partial J} \oint \frac{\partial W}{\partial q} \, dq = \frac{\partial}{\partial J} \oint p \, dq = 1 \tag{4.177}$$

where in the last equality we used the definition of J in Eq.(4.173). Also, we have $\Delta\omega = \nu\tau$ where τ is the period. Thus

$$\nu = \frac{1}{\tau} \tag{4.178}$$

is the frequency of periodic motion. If we find H = H(J) then

$$\nu = \frac{\partial H(J)}{\partial J} \tag{4.179}$$

immediately gives the frequency $\nu = \nu(J)$ for the system. Often, we then J = J(E) to get $\nu = \nu(E)$ the frequency at a given energy. This is a very efficient way of finding the frequency of the motion without solving for extraneous information.

Example: let us consider a pendulum with action-angle variables. We define:

$$\tilde{E} = \frac{E}{mga} \tag{4.180}$$

so that $\tilde{E} > 1$ corresponds to rotation and $\tilde{E} < 1$ corresponds to oscillation. This means

$$p_{\theta} = \pm \sqrt{2m^2 g a^3} \sqrt{\tilde{E} + \cos \theta}. \tag{4.181}$$

For $\tilde{E} > 1$:

$$J = \sqrt{2m^2ga^3} \int_{-\pi}^{\pi} d\theta \sqrt{\tilde{E} + \cos\theta}, \qquad (4.182)$$

which is an elliptic integral. For $\tilde{E} < 1$:

$$J = \sqrt{2m^2ga^3} \int_{-\theta_0}^{+\theta_0} d\theta \sqrt{\tilde{E} + \cos\theta} + \sqrt{2m^2ga^3} \int_{\theta_0}^{-\theta_0} d\theta \left[-\sqrt{\tilde{E} + \cos\theta} \right]$$
$$= 4\sqrt{2m^2ga^3} \int_0^{\theta_0} d\theta \sqrt{\tilde{E} + \cos\theta} , \qquad (4.183)$$

as the contributions from the four intervals that the pendulum swings through in one period are all equivalent. Here θ_0 is the turning point of the oscillation, and $\tilde{E} = -\cos(\theta_0)$.

From this:

$$\nu = \frac{\partial E}{\partial J} = \left(\frac{\partial J}{\partial E}\right)^{-1} \tag{4.184}$$

which we can solve graphically by making a plot of J vs E, then dJ/dE versus E, and finally the inverse $\nu = dE/dJ$ versus E.

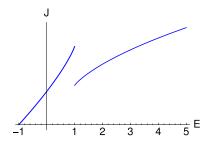


Figure 4.6: Plot of J(E) versus \tilde{E} . The discontinuity corresponds to the transition from Oscilation to Rotation.

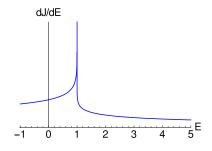


Figure 4.7: Plot of $\frac{dJ}{dE}$. The discontinuity is logarithmic divergent so it is integrable.

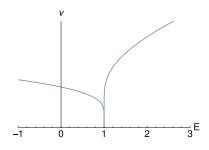


Figure 4.8: Plot of the frequency of oscillation $\nu(E)$ versus \tilde{E} . As $\tilde{E} \to -1$ we approach the small amplitude limit, where $\nu = (2\pi)^{-1} \sqrt{g/a}$.

Example: let us consider the limit $|\theta| \ll 1$ of a pendulum, so:

$$H = \frac{p_{\theta}^2}{2ma^2} + \frac{mga}{2}\theta^2 - mga \tag{4.185}$$

We can actually consider this in the context of a general harmonic oscillator

$$H = \frac{p^2}{2m} + \frac{m\Omega^2}{2}x^2,\tag{4.186}$$

where:

$$\Omega = \sqrt{\frac{g}{a}}$$
 , $x = a\theta$ and $p = \frac{p_{\theta}}{a}$ (4.187)

Notationally, Ω is used for the harmonic oscillator frequency to distinguish from the transformed angle variable ω . We then have:

$$J = \oint p \, dq = \oint \pm \sqrt{2mE - m^2 \Omega^2 x^2} \, dx \tag{4.188}$$

Note that the coordinate does not need to be an angle, as may be the case for general x. This gives:

$$J = 4\sqrt{2mE} \int_0^{x_0} \sqrt{1 - \frac{x^2}{x_0^2}} dx$$
 where $x_0 \equiv \frac{\sqrt{2mE}}{m\Omega}$. (4.189)

Solving the integral yields

$$J = \pi \sqrt{2mE}x_0 = \frac{2\pi mE}{m\Omega} = \frac{2\pi E}{\Omega},\tag{4.190}$$

which gives us

$$\frac{\partial E}{\partial J} = \frac{\Omega}{2\pi},\tag{4.191}$$

the expected cyclic frequency for the harmonic oscillator.

Multiple Variables: We can treat multiple variables $(q_1, \ldots, q_n, p_1, \ldots, p_n)$ with the actionangle formalism if each pair (q_i, p_i) has an oscillatory or rotating orbit. Lets also assume that the H-J equation is completely separable into:

$$W = \sum_{j} W_j(q_j, \alpha). \tag{4.192}$$

Here we have

$$p_i = \frac{\partial W_i}{\partial q_i} = p_i(q_i, \alpha_1, \dots, \alpha_n) \Rightarrow J_i = \oint p_i \, dq_i = J_i(\alpha_1, \dots, \alpha_n)$$
 (4.193)

where repeated indices do not correspond to implicit sums here. This implies that the inverse will be $\alpha_j = \alpha_j(J_1, \ldots, J_n)$ and thus $\alpha_1 = H = H(J_1, \ldots, J_n)$. Likewise:

$$\omega_i = \frac{\partial W}{\partial J_i} = \sum_j \frac{\partial W_j}{\partial J_i} = \omega_i(q_1, \dots, q_n, J_1, \dots, J_n). \tag{4.194}$$

Just as in the one dimensional case the time derivative of the angle variables is a constant

$$\dot{\omega}_i = \frac{\partial H}{\partial J_i} = \nu_i(J_1, \dots, J_n) \tag{4.195}$$

which are the frequencies describing motion in this "multi-periodic" system. Due to the presence of multiple frequencies, the motion through the whole 2n-dimensional phase space need not be periodic in time.

Example: in the 2-dimensional harmonic oscillator:

$$x = A\cos(2\pi\nu_1 t) \text{ and } y = B\cos(2\pi\nu_2 t)$$
 (4.196)

$$p_x = m\dot{x} \text{ and } p_y = m\dot{y} \tag{4.197}$$

The overall motion is not periodic in time unless $\frac{\nu_1}{\nu_2}$ is a rational number.

Kepler Problem Example:

Let us do a more extended and detailed example. Returning to the Kepler problem:

$$V(r) = -\frac{k}{r} \tag{4.198}$$

with its separable W:

$$W = W_r(r,\alpha) + W_\theta(\theta,\alpha) + W_\varphi(\varphi,\alpha). \tag{4.199}$$

If we take E < 0, we have oscillation in r and θ , along with a rotation in φ . In particular from solving our earlier differential equations for W_{θ} and W_{r} , we have

$$W_{\varphi} = \alpha_{\varphi} \varphi$$

$$W_{\theta} = \pm \int d\theta \sqrt{\alpha_{\theta}^2 - \frac{\alpha_{\varphi}^2}{\sin^2(\theta)}}$$

$$W_r = \pm \int dr \sqrt{2m(\alpha_1 - V(r)) - \frac{\alpha_{\theta}^2}{r^2}}$$

Here we have

$$J_{\varphi} = \oint p_{\varphi} \, d\varphi = \oint \frac{\partial W}{\partial \varphi} \, d\varphi = \oint \alpha_{\varphi} \, d\varphi \tag{4.200}$$

For the cyclic variable φ , we still call the constant p_{φ} periodic and will take the period to be 2π (arbitrarily since any period would work), which corresponds to particle returning to the original point in space. Thus

$$J_{\varphi} = 2\pi\alpha_{\varphi},\tag{4.201}$$

where α_{φ} is the angular momentum about \hat{z} .

Continuing, in a similar manner we have

$$J_{\theta} = \oint p_{\theta} d\theta = \oint \frac{\partial W}{\partial \theta} d\theta = \oint \pm \sqrt{\alpha_{\theta}^2 - \frac{\alpha_{\varphi}^2}{\sin^2(\theta)}} d\theta \tag{4.202}$$

Let us call:

$$\cos(\gamma) \equiv \frac{\alpha_{\varphi}}{\alpha_{\theta}} \tag{4.203}$$

the angular momentum fraction. Then:

$$J_{\theta} = \alpha_{\theta} \oint \pm \sqrt{1 - \frac{\cos^2(\gamma)}{\sin^2(\theta)}} \, d\theta \tag{4.204}$$

If we let $\sin(\theta_0) = \cos(\gamma)$, then $p_{\theta} = 0$ at the turning points, $\theta \in \{\theta_0, \pi - \theta_0\}$, as expected. Here one oscillator goes from $\pi - \theta_0 \to \theta_0$ when $p_{\theta} > 0$, and in reverse for $p_{\theta} < 0$. Moreover, $\sin(\theta)^{-2}$ is even about $\theta = \frac{\pi}{2}$. This gives

$$J_{\theta} = 4\alpha_{\theta} \int_{\frac{\pi}{2}}^{\theta_0} \sqrt{1 - \frac{\cos^2(\gamma)}{\sin^2(\theta)}} d\theta. \tag{4.205}$$

Making two more substitutions

$$\cos(\theta) \equiv \sin(\gamma)\sin(\psi)$$
, and then $u \equiv \tan(\psi)$, (4.206)

after some work the expression becomes

$$J_{\theta} = 4\alpha_{\theta} \int_{0}^{\infty} \left(\frac{1}{1+u^{2}} - \frac{\cos^{2}(\gamma)}{1+u^{2}\cos^{2}(\gamma)} \right) du = 2\pi\alpha_{\theta} (1-\cos(\gamma))$$
$$= 2\pi(\alpha_{\theta} - \alpha_{\varphi}), \tag{4.207}$$

which gives

$$J_{\theta} + J_{\varphi} = 2\pi\alpha_{\theta}. \tag{4.208}$$

Finally we can consider

$$J_r = \oint \sqrt{2mE - 2mV(r) - \frac{(J_\theta + J_\varphi)^2}{4\pi^2 r^2}} dr$$
 (4.209)

We can immediately make some observations. We observe that $J_r = J_r(E, J_\theta + J_\varphi)$ is a function of two variables for any V = V(r), and thus if we invert $E = E(J_r, J_\theta + J_\varphi)$. This implies:

$$\frac{\partial E}{\partial J_{\theta}} = \frac{\partial E}{\partial J_{\varphi}} \quad \Rightarrow \quad \nu_{\theta} = \nu_{\varphi} \tag{4.210}$$

The two frequencies are degenerate for any V = V(r).

For the $V(r) = -kr^{-1}$ potential, the integration can be performed (for example, by contour integration) to give (for E < 0)

$$J_r = -(J_\theta + J_\varphi) + \pi k \sqrt{\frac{2m}{-E}} \Rightarrow J_r + J_\theta + J_\varphi = \pi k \sqrt{\frac{2m}{-E}}.$$
 (4.211)

This means:

$$E = -\frac{2\pi^2 k^2 m}{(J_r + J_\theta + J_\varphi)^2} \quad \Rightarrow \quad \nu_\theta = \nu_\varphi = \nu_r \tag{4.212}$$

In particular:

$$\nu_r = \frac{\partial E}{\partial J_r} = 4\pi^2 k^2 (J_r + J_\theta + J_\varphi)^{-3} = \frac{1}{\pi k} \sqrt{\frac{-2E^3}{m}}$$
(4.213)

which is the correct orbital frequency in a bound Kepler orbit.

Using the relations between $\{\alpha_1 = E, \alpha_{\theta}, \alpha_{\varphi}\}\$ and $\{J_r, J_{\theta}, J_{\varphi}\}\$, we can also get Hamilton's characteristic function for this system as

$$W = W_{\varphi} + W_{\theta} + W_{r}$$

$$= \frac{\varphi J_{\varphi}}{2\pi} \pm \int \sqrt{(J_{\theta} + J_{\varphi})^{2} - \frac{J_{\varphi}^{2}}{\sin^{2}(\theta)}} \frac{d\theta}{2\pi} \pm \int \sqrt{\frac{-(2\pi mk)^{2}}{(J_{r} + J_{\theta} + J_{\varphi})^{2}} + \frac{2mk}{r} - \frac{(J_{\theta} + J_{\varphi})^{2}}{(2\pi r)^{2}}} dr.$$

This then gives the angle variables:

$$\omega_{r} = \frac{\partial W}{\partial J_{r}} = \omega_{r}(r, J_{r} + J_{\theta} + J_{\varphi}, J_{\theta} + J_{\varphi})$$

$$\omega_{\theta} = \frac{\partial W}{\partial J_{\theta}} = \omega_{\theta}(r, \theta, J_{r} + J_{\theta} + J_{\varphi}, J_{\theta} + J_{\varphi}, J_{\varphi})$$

$$\omega_{\varphi} = \frac{\partial W}{\partial J_{\varphi}} = \omega_{\varphi}(r, \theta, \varphi, J_{r} + J_{\theta} + J_{\varphi}, J_{\theta} + J_{\varphi}, J_{\varphi})$$

$$(4.214)$$

where $\dot{\omega}_r = \nu_r$, $\dot{\omega}_{\theta} = \nu_{\theta}$, and $\dot{\omega}_{\varphi} = \nu_{\varphi}$. Of course, in this case, $\nu_r = \nu_{\theta} = \nu_{\varphi}$.

At this point we can identify five constants for the Kepler problem

$$J_{1} = J_{\varphi}$$

$$J_{2} = J_{\theta} + J_{\varphi}$$

$$J_{3} = J_{r} + J_{\theta} + J_{\varphi}$$

$$\omega_{1} = \omega_{\varphi} - \omega_{\theta}$$

$$\omega_{2} = \omega_{\theta} - \omega_{r}.$$

$$(4.215)$$

(These 5 constants could also be identified from the angular momentum \vec{L} , energy E, and Laplace-Runge-Lenz vector \vec{A} .) What are they? There are two constants specifying the plane of the orbit (the x'y'-plane), which are the inclination angle i and the longitude of the ascending node Ω . There are three constants specifying the form of the ellipse, which are the semi-major axis a (giving the size), the eccentricity ε (giving the shape), and the angle ω (giving the orientation within the plane). These are all shown in Fig. 4.9.

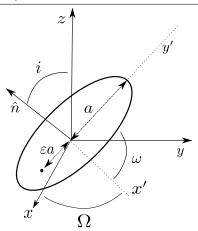


Figure 4.9: Picture of an orbit in 3d and the five parameters necessary to fully specify it. The angles i, Ω and ω provide the orientation in space while a and ε provide the size and shape of the conic section.

It can be shown that the relations between these constants and the ones above are

$$\cos(i) = \frac{J_1}{J_2} \qquad a = -\frac{k}{2E} = \frac{J_3^2}{4\pi^2 mk} \qquad \epsilon = \sqrt{1 - \left(\frac{J_2}{J_3}\right)^2}$$

$$\Omega = 2\pi\omega_1 \qquad \omega = 2\pi\omega_2 \qquad (4.216)$$

providing a fairly simple physical interpretations to $(J_1, J_2, J_3, \omega_1, \omega_2)$. Also recall that $J_2 = 2\pi\alpha_{\theta} = 2\pi\ell$.

When the orbit is perturbed by additional forces (like a moon, another planet, general relativistic corrections, and so on), these action-angle variables provide a natural way to describe the modified orbit. We will soon see that they become functions of time. For example, from the above list of constants, the perturbed $\omega = \omega(t)$ is the precession of the perihelion of an orbit. We will learn how to derive this time dependence next, in our chapter on Perturbation Theory.

Chapter 5

Perturbation Theory

In this chapter we will discuss time dependent perturbation theory in classical mechanics. Many problems we have encountered yield equations of motion that cannot be solved analytically. Here, we will consider cases where the problem we want to solve with Hamiltonian H(q, p, t) is "close" to a problem with Hamiltonian $H_0(q, p, t)$ for which we know the exact solution. We say

$$H(q, p, t) = H_0(q, p, t) + \Delta H(q, p, t),$$
 (5.1)

where ΔH is small. The general idea is to expand variables

$$z(t) = z_0(t) + \epsilon z_1(t) + \epsilon^2 z_2(t) + \dots,$$
 (5.2)

for $z \in \{q, p\}$ and use the expanded equations of motion to determine the series

$$z_0(t) \to z^{(1)}(t) \to z^{(2)}(t) \to \dots$$
, where $z^{(k)}(t) = \sum_{j=0}^k \epsilon^j z_j(t)$. (5.3)

We can do this with any of our methods for solving problems in classical mechanics, including the Euler-Lagrange equations, Hamilton equations, Poisson bracket equations, or Hamilton-Jacobi equations. Since there are some practical benefits, our focus will be on doing this for the Hamilton-Jacobi equations, but lets first start with an example where we carry out an expansion for the Hamilton equations.

Example Consider $H_0 = \frac{p^2}{2m}$ the free Hamiltonian, and $\Delta H = \frac{m\omega^2}{2}x^2$. Here ω is an oscillator frequency. The full Hamiltonian $H = H_0 + \Delta H$ in this case is just a Harmonic oscillator where we already know the solution, so we have the opportunity to see how this solution is built up perturbatively. Without any approximation, the Hamilton equations are

$$\dot{x} = \frac{p}{m} \,, \qquad \qquad \dot{p} = -m\omega^2 x \,. \tag{5.4}$$

To carry out perturbation theory we are going to count w^2 as $\mathcal{O}(\epsilon)$ and then at each order we balance the number of ϵ s on each side of the equations of motion. For H_0 , we have

 $\dot{p}_0 = 0$ so the momentum p_0 is a constant which we fix as the initial condition value. We also have $\dot{x}_0 = \frac{p_0}{m}$ (since w^2 does not appear we have not dropped anything in this equation). Integrating we get $x_0 = \frac{p_0}{m}t$, where we have taken the initial condition x(t=0) = 0 for simplicity.

Having setup the 0'th order solution, lets now consider determining the solution at 1'st order. At first order the RHS of the equations of motion should be $\mathcal{O}(\epsilon)$. Therefore

$$\dot{p}^{(1)} = -m\omega^2 x^{(0)} = -m\omega^2 \frac{p_0}{m} t = -\omega^2 t \, p_0,$$

$$p^{(1)}(t) = p_0 - \frac{1}{2} p_0 \omega^2 t^2.$$
(5.5)

For the other equation of motion at this order we then have

$$\dot{x}^{(1)} = \frac{p^{(1)}}{m} = \frac{p_0}{m} - \frac{p_0 \omega^2 t^2}{2m},$$

$$x^{(1)}(t) = \frac{p_0}{m} t - \frac{p_0 \omega^2 t^3}{6m}.$$
(5.6)

These are precisely the 1st order terms in the full solution

$$p(t) = p_0 \cos(\omega t), \qquad x(t) = \frac{p_0}{m\omega} \sin(\omega t).$$
 (5.7)

5.1 Time Dependent Perturbation Theory for the Hamilton-Jacobi Equations

From here on we will focus on using H-J methods. If $H = H_0 + \Delta H$, then the solution for H_0 has a principal function $S(q, \alpha, t)$ that is the generating function that makes a canonical transformation $(q, p) \to (\alpha, \beta)$, so that

$$H_0\left(q, \frac{\partial S}{\partial q}, t\right) + \frac{\partial S}{\partial t} = 0.$$
 (5.8)

For the dynamics generated by H_0 the variables (α, β) are constants. However, the resulting canonical transformation provides a new set of variables that is valid for use with any Hamiltonian, they are just particularly simple variables for H_0 . Therefore, for H, we can still use the canonical transformation generated by S, but now the new variables

$$P_i = \alpha_i = \alpha_i(p, q), \qquad Q_i = \beta_i = \beta_i(p, q), \qquad (5.9)$$

will no longer be constant in time. The new Hamiltonian is

$$K = H_0 + \Delta H + \frac{\partial S}{\partial t} = \Delta H = \Delta H(\alpha, \beta, t)$$
 (5.10)

The new Hamilton equations $\dot{Q}_i = \frac{\partial K}{\partial P_i}$ and $\dot{P}_i = -\frac{\partial K}{\partial Q_i}$ now yield exact equations of motion for these variables

$$\dot{\alpha}_i = -\frac{\partial \Delta H}{\partial \beta_i}, \qquad \qquad \dot{\beta}_i = \frac{\partial \Delta H}{\partial \alpha_i}.$$
 (5.11)

The idea of perturbation theory is to solve these equations with an expansion. Since here the small $\Delta H \sim \epsilon$ appears on the RHS of both equations, we will always use lower order solutions on the RHS to obtain the higher order results on the LHS. Thus we use $\alpha^{(0)}$ and $\beta^{(0)}$ to get the first order $\alpha^{(1)}$ and $\beta^{(1)}$:

$$\dot{\alpha}_{i}^{(1)} = -\frac{\partial \Delta H}{\partial \beta_{i}} \bigg|_{\substack{\alpha_{i} = \alpha_{i}^{(0)} \\ \beta_{i} = \beta_{i}^{(0)}}} = -\frac{\partial \Delta H}{\partial \beta_{i}} \bigg|_{0}, \qquad (5.12)$$

$$\dot{\beta}_{i}^{(1)} = \frac{\partial \Delta H}{\partial \alpha_{i}} \bigg|_{\substack{\alpha_{i} = \alpha_{i}^{(0)} \\ \beta_{i} = \beta_{i}^{(0)}}} = \frac{\partial \Delta H}{\partial \alpha_{i}} \bigg|_{0}, \qquad (5.12)$$

where the $|_0$ is a shorthand notation. We then use $\alpha^{(1)}$ and $\beta^{(1)}$ to get $\alpha^{(2)}$ and $\beta^{(2)}$, and so on. At nth order we have

$$\dot{\alpha}_{i}^{(n)} = -\frac{\partial \Delta H}{\partial \beta_{i}} \bigg|_{\substack{\alpha_{i} = \alpha_{i}^{(n-1)} \\ \beta_{i} = \beta_{i}^{(n-1)}}} = -\frac{\partial \Delta H}{\partial \beta_{i}} \bigg|_{n-1}, \qquad (5.13)$$

$$\dot{\beta}_{i}^{(n)} = \frac{\partial \Delta H}{\partial \alpha_{i}} \bigg|_{\substack{\alpha_{i} = \alpha_{i}^{(n-1)} \\ \beta_{i} = \beta_{i}^{(n-1)}}} = \frac{\partial \Delta H}{\partial \alpha_{i}} \bigg|_{n-1}.$$

Example Lets once again consider $H_0 = \frac{p^2}{2m}$ and $\Delta H = \frac{m\omega^2}{2}x^2$. For H_0 , the H-J equation is $\frac{1}{2m} \left(\frac{\partial S}{\partial x}\right)^2 + \frac{\partial S}{\partial t} = 0$. As x is cyclic, the solution is $S = \alpha x - \frac{\alpha^2}{2m}t$. Here,

$$P = \alpha$$
, $Q = \beta = \frac{\partial S}{\partial \alpha} = x - \frac{\alpha}{m}t$, (5.14)

giving the exact transformation equations

$$x = \frac{\alpha}{m}t + \beta$$
, $p = \frac{\partial S}{\partial x} = \alpha$.

For simplicity, we can take the initial constants as $\alpha^{(0)} = \alpha_0$ and $\beta^{(0)} = \beta_0 = 0$. In terms of the new variables our perturbing Hamiltonian is $\Delta H = \frac{m\omega^2}{2} \left(\frac{\alpha}{m}t + \beta\right)^2$, so prior to expanding the full equations of motion are

$$\dot{\alpha} = -\frac{\partial \Delta H}{\partial \beta} = -m\omega^2 \left(\frac{\alpha}{m}t + \beta\right) ,$$

$$\dot{\beta} = \frac{\partial \Delta H}{\partial \alpha} = \omega^2 t \left(\frac{\alpha}{m}t + \beta\right) .$$
(5.15)

Plugging in 0th order solutions on the RHS, to 1st order we have

$$\dot{\alpha}^{(1)} = -\omega^2 \alpha_0 t \quad \Rightarrow \quad \alpha^{(1)}(t) = \alpha_0 - \frac{1}{2} \omega^2 \alpha_0 t^2 \,, \tag{5.16}$$

$$\dot{\beta}^{(1)} = \frac{\omega^2}{m} \alpha_0 t^2 \quad \Rightarrow \quad \beta^{(1)}(t) = \frac{\omega^2 \alpha_0 t^3}{3m} \,.$$

If we change back to our original variables with the inverse transformation (which we may wish to do at any point) this gives

$$p^{(1)} = \alpha^{(1)} = \alpha_0 - \frac{1}{2}\omega^2 \alpha_0 t^2, \tag{5.17}$$

and

$$x^{(1)}(t) = \frac{\alpha^{(1)}(t)}{m}t + \beta^{(1)}(t) = \frac{\alpha_0}{m}t - \frac{\omega^2\alpha_0}{m}\frac{t^3}{2} + \frac{\omega^2\alpha_0}{m}\frac{t^3}{3} = \frac{\alpha_0}{m}t - \frac{\omega^2\alpha_0}{m}\frac{t^3}{3!},$$
 (5.18)

which are the same results we previously obtained by solving Hamilton's equations perturbatively.

5.2 Periodic and Secular Perturbations to Finite Angle Pendulum

Example Let us consider a case where we do not have a simple solution. Consider a pendulum, with

$$H = \frac{p^2}{2ma^2} - mga\cos(\theta) \tag{5.19}$$

with $\theta \ll 1$. Expanding the cosine term we have

$$H = -mga + \frac{p^2}{2ma^2} + \frac{mga}{2}\theta^2 + \frac{mga}{2}\theta^2 \left(-\frac{\theta^2}{12} + \frac{\theta^4}{360} + \dots\right).$$
 (5.20)

In this case, the first term is a constant that will not play a role in our equations of motion, so we can identify $H_0 = \frac{p^2}{2ma^2} + \frac{mga}{2}\theta^2$. If we are only interested in applying first order perturbation theory we can simply take $\Delta H = -\frac{mga}{24}\theta^4$ and drop terms of $\mathcal{O}(\theta^6)$ and higher. The Hamiltonian H_0 is just a harmonic oscillator with 'moment of inertia $I = ma^2$ and frequency $\Omega^2 = \frac{g}{a}$. Again we use Ω here for angular frequency of the H_0 harmonic oscillator, to avoid confusion with the angle variable ω .

The action-angle variables for H_0 are

$$\alpha = H_0 = \frac{\Omega}{2\pi} J, \qquad \omega = \nu t + \beta$$
 (5.21)

where $\nu = \frac{\Omega}{2\pi}$, J is the action variable, and ω is the angle variable. This gives

$$\theta = \sqrt{\frac{2\alpha}{I\Omega^2}} \sin(\Omega t + \delta) = \sqrt{\frac{J}{\pi I\Omega}} \sin\left[2\pi(\nu t + \beta)\right],$$

$$p = \sqrt{2I\alpha} \cos(\Omega t + \delta) = \sqrt{\frac{IJ\Omega}{\pi}} \cos\left[2\pi(\nu t + \beta)\right].$$
(5.22)

Since ω and β are linearly related, the fact that (J, ω) are canonical implies that (J, β) are also a canonical pair. So we are free to take (J, β) as our new canonical variables when using the transformation in Eq. (5.22).

If we use (J, β) as the new variables, with $J^{(0)} = J_0$ and $\beta^{(0)} = \beta_0$ as given constants fixed by the initial conditions, then in terms of the new variables

$$\Delta H = -\frac{mga}{24}\theta^4 = -\frac{J^2}{24\pi^2 I}\sin^4(2\pi(\nu t + \beta)). \tag{5.23}$$

Expanding by using the 0th order solution gives

$$\dot{\beta}^{(1)} = \frac{\partial \Delta H}{\partial J} \Big|_{0} = -\frac{J_{0}}{12\pi^{2}I} \sin^{4}(2\pi(\nu t + \beta_{0})),
\dot{J}^{(1)} = -\frac{\partial \Delta H}{\partial \beta} \Big|_{0} = \frac{J_{0}^{2}}{3\pi I} \sin^{3}(2\pi(\nu t + \beta_{0})) \cos(2\pi(\nu t + \beta_{0})).$$
(5.24)

These results can be integrated to give $\beta^{(1)} = \beta^{(1)}(J_0, \beta_0, \nu, t)$ and $J^{(1)} = J^{(1)}(J_0, \beta_0, \nu, t)$. Before we consider computing these functions, lets pause to characterize two types of solution that occur in a more general context than simply this example.

Often we can characterize the nature of the perturbative solution without requiring a full study of the analytic form of a solution. A common situation where this is the case is when H_0 exhibits periodic orbits (as in the harmonic oscillator) with some frequency ν . In this case a relevant question is the following: what cumulative effect does the small perturbation have after going through one or more periods $T = \frac{1}{\nu}$? There are two possibilities:

- The perturbation itself could be *periodic*, where the parameter returns to its initial value. Here the perturbed trajectory looks much like the unperturbed one.
- Alternatively, we could have a net increment in the parameter after each orbit, called a *secular* change. After many periods, the parameter will be quite different from its value in H_0 .

Example Returning to our pendulum from before, the interesting quantity to study is the average over one period of the time rate of change of the variable,

$$\overline{\dot{J}^{(1)}} = \frac{1}{T} \int_0^T \dot{J}^{(1)}(t) dt = \frac{J^{(1)}(T) - J^{(1)}(0)}{T},$$
 (5.25)

since this tells us how much the variable changes over one period. For our example $\overline{\dot{J}^{(1)}} = 0$ because $\int_0^{2\pi} \sin^3(\theta) \cos(\theta) d\theta = 0$, and therefore the perturbation to J is periodic. Actually, from integrating Eq. (5.24) we have

$$J^{(1)}(t) = J_0 + \frac{J_0^2}{24\pi^2 I \nu} \sin^4(2\pi(\nu t + \beta_0)).$$
 (5.26)

Note from Eq. (5.22) that J determines the amplitude for $\theta(t)$ and p(t). A comparison between the trajectory with J_0 and with $J^{(1)}(t)$ is made in Fig. 5.1, where for this figure we set $\beta_0 = 0$.

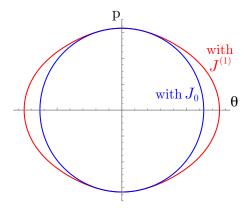


Figure 5.1: Comparison of the pendulum's periodic phase space trajectory using J_0 and $J^{(1)}(t)$.

In contrast, using $\int_0^{2\pi} \sin^4(\theta) \frac{d\theta}{2\pi} = \frac{3}{8}$, we find

$$\overline{\dot{\beta}^{(1)}} = \frac{1}{T} \int_0^T \dot{\beta}^{(1)}(t) dt = \frac{\beta^{(1)}(T) - \beta^{(1)}(0)}{T} = -\frac{J_0}{32\pi^2 I},$$
 (5.27)

which means β experiences a secular change. After many periods $(t \gg T)$ the change continues to build up, and we have on average that $\beta^{(1)}(t) \approx \overline{\dot{\beta}^{(1)}} t + \beta_0$. (If we look at the exact solution for $\beta^{(1)}(t)$ then it has precisely this linear term in t, plus terms that are periodic over the period τ , and that is what we mean by the \approx here.) Looking back at how the $\beta(t)$ dependence appears in $\theta = \theta(J, \beta, t)$ and $p = p(J, \beta, t)$ from Eq. (5.22), we see that on average the 1st order perturbation simply shifts the frequency to $\nu' = \nu + \dot{\beta}^{(1)}$.

Recall that we determined the full frequency $\nu_{\text{full}}(E)$ numerically as an example in our study of action-angle variables, which is shown below in Figure 5.2. Recalling that $J_0 =$

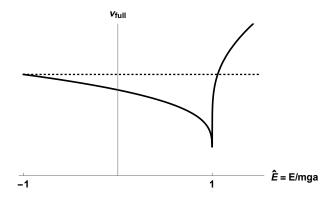


Figure 5.2: The full frequency ν_{full} vs. \hat{E}

 $\frac{H_0}{\nu} = \frac{E + mga}{\nu}$, we can write our perturbative shift to the frequency as a function of energy

$$\nu' - \nu = \overline{\dot{\beta}^{(1)}} = -\frac{(E + mga)}{32\pi^2 ma^2 \nu}.$$
 (5.28)

This is the first order correction to $\nu_{\text{full}}(E)$ when it is expanded about the simple harmonic oscillator minimum at E = -mga, which in Fig. 5.2 gives the negative linear correction to the frequency that occurs just above $\hat{E} = E/(mga) = -1$.

5.3 Perihelion Precession from Perturbing a Kepler Orbit

Kepler Example: Consider a central force perturbation

$$H = \underbrace{\frac{1}{2m} \left(p_r^2 + \frac{p_\psi^2}{r^2} \right) - \frac{k}{r}}_{H_0} \underbrace{-\frac{h}{r^n}}_{\Delta H}$$
 (5.29)

where the coupling parameter h is small and $n \geq 2$. The action-angle variables for H_0 are

$$J_1 = J_{\phi}$$

$$J_2 = J_{\theta} + J_{\phi}$$

$$J_3 = J_r + J_{\theta} + J_{\phi}$$

$$\omega_1 = \omega_{\phi} - \omega_{\theta}$$

$$\omega_2 = \omega_{\theta} - \omega_r$$

$$\omega_3 = \omega_r$$

where only $\dot{\omega}_3 = \nu_r \neq 0$, and all the others are constant in time. The pairs (ω_i, J_i) for $i \in \{1, 2, 3\}$ are all canonically conjugate. One way to see this is to note that we can implement a change of variables from the canonical pairs $\{(\omega_r, J_r), (\omega_\theta, J_\theta), (\omega_\phi, J_\phi)\}$ that we considered earlier, to these variables by using the generating function

$$F_2 = (\omega_\phi - \omega_\theta)J_1 + (\omega_\theta - \omega_r)J_2 + \omega_r J_3. \tag{5.30}$$

Let us study the perihelion precession, with the parameter $\omega = 2\pi\omega_2$ determining the perihelion angle. Some examples of precession are shown in Fig. 5.3, where in the case of a planet, the precession is like that of Fig. 5.3(b) with the sun at the focus of the ellipse. We

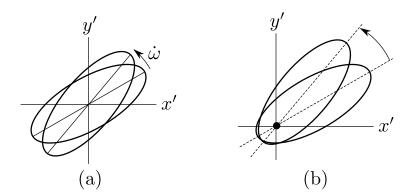


Figure 5.3: Precession of the perihelion from the point of view of coordinates centered on (a) the center of the ellipse, and (b) the focus of the ellipse. The latter is relevant for a planet like mercury orbiting the sun (shown with exaggerated eccentricity here).

know that

$$\dot{\omega}_2 = \frac{\partial \Delta H}{\partial J_2}, \qquad J_2 = J_\theta + J_\phi = 2\pi\alpha_\theta = 2\pi\ell \tag{5.31}$$

where $\ell = |\mathbf{L}|$ is the magnitude of the angular momentum. From the equation of motion $\dot{w}_2 = \partial \Delta H / \partial J_2$ we therefore have

$$\dot{\omega} = 2\pi \frac{\partial \Delta H}{\partial J_2} = \frac{\partial \Delta H}{\partial \ell},\tag{5.32}$$

and perturbatively, $\dot{\omega}^{(1)} = \frac{\partial \Delta H}{\partial \ell} \Big|_{0}$. We can average over the orbit to find the secular change:

$$\overline{\dot{\omega}^{(1)}} = \frac{1}{T} \int_0^T \frac{\partial \Delta H}{\partial \ell} \bigg|_0 dt. \tag{5.33}$$

Recall for the Kepler problem that $E^{(0)} = -(2\pi^2 k^2 m)/J_3^2$, and that at 0th order the period is

$$T = \frac{1}{\nu_r} = \pi k \sqrt{\frac{m}{-2E^{(0)3}}} = T(J_3). \tag{5.34}$$

Therefore $\frac{\partial}{\partial \ell} = 2\pi \frac{\partial}{\partial J_2}$ does not act on $T = T(J_3)$, so we can pull the partial derivative outside the integral,

$$\overline{\dot{\omega}^{(1)}} = \frac{\partial}{\partial \ell} \left(\frac{1}{T} \int_0^T \Delta H|_0 \, dt \right) = \frac{\partial}{\partial \ell} \, \overline{\Delta H|_0} \,. \tag{5.35}$$

Thus we must calculate the average of ΔH over one period.

$$\overline{\Delta H|_0} = -\frac{h}{T} \int_0^T \frac{dt}{r^n(t)}.$$
 (5.36)

Using $\ell = mr^2\dot{\psi}$ so that $dt = \frac{mr^2}{\ell}d\psi$ to switch variables from t to ψ , and then using the orbital equation for $r(\psi)$ we have

$$\overline{\Delta H|_{0}} = -\frac{hm}{\ell T} \int_{0}^{2\pi} \frac{d\psi}{r^{n-2}(\psi)} = -\frac{hm}{\ell T} \left(\frac{mk}{\ell^{2}}\right)^{n-2} \int_{0}^{2\pi} \left[1 + \epsilon \cos(\psi - \psi')\right]^{n-2} d\psi, \quad (5.37)$$

where the eccentricity $\epsilon = \sqrt{1 + \frac{2E\ell^2}{mk^2}}$ also depends on ℓ . There are two simple cases where can perform this integral:

- If n=2, then $\overline{\Delta H}=-\frac{2\pi hm}{\ell T}$, so $\overline{\dot{\omega}^{(1)}}=\frac{2\pi hm}{\ell^2 T}$.
- If n = 3, then $\overline{\Delta H} = -\frac{hkm^2}{\ell^3 T} \int_0^{2\pi} d\psi (1 + \epsilon \cos(\psi \psi'))$, where the cos term vanishes upon integration, so $\overline{\dot{\omega}^{(1)}} = \frac{\partial \overline{\Delta H}|_0}{\partial \ell} = \frac{6\pi m^2 hk}{\ell^4 T}$.

The latter type of potential (n = 3) is induced by corrections from general relativity to the Newtonian potential. The Schwarzschild metric is

$$ds^{2} = -c^{2}(d\tau)^{2} = \left(1 - \frac{r_{S}}{r}\right)c^{2}(dt)^{2} - \frac{(dr)^{2}}{1 - \frac{r_{S}}{r}} - r^{2}(d\theta)^{2} - r^{2}\sin^{2}(\theta)(d\phi)^{2}$$
 (5.38)

for $r_s = \frac{2GM}{c^2}$ where M is the central mass (say of the sun), G is Newton's gravitational constant, and c is the speed of light. The geodesic equation for radial motion is given by

$$E = \frac{m}{2} \left(\frac{dr}{d\tau}\right)^2 + V_{\text{eff}}(r) \tag{5.39}$$

where $V_{\text{eff}}(r) = -\frac{mc^2r_{\text{S}}}{2r} + \frac{\ell^2}{2mr^2} - \frac{r_{\text{S}}\ell^2}{2mr^3}$. (More background details on the Schwarzchild metric and the derivation of this geodesic equation are discussed below in the Side Note on page 103.) Defining k = GMm, then the effective potential can be rewritten as

$$V_{\text{eff}}(r) = -\frac{k}{r} + \frac{\ell^2}{2mr^2} - \frac{k\ell^2}{c^2m^2r^3},\tag{5.40}$$

from which we can identify $h = \frac{k\ell^2}{c^2m^2}$. Note that h must be treated as a constant independent of the canonical variable ℓ for the purpose of the above perturbative analysis (we simply substitute this value for h at the end).

Plugging in the h for the general relativistic correction we have

$$\overline{\dot{\omega}^{(1)}} = \frac{6\pi m^2 h k}{\ell^4 T} = \frac{6\pi}{T} \frac{k^2}{\ell^2 c^2} = \frac{6\pi}{T} \frac{k}{mc^2} \frac{1}{a(1-\varepsilon^2)}, \tag{5.41}$$

where in the last equality we used $\ell^2 = mka(1 - \varepsilon^2)$, which follows from Eq. (4.216). For Mercury, the period $T = 0.2409T_{\rm Earth}$, eccentricity $\varepsilon = 0.2056$, and semi-major axis $a = 5.79 \times 10^7 \, {\rm km}$, while $\frac{k}{mc^2} = \frac{GM_{\rm sun}}{c^2} = 1.4766 \, {\rm km}$, so we get a precession rate of

$$\overline{\dot{\omega}^{(1)}} = 2.0836 \times 10^{-6} \,\text{radians/year} = 42.98 \,\text{arcseconds/century}$$
 (5.42)

from general relativity. (An arcsecond is 1/3600'th of a degree.) After removing other contributions, such as a shift of 531.54 arcseconds/century from perturbations by other planets, the data on mercury's orbit shows a shift of 43.1 arcseconds/century (excellent agreement!). This was historically one of the first tests of general relativity, and still remains an important one.

We could also consider perturbations involving momentum variables rather than coordinates, and things work out in a very similar manner. To see this lets do one example.

Example Consider the relativistic correction to harmonic oscillator where the relativistic energy

$$E = \sqrt{c^4 m^2 + c^2 p^2} = mc^2 + \frac{p^2}{2m} - \frac{p^4}{8m^3 c^2} + \dots$$
 (5.43)

In this case to analyze the first order perturbative correction we take

$$H = \underbrace{\frac{1}{2m} \left(p^2 + m^2 \Omega^2 q^2 \right)}_{H_0} \underbrace{-\frac{p^4}{8c^2 m^3}}_{\Delta H}.$$
 (5.44)

From H_0 , the variables have a canonical transformation from the H-J analysis that gives

$$q = \sqrt{\frac{J}{\pi m \Omega}} \sin(2\pi(\nu t + \beta)), \qquad p = \sqrt{\frac{Jm\Omega}{\pi}} \cos(2\pi(\nu t + \beta)). \qquad (5.45)$$

This gives

$$\Delta H = -\frac{J^2 \Omega^2}{8\pi^2 c^2 m} \cos^4 \left[2\pi (\nu t + \beta) \right]. \tag{5.46}$$

Since $\dot{J}^{(1)}$ is odd over one period, it turns out that J is periodic once again

$$\dot{J}^{(1)} = -\frac{\partial \Delta H}{\partial \beta} \bigg|_{0} \quad \Rightarrow \quad \overline{\dot{J}^{(1)}} = 0.$$
(5.47)

Meanwhile, the change for $\dot{\beta}^{(1)}$ is secular,

$$\dot{\beta}^{(1)} = \frac{\partial \Delta H}{\partial J} \bigg|_{0} = -\frac{J_0 \Omega^2}{4\pi^2 mc^2} \cos^4 \left[2\pi (\nu t + \beta) \right] \quad \Rightarrow \quad \overline{\dot{\beta}^{(1)}} = -\frac{3J_0 \Omega^2}{32\pi^2 mc^2} \,. \tag{5.48}$$

Thus, $\dot{\beta}^{(1)}$ from the relativistic correction ΔH is again a negative shift to the frequency of the oscillator.

This ends our discussion of perturbation theory in classical mechanics.

Side Note: The Schwarzchild Geodesic from Action Minimization

The Schwarzchild metric is given by

$$ds^{2} = g_{\mu\nu}dx^{\mu}dx^{\nu} = \left(1 - \frac{r_{s}}{r}\right)c^{2}t^{2} - \frac{dr^{2}}{\left(1 - \frac{r_{s}}{r}\right)} - r^{2}d\theta^{2} - r^{2}\sin^{2}\theta d\phi^{2}, \qquad (5.49)$$

where $r_s = \frac{2GM}{c^2}$ is the Schwartzchild radius. The geodesic orbit for a test particle is a curve which minimizes proper distance with this metric. In this case, we have

$$0 = \delta s = \delta \int ds = \delta \int \left(g_{\mu\nu} \frac{dx^{\mu}}{d\tau} \frac{dx^{\nu}}{d\tau} \right)^{\frac{1}{2}} d\tau \tag{5.50}$$

where τ is the proper time and $ds^2=c^2d\tau^2$. (One method of determining the geodesic equations is to use $0=\frac{d^2x^\lambda}{d\tau^2}+\Gamma^\lambda_{\mu\nu}\frac{dx^\mu}{d\tau}\frac{dx^\nu}{d\tau}$ with the Christoffel symbols $\Gamma^\lambda_{\mu\nu}$ determined from the metric, but we will follow a different approach.)

The minimization in Eq. (5.50) is equivalent to applying the minimal action principal for the Lagrangian

$$L = \frac{m}{2} g_{\mu\nu} \frac{dx^{\mu}}{d\tau} \frac{dx^{\nu}}{d\tau}$$

$$= \frac{m}{2} \left[\left(1 - \frac{r_s}{r} \right) c^2 \left(\frac{dt}{d\tau} \right)^2 - \frac{1}{\left(1 - \frac{r_s}{r} \right)} \left(\frac{dr}{d\tau} \right)^2 - r^2 \left(\frac{d\theta}{d\tau} \right)^2 - r^2 \sin^2 \theta \left(\frac{d\phi}{d\tau} \right)^2 \right]$$
(5.51)

since the presence of the extra square root $(...)^{\frac{1}{2}}$ does not matter for this minimization. Here we have the generalized coordinates $x^{\mu} = (t, r, \theta, \phi)$ which are to be considered as functions of the proper time variable τ . Also, the mass m is a test mass (which also gives us the proper units).

Because t and ϕ are cyclic variables in L, we have

$$p_{t} = \frac{\partial L}{\partial \dot{t}} = m \left(1 - \frac{r_{s}}{r} \right) c^{2} \frac{dt}{d\tau} = E^{\text{tot}}$$
 energy (5.52)
$$p_{\phi} = -\frac{\partial L}{\partial \dot{\phi}} = mr^{2} \sin^{2} \theta \frac{d\phi}{d\tau} = \ell$$
 angular momentum

Using the E-L equation on θ we obtain

$$0 = \frac{d}{d\tau} \frac{\partial L}{\partial \dot{\theta}} - \frac{\partial L}{\partial \theta}$$

$$= \frac{d}{d\tau} \left(mr^2 \frac{d\theta}{d\tau} \right) - mr^2 \sin \theta \cos \theta \dot{\phi}^2$$
(5.53)

where planar motion with $\theta = \frac{\pi}{2}$ is a solution that suffices for our purposes. Now

$$\frac{ds^2}{d\tau^2} = c^2 = \left(1 - \frac{r_s}{r}\right)c^2 \frac{E_{\text{tot}}^2}{m^2\left(1 - \frac{r_s}{r}\right)^2c^4} - \frac{1}{\left(1 - \frac{r_s}{r}\right)}\left(\frac{dr}{d\tau}\right)^2 - r^2 \frac{\ell^2}{m^2r^4}$$
(5.54)

gives a radial equation. Separating out the rest mass and expanding $E=E^{\rm tot}-mc^2\ll mc^2$ we have:

$$E^{\text{tot}} = mc^2 + E \quad \Rightarrow \quad (E^{\text{tot}})^2 \approx m^2 c^4 + 2mc^2 E \tag{5.55}$$

Putting this together we have

$$\underbrace{\frac{m}{2} \left(c^2 + \frac{l^2}{m^2 r^2}\right) \left(1 - \frac{r_s}{r}\right) - \frac{m}{2}c^2}_{V_{\text{eff}}(r)} + \frac{m}{2} \left(\frac{dr}{d\tau}\right)^2 = E$$
(5.56)

Note that the rest mass $mc^2/2$ terms cancel. Therefore the effective potential is

$$V_{\text{eff}}(r) = -\frac{r_s mc^2}{2r} + \frac{l^2}{2mr^2} - \frac{r_s l^2}{2mr^3}$$
$$= -\frac{k}{r} + \frac{l^2}{2mr^2} - \frac{kl^2}{m^2 c^2 r^3}$$

where $r_s = \frac{2GM}{c^2}$ and k = GMm, so $mr_sc^2 = 2k$. This is the result that was quoted above in Eq. (5.40).

Chapter 6

Fluid Mechanics

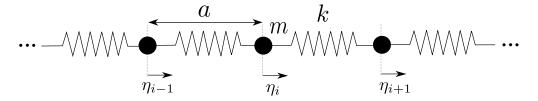
So far, our examples of mechanical systems have all been discrete, with some number of masses acted upon by forces. Even our study of rigid bodies (with a continuous mass distribution) treated those bodies as single objects. In this chapter, we will treat an important continuous system, which is that of fluids.

Fluids include both liquids and gases. (It also includes plasmas, but for them a proper treatment requires the inclusion of electromagnetic effects, which will not be discussed here.) For our purposes, a fluid is a material which can be treated as *continuous*, which has the ability to flow, and which has very little resistance to deformation (that is, it has only a small support for shear stress, which refers to forces parallel to an applied area). Applications include meteorology, oceanography, astrophysics, biophysics, condensed matter physics, geophysics, medicine, aerodynamics, plumbing, cosmology, heavy-ion collisions, and so on.

The treatment of fluids is an example of classical field theory, with continuous field variables as the generalized coordinates, as opposed to the discrete set of variables q_i that we have treated so far. Therefore the first step we have to take is understanding how to make the transition from discrete to continuum.

6.1 Transitioning from Discrete Particles to the Continuum

Rather than starting with fluids, lets instead consider a somewhat simpler system, that of an infinite one dimensional chain of masses m connected by springs with spring constant k, which we take as an approximation for an infinite continuous elastic one-dimensional rod.



If the equilibrium separation of masses is a and the distance the i'th mass is translated from equilibrium is η_i , then

$$V = \frac{k}{2} \sum_{i} (\eta_{i+1} - \eta_i)^2 \qquad T = \frac{m}{2} \sum_{i} \dot{\eta}_i^2, \tag{6.1}$$

where V is the potential energy from the springs, and T is the kinetic energy. It is convenient to write the Lagrangian as

$$L = T - V = \frac{1}{2} \sum_{i} a \left(\frac{m}{a} \dot{\eta}_i^2 - ka \left(\frac{\eta_{i+1} - \eta_i}{a} \right)^2 \right), \tag{6.2}$$

and the corresponding equations of motion obtained from the Euler-Lagrange equations as

$$\frac{m}{a}\ddot{\eta}_i - ka\left(\frac{\eta_{i+1} - \eta_i}{a^2}\right) + ka\left(\frac{\eta_i - \eta_{i-1}}{a^2}\right) = 0.$$
(6.3)

Technically both the Lagrangian and the equations of motion are independent of a, but we have introduced factors of a to facilitate taking the continuous limit $a \to 0$. In this limit the masses become progressively more densely distributed along the line. The important question when taking this limit is which quantities do we hold fixed.

Lets define $\mu \equiv \frac{m}{a}$ as the mass density and Y = ka as the Young's modulus. Here Y is equivalent to the spring constant for a continuous rod. (For a rod, the force $F = Y\xi$ where ξ is the longitudinal extension per unit length, or in other words, the strain.) We intend to hold μ and Y fixed when taking the continuous limit.

The key change in the continuous limit is that the discrete position index i becomes a continuous position label x, so instead of $\eta_i = \eta_i(t)$, now $\eta_x = \eta_x(t)$, or with more conventional notation, $\eta = \eta(x, t)$. This also means that

$$\frac{\eta(x+a,t) - \eta(x,t)}{a} \to \frac{\partial \eta}{\partial x},\tag{6.4}$$

$$\frac{1}{a} \left(\frac{\eta(x+a,t) - \eta(x,t)}{a} - \frac{\eta(x,t) - \eta(x-a,t)}{a} \right) \to \frac{\partial^2 \eta}{\partial x^2},\tag{6.5}$$

$$\sum_{i} a \to \int dx \,. \tag{6.6}$$

Using these results in Eq. (6.2) gives $L = \int \mathcal{L} dx$ where

$$\mathcal{L} = \frac{1}{2} \left(\mu \left(\frac{\partial \eta(x,t)}{\partial t} \right)^2 - Y \left(\frac{\partial \eta(x,t)}{\partial x} \right)^2 \right)$$
 (6.7)

is the Lagrangian density. Likewise, using them in Eq. (6.3) gives the equations of motion

$$\mu \frac{\partial^2 \eta}{\partial t^2} - Y \frac{\partial^2 \eta}{\partial r^2} = 0, \qquad (6.8)$$

which we recognize as the wave equation.

The results for the Lagrange density and Euler Lagrange equations are of course not indepedent. We can also use

$$\mathcal{L} = \mathcal{L}\left(\eta, \frac{\partial \eta}{\partial x}, \frac{\partial \eta}{\partial t}, x, t\right) \tag{6.9}$$

with Hamilton's principle,

$$\delta S = \delta \int_{t_1}^{t_2} \int_{-\infty}^{\infty} \mathcal{L} \, dx \, dt = 0. \tag{6.10}$$

to formulate the dynamics, and thus derive the Euler-Lagrange equations. Because $\eta = \eta(x,t)$ has two parameters, if we follow the standard procedure of varying the path η takes between the two endpoints, we get variations from the dependence of the Lagrange density on its first three arguments. Integrating by parts in each of t and x, and setting the surface terms to zero, then yields

$$\frac{\partial}{\partial t} \left(\frac{\partial \mathcal{L}}{\partial \left(\frac{\partial \eta}{\partial t} \right)} \right) + \frac{\partial}{\partial x} \left(\frac{\partial \mathcal{L}}{\partial \left(\frac{\partial \eta}{\partial x} \right)} \right) - \frac{\partial \mathcal{L}}{\partial \eta} = 0 \tag{6.11}$$

as the continuum Euler-Lagrange equation. Recall that for N particles we expect N E-L equations for the time dependence, but here we have just one equation. However actually by that counting, this result corresponds to an infinite number of equations, one for each value of x. From this point of view, the derivatives with respect to x are what couples these equations together.

Example For the Lagrangian density of the elastic rod,

$$\frac{\partial \mathcal{L}}{\partial \left(\frac{\partial \eta}{\partial t}\right)} = \mu \frac{\partial \eta}{\partial t}, \qquad \frac{\partial \mathcal{L}}{\partial \left(\frac{\partial \eta}{\partial x}\right)} = -Y \frac{\partial \eta}{\partial x}, \qquad \frac{\partial \mathcal{L}}{\partial \eta} = 0. \tag{6.12}$$

Putting these results together in Eq. (6.11) gives the wave equation $\mu \frac{\partial^2 \eta}{\partial t^2} - Y \frac{\partial^2 \eta}{\partial x^2} = 0$ as anticipated.

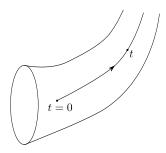
In our above analysis, $\eta = \eta(x, t)$ is a continuum generalized coordinate called a *classical field*. Here t is a parameter and x is a continuous label as well.

Although we have been talking about one dimension so far, it is easy to generalize the above discussion to a higher number of dimensions. For example, in three dimensions we simply have dependence on three continuous label parameters, $\eta = \eta(x, y, z, t)$ or $\eta = \eta(\mathbf{r}, t)$. The field $\eta(\mathbf{r}, t)$ is called a scalar field because the output is a single number. With multiple dimensions we also have vector fields $\eta(\mathbf{r}, t)$, where the output is a vector. An example of vector fields that you are familiar with are the electromagnetic fields $\mathbf{E}(\mathbf{r}, t)$ and $\mathbf{B}(\mathbf{r}, t)$. In fact, classical fields of exactly this sort are also the starting point for formulating quantum field theory. One formulates a classical Lagrangian density \mathcal{L} (that is most often Lorentz invariant) which depends on fields like the electromagnetic scalar and vector potentials $\phi(\mathbf{r}, t)$ and $\mathbf{A}(\mathbf{r}, t)$. Then one quantizes these fields.

Our description of fluids will make use of classical field variables in 3-dimensions without considering quantization.

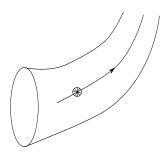
Eulerian & Lagrangian Variables

Let us consider fluid flowing in a tube. A natural set of variables would be to take some initial time t = 0 and label all fluid elements by their coordinates at this time: \mathbf{r}_0 . We will often refer to fluid element, which is a point in the fluid, or sometimes an infinitesimally small region of the fluid. The motion of the fluid could then be described by $\mathbf{r} = \mathbf{r}(\mathbf{r}_0, t)$, which determines the subsequent coordinates of the fluid element labeled by \mathbf{r}_0 at the later time t.



If we continue in this direction it leads to the Lagrangian formulation of fluid dynamics. The advantage is that the usual laws of classical mechanics apply to fluid particles. The disadvantage is that it is more work to make the connection to measurable quantities that characterize the fluid.

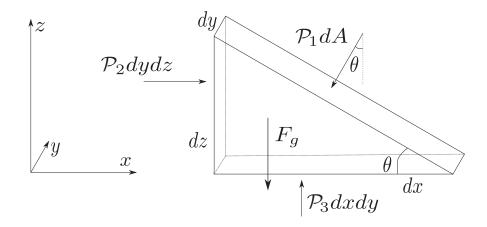
Instead, we will explore an alternate formulation. Pretend we sit at a fixed point \mathbf{r} in the fluid and ask what happens there as a function of the time t. We can think of it like placing a measuring device at \mathbf{r} and measuring, at any time t, quantities like the fluid velocity \mathbf{v} , density ρ , or pressure \mathcal{P} . In fact, these 5 variables (as density and pressure are scalars, while velocity is a vector) are enough to describe the state of the moving fluid. This is the Eulerian formulation of fluid dynamics. Additionally, ρ and \mathcal{P} are enough to determine all thermodynamic quantities (assuming we know the equation of state).



Example for an ideal gas at temperature T, we have $\mathcal{P}V = nRT$, where V is volume, n is the number of moles, R is contant, and T is temperature. Dividing by V we have that pressure is simply proportional to density, $\mathcal{P} = \rho R'T$ for a rescaled gas constant R'.

Pressure is *isotropic* in a fluid. This means it is the same from all directions, so there is only 1 number for each (\mathbf{r}, t) ; thus, pressure is a scalar field, $\mathcal{P}(\mathbf{r}, t)$.

To prove this consider the infinitesimal wedge-shaped fluid element below, which we take to be at rest in a gravitational field. Recall that pressure is force per unit area, $\mathcal{P} = \hat{n} \cdot \mathbf{F}/A$, where \hat{n} is a unit vector perpendicular to the area A that the force \mathbf{F} is acting on. In the figure we label three pressures \mathcal{P}_1 , \mathcal{P}_1 , \mathcal{P}_1 acting on three of the sides, that have corresponding force components in the x and z directions. (There is also pressure on the other two faces with forces in the y direction, but the above will suffice to argue why the pressure is isotropic.)



Next we balance forces for the wedge at rest. Gravity pulls down, and the volume of the wedge is dxdydz/2 so the force of gravity is

$$F_g = \rho g \, \frac{dx \, dy \, dz}{2} \,. \tag{6.13}$$

Also by simple trigonometry the area dA of the slanted face can be written in two different ways

$$dA = dy \frac{dz}{\sin(\theta)} = dy \frac{dx}{\cos(\theta)}.$$

Balancing forces in the x and z directions then means

$$0 = dF_x = \mathcal{P}_2 \, dy \, dz - \mathcal{P}_1 \sin(\theta) \, dA = (\mathcal{P}_2 - \mathcal{P}_1) \, dy \, dz,$$

$$0 = dF_z = \mathcal{P}_3 \, dx \, dy - \mathcal{P}_1 \cos(\theta) \, dA - \frac{1}{2} \rho g \, dx \, dy \, dz.$$

$$(6.14)$$

The first equation implies $\mathcal{P}_1 = \mathcal{P}_2$. In the second equation we can pull out a common dxdy to give

$$\mathcal{P}_3 = \mathcal{P}_1 + \frac{1}{2}\rho g \, dz,\tag{6.15}$$

then as the infinitesimal distance $dz \to 0$ we have

$$\mathcal{P}_1 = \mathcal{P}_3. \tag{6.16}$$

Thus, pressure is the same in all directions. Even if the fluid is moving or even accelerating we would come to the same conclusion. For example, if we had to balance the force against acceleration this would lead to adding term

$$\rho \mathbf{a} \, dx \, dy \, dz \tag{6.17}$$

for acceleration \mathbf{a} , which again drops out for an infinitesimal fluid element just like the gravitational force did.

Time Derivatives:

The total time derivative $\frac{d}{dt}$ tells us the rate at which a quantity changes as we move with a fluid element. The partial time derivative $\frac{\partial}{\partial t}$ tells us the rate of change of a quantity at a fixed position **r**. We can work out a relation between them.

Example For \mathcal{P}

$$\frac{d\mathcal{P}}{dt} = \frac{\partial \mathcal{P}}{\partial t} + \frac{\partial \mathcal{P}}{\partial x}\dot{x} + \frac{\partial \mathcal{P}}{\partial y}\dot{y} + \frac{\partial \mathcal{P}}{\partial z}\dot{z}
= \frac{\partial \mathcal{P}}{\partial t} + \mathbf{v} \cdot \nabla \mathcal{P}.$$
(6.18)

In general, the time derivative acts as

$$\frac{d}{dt} = \frac{\partial}{\partial t} + \mathbf{v} \cdot \nabla \tag{6.19}$$

on any fluid field (scalar, vector, or tensor) that is a function of (x, y, z, t).

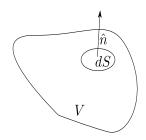
6.2 Fluid Equations of Motion

6.2.1 Continuity Equations

Let us consider a volume V of fluid. Then from Gauss' divergence theorem (which you may recall from electomagnetism, but which is a general result of vector calculus for any vector field):

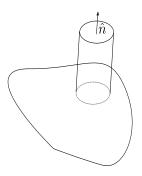
$$\int_{V} dV \nabla \cdot \mathbf{v} = \int_{\partial V} d\mathcal{S} \,\hat{n} \cdot \mathbf{v} = \int_{\partial V} d\mathbf{S} \cdot \mathbf{v}, \quad (6.20)$$

where ∂V is the closed area that bounds the volume V, \hat{n} is unit vector orthogonal to the surface element dS, and $d\mathbf{S} \equiv dS\hat{n}$.



Lets ask: As the fluid moves, how does V change?

The quantity $\hat{n} \cdot \mathbf{v}$ is the outward velocity of the surface $d\mathcal{S}$, so $\hat{n} \cdot \mathbf{v} d\mathcal{S} dt$ is the volume added in a time interval dt, as illustrated on the right.



This means that the change in volume can be determined by adding up all the changes from integrating over the entire surface

$$\frac{dV}{dt} = \int_{\partial V} \mathbf{v} \cdot d\mathbf{S} = \int_{V} \nabla \cdot \mathbf{v} \, dV. \tag{6.21}$$

This result becomes even simpler if we consider an infinitesimal volume δV over which $\nabla \cdot \mathbf{v}$ does not vary, then we can pull $\nabla \cdot \mathbf{v}$ outside the integral to obtain simply

$$\frac{d\delta V}{dt} = \delta V \nabla \cdot \mathbf{v}. \tag{6.22}$$

Thus the divergence of the velocity, $\nabla \cdot \mathbf{v}$, controls how the fluid volume expands with time. If $\nabla \cdot \mathbf{v} = 0$ everywhere then we say the fluid is *incompressible* because for every volume element $\frac{dV}{dt} = 0$.

Even if the volume changes, the mass of the fluid element will not,

$$\frac{d\delta m}{dt} = \frac{d}{dt} \left(\rho \delta V \right) = 0. \tag{6.23}$$

This implies that

$$0 = \delta V \frac{d\rho}{dt} + \rho \frac{d\delta V}{dt} = \delta V \left(\frac{d\rho}{dt} + \rho \nabla \cdot \mathbf{v} \right), \tag{6.24}$$

SO

$$\frac{d\rho}{dt} + \rho \nabla \cdot \mathbf{v} = 0. \tag{6.25}$$

Expanding out the time derivative into partial derivatives this yields

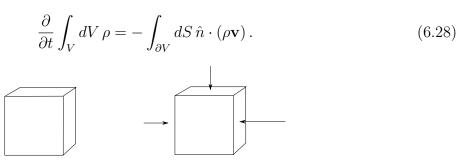
$$\frac{\partial \rho}{\partial t} + \mathbf{v} \cdot \nabla \rho + \rho \nabla \cdot \mathbf{v} = 0, \tag{6.26}$$

and simplifying the result gives

$$\frac{\partial \rho}{\partial t} + \nabla \cdot (\rho \mathbf{v}) = 0. \tag{6.27}$$

This is an important result the *continuity equation* for mass of the fluid, which is a partial differential equation in the fluid variables.

Here $\rho \mathbf{v}$ is the mass density flux (the flow of mass density), and this equation says that nowhere is the matter making up the fluid created or destroyed. If the density changes, then there must be an incoming or outgoing flux. This is easier to see if we use the divergence theorem for the vector $(\rho \mathbf{v})$ to write this result in integral form,



Mass increase equals the inflowing mass

Here $\frac{\partial}{\partial t} \int_V dV \rho$ is the increase of mass in the volume V, while $\int_{\partial V} dS \hat{n} \cdot (\rho \mathbf{v})$ is the outflow of mass through the surface surrounding this volume (which becomes an inflow with the minus sign).

We've talked about $\nabla \cdot \mathbf{v}$ thus far, so it's natural to ask: is there a physical interpretation to $\nabla \times \mathbf{v}$? There is indeed.

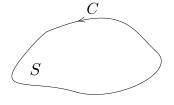
The quantity $\Omega = \nabla \times \mathbf{v}$ is the *vorticity*. If the velocity is constant or uniform in a single direction $\vec{v} = v_x(x)\hat{x}$, then its obvious that $\Omega = 0$. To consider what it measures we can use Stoke's theorem for the velocity vector field,

$$\int_{\mathcal{S}} (\nabla \times \mathbf{v}) \cdot \hat{n} \, d\mathcal{S} = \oint_{\partial \mathcal{S}} \mathbf{v} \cdot d\mathbf{l}, \qquad (6.29)$$

where \mathcal{S} is now an open surface and $\partial \mathcal{S}$ is its closed boundary curve. We can use this result to determine when $\Omega = \nabla \times \mathbf{v}$ is non-zero.

Consider a rotating fluid and the circular curve shown. When we integrate the velocity along this circular curve the terms always add up, so $\Omega \neq 0$. It thus provides a measures of the rate of rotation of the fluid.

If we consider flow in one direction with a velocity gradient, again $\Omega \neq 0$. Here a larger positive contribution to the integral is obtained from the right side of the circle, relative to the negative contribution from the left side.







If $\nabla \times \mathbf{v} = 0$ everywhere in a moving fluid, we say the flow is *irrotational*.

We can determine more precisely what Ω is as follows. Consider a rotating coordinate

system with constant angular velocity $\boldsymbol{\omega}$, so $\mathbf{v} = \mathbf{v}' + \boldsymbol{\omega} \times \mathbf{r}$. Then

$$\nabla \times \mathbf{v} = \nabla \times \mathbf{v}' + \nabla \times (\boldsymbol{\omega} \times \mathbf{r})$$
$$= \nabla \times \mathbf{v}' + \boldsymbol{\omega} \nabla \cdot \mathbf{r} - \boldsymbol{\omega} \cdot \nabla \mathbf{r}$$
$$= \nabla \times \mathbf{v}' + 2\boldsymbol{\omega}.$$

Now ω was constant, but we have not yet specified its value. If we pick $\omega = \frac{1}{2}\nabla \times \mathbf{v}$ at a point \mathbf{r} , then $\nabla \times \mathbf{v}' = 0$ and the fluid is irrotational at \mathbf{r} in the rotating frame. This argument can be repeated for other points \mathbf{r} . Considering this from the point of view of the original frame we thus see that $\Omega/2$ is the angular velocity of the fluid at position \mathbf{r} .

6.2.2 Ideal Fluid: Euler's Equation and Entropy Conservation

Let us consider an *ideal fluid* which has no energy dissipation due to internal friction (meaning no viscosity) and no heat exchange between different parts of the fluid (meaning no thermal conductivity).

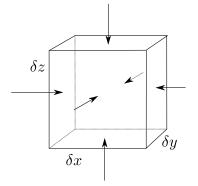
The force on a fluid element δV from pressure in the \hat{x} direction is

$$\delta F_x = F_x(x) - F_x(x + \delta x) =$$

$$= \frac{\mathcal{P}(x) - \mathcal{P}(x + \delta x)}{\delta x} \, \delta y \, \delta z \, \delta x = -\frac{\partial \mathcal{P}}{\partial x} \delta V \,. \tag{6.30}$$

More generally accounting for all directions we have

$$\delta \mathbf{F} = -\nabla \mathcal{P} \, \delta V. \tag{6.31}$$



From external forces $\delta \mathbf{F}$ it is useful to define the force density \mathbf{f} by $\delta \mathbf{F} = \mathbf{f} \delta V$. This means that Newton's law, $m\mathbf{a} = \mathbf{F}$ becomes

$$\rho \delta V \frac{d\mathbf{v}}{dt} = (-\nabla \mathcal{P} + \mathbf{f})\delta V. \tag{6.32}$$

Writing out the total time derivatives this becomes

$$\frac{\partial \mathbf{v}}{\partial t} + \mathbf{v} \cdot \nabla \mathbf{v} + \frac{\nabla \mathcal{P}}{\rho} = \frac{\mathbf{f}}{\rho}$$
 (6.33)

which is the Euler equation for fluid dynamics. There are two special cases which are useful to consider.

1. Often **f** can be derived from a potential: $\mathbf{f} = -\rho \nabla \Phi$. Note that Φ here is a potential energy per unit mass. For example, with gravity, $\Phi = gz$, so $\mathbf{f} = -\rho g\hat{z}$. In general, then

$$\frac{\partial \mathbf{v}}{\partial t} + \mathbf{v} \cdot \nabla \mathbf{v} + \frac{\nabla \mathcal{P}}{\rho} + \nabla \Phi = 0 \tag{6.34}$$

is a rewriting of the Euler equation.

2. We can use $\mathbf{v} \times \mathbf{\Omega} = \mathbf{v} \times (\nabla \times \mathbf{v}) = \nabla \left(\frac{\mathbf{v}^2}{2}\right) - \mathbf{v} \cdot \nabla \mathbf{v}$. This gives

$$\frac{\partial \mathbf{v}}{\partial t} + \nabla \left(\frac{\mathbf{v}^2}{2} + \Phi\right) + \frac{\nabla \mathcal{P}}{\rho} - \mathbf{v} \times \mathbf{\Omega} = 0 \tag{6.35}$$

as another rewriting of the Euler equation. For constant ρ , we can take the curl and use the fact that $\nabla \times (\nabla h) = 0$ for any h to obtain

$$\frac{\partial \mathbf{\Omega}}{\partial t} = \nabla \times (\mathbf{v} \times \mathbf{\Omega}). \tag{6.36}$$

We will have occasion to use this result later on.

To solve the Euler and continuity partial differential equations we need boundary conditions. The boundary conditions for ideal fluids are simply that they cannot penetrate solid surfaces, so if a surface with normal vector \hat{n} is stationary then

$$\mathbf{v} \cdot \hat{n} \Big|_{\text{surface}} = 0. \tag{6.37}$$

If the surface does move with velocity $\mathbf{v}_{\text{surface}}$ then

$$\mathbf{v} \cdot \hat{n} \Big|_{\text{surface}} = \mathbf{v}_{\text{surface}} \cdot \hat{n}. \tag{6.38}$$

So far we have four equations (continuity for the scalar density ρ , and the Euler equation for the vector velocity \mathbf{v}) for five unknowns. For an ideal fluid the 5th equation,

$$\frac{dS}{dt} = 0, (6.39)$$

is the statement that the entropy S is conserved, so there is no heat exchange. Effectively, this provides a relationship between pressure and density through $\mathcal{P} = \mathcal{P}(\rho, S)$. A simple example is an ideal gas at constant temperature, where $\mathcal{P} = \rho R'T$.

6.2.3 Conservation of Momentum and Energy

Due to the term $(\mathbf{v} \cdot \nabla)\mathbf{v}$ which has a $v_i v_j$, the Euler equation is nonlinear. For simpler situations it is therefore very useful to consider conservation laws.

Let us start by considering *Momentum Conservation*. The quantity $\rho \mathbf{v}$ is the flux of mass density, which is also the density of momentum (in direct analogy to $\mathbf{p} = m\mathbf{v}$). Consider $\frac{\partial}{\partial t}(\rho v_i) = \frac{\partial \rho}{\partial t}v_i + \rho \frac{\partial v_i}{\partial t}$. Using the continuity and Euler equations to replace these two partial derivatives, and once again implicitly summing over repeated indices until further notice, this becomes

$$\frac{\partial}{\partial t} (\rho v_i) = -v_i \frac{\partial}{\partial x_j} (\rho v_j) + \rho \left(-v_j \frac{\partial v_i}{\partial x_j} - \frac{1}{\rho} \frac{\partial \mathcal{P}}{\partial x_i} \right) + f_i. \tag{6.40}$$

This is rearranged to give

$$\frac{\partial}{\partial t} (\rho v_i) + \frac{\partial \mathcal{P}}{\partial x_i} + \frac{\partial}{\partial x_j} (\rho v_i v_j) = f_i. \tag{6.41}$$

We define the *stress tensor* for an ideal fluid as

$$T_{ij} = \mathcal{P}\delta_{ij} + \rho v_i v_j \tag{6.42}$$

which gives the momentum flux density (which is to say, the density of momentum in the direction \mathbf{e}_i flowing in the direction of \mathbf{e}_j). Note that T_{ij} is symmetric. Then, in vector form, the equation above becomes

$$\frac{\partial}{\partial t} (\rho \mathbf{v}) + \nabla \cdot \mathbb{T} = \mathbf{f}. \tag{6.43}$$

This describes the conservation of linear momentum density with sources given by the external force densities \mathbf{f} . Comparing to the continuity equation where the density ρ is a scalar so its flux $\rho \mathbf{v}$ is a vector, here the momentum density $\rho \mathbf{v}$ is a vector so its flux \mathbb{T} is a tensor.

Next consider Conservation of Energy. Conservation of energy can be said to arise from the Euler equation $d\mathbf{v}/dt + \nabla \mathcal{P}/\rho + \nabla \Phi = 0$. Note that here we are switching back to the total time derivative since this is more useful for our discussion of energy. For a volume element δV , we take the inner product of the Euler equation with $\rho \mathbf{v} \delta V$ to obtain

$$\delta V \rho \mathbf{v} \cdot \frac{d\mathbf{v}}{dt} + \delta V \mathbf{v} \cdot \nabla \mathcal{P} + \delta V \rho \mathbf{v} \cdot \nabla \Phi = 0.$$
 (6.44)

Given that $\frac{d}{dt}(\rho\delta V) = 0$ by the conservation of mass we can move this combination inside of total time derivatives. Also recall that $\mathbf{v} \cdot \nabla = d/dt - \partial/\partial t$. Using these two facts we obtain

$$\frac{d}{dt}\left(\frac{1}{2}\rho\mathbf{v}^2\delta V\right) + (\mathbf{v}\cdot\nabla\mathcal{P})\delta V + \frac{d}{dt}\left(\rho\Phi\delta V\right) - \rho\frac{\partial\Phi}{\partial t}\delta V = 0.$$
 (6.45)

Next we try to convert the second term to time derivatives. Consider using

$$\frac{d}{dt}(\mathcal{P}\delta V) = \frac{\partial \mathcal{P}}{\partial t}\delta V + (\mathbf{v}\cdot\nabla\mathcal{P})\delta V + \mathcal{P}(\nabla\cdot\mathbf{v})\delta V, \tag{6.46}$$

where we recalled that $d\delta V/dt = (\nabla \cdot \mathbf{v})\delta V$. Using this to eliminate $(\mathbf{v} \cdot \nabla \mathcal{P})\delta V$ gives

$$\frac{d}{dt} \left(\frac{1}{2} \rho \mathbf{v}^2 \delta V + \rho \Phi \delta V + \mathcal{P} \delta V \right) = \left(\frac{\partial \mathcal{P}}{\partial t} + \rho \frac{\partial \Phi}{\partial t} \right) \delta V + \mathcal{P} \nabla \cdot \mathbf{v} \delta V, \tag{6.47}$$

where $\frac{1}{2}\rho \mathbf{v}^2 \delta V$ is the kinetic energy, $\rho \Phi \delta V$ is the external potential energy, and $\mathcal{P} \delta V$ is the internal potential energy due to pressure. The terms with partial time derivatives act like sources. Unfortunately there is still a term without a total or partial time derivative, however

this term is easier to interpret. It is related to the work W_u done by δV when it expands and exerts pressure on the surrounding fluid. In particular

$$\frac{dW_u}{dt} = \mathcal{P}\frac{d(\delta V)}{dt} = \mathcal{P}\nabla \cdot \mathbf{v}\,\delta V \equiv -\frac{d}{dt}(U\delta m) \tag{6.48}$$

where in the last step we have defined the work as a negative potential energy U (per unit mass). If the equation of state is given, this U can be calculated either as an integral in density or pressure,

$$U = \int d(\delta V) \frac{\mathcal{P}}{\delta m} = \int_{\rho_0}^{\rho} d\rho' \frac{\mathcal{P}(\rho')}{{\rho'}^2} = \int_{\mathcal{P}_0}^{\mathcal{P}} \mathcal{P}' \left(\frac{1}{\rho^2} \frac{d\rho}{d\mathcal{P}'}\right) (\mathcal{P}') d\mathcal{P}'$$
 (6.49)

where we used the fixed $\delta m = \delta V \rho$ to switch variables between δV and ρ . Using Eq. (6.48) means we can now write everything in terms of time derivatives,

$$\frac{d}{dt}\left(\left(\frac{1}{2}\rho\mathbf{v}^2 + \rho\Phi + \mathcal{P} + \rho U\right)\delta V\right) = \left(\frac{\partial\mathcal{P}}{\partial t} + \rho\frac{\partial\Phi}{\partial t}\right)\delta V \tag{6.50}$$

This is the equation for energy conservation in an ideal fluid. If \mathcal{P} and Φ are not explicitly dependent on time at any point in space (which is often the case), then any fluid element has a constant total energy as it moves along (recall that this is the meaning of d/dt). For applications to fluids it is more convenient to divide this result by $\delta m = \rho \delta V$ to give

$$\frac{d}{dt}\left(\frac{\mathbf{v}^2}{2} + \Phi + \frac{\mathcal{P}}{\rho} + U\right) = \frac{1}{\rho}\frac{\partial \mathcal{P}}{\partial t} + \frac{\partial \Phi}{\partial t}.$$
 (6.51)

This is Bernoulli's equation.

6.3 Static Fluids & Steady Flows

Having derived the equations of motion and conservation laws for ideal fluids, let us now consider some important special cases.

Static Fluids

Static fluids have $\mathbf{v} = 0$ everywhere, so the fluid is at rest (implying mechanical equilibrium). Continuity then says $\frac{\partial \rho}{\partial t} = 0$, so ρ and \mathcal{P} are independent of time. If $\Phi = gz$ for gravity, the Euler equation says

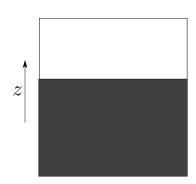
$$\frac{\nabla \mathcal{P}}{\rho} = -g\hat{z},\tag{6.52}$$

so \mathcal{P} and ρ can each only be functions of z, while in fact $\frac{1}{\rho} \frac{\partial \mathcal{P}}{\partial z}$ is independent of z.

Example if the density ρ is constant, then if $\Phi = gz$, then

$$\mathcal{P}(z) = \mathcal{P}_0 - \rho g z,\tag{6.53}$$

so the pressure decreases with height.



Example let us pretend the atmosphere is an ideal gas at a uniform constant temperature T. As $\mathcal{P} = \rho R'T$, then

$$\frac{\partial \mathcal{P}}{\partial z} = -\frac{g}{R'T}\mathcal{P},\tag{6.54}$$

SO

$$\mathcal{P}(z) = \mathcal{P}_0 e^{-\frac{gz}{R'T}}, \tag{6.55}$$

and the pressure falls exponentially.

Example the Archimedes principle says that the pressure balances the weight of any displaced fluid. This follows from our first example above. The pressure on the bottom of the object displacing the fluid is ρgz , where z is the distance from the surface. The force is pressure times area, so that is the volume displaced times ρg , or the mass displaced times g.

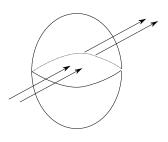


Note that if the temperature is not uniform, then the mechanical equilibrium is not stable, as the temperature gradients result in *convection* currents which mix the fluid. Therefore we have used the fact that we are discussing an ideal fluid.

Steady Flows

Steady flows are ones in which $\frac{\partial \rho}{\partial t} = 0$, $\frac{\partial \mathcal{P}}{\partial t} = 0$, and $\frac{\partial \mathbf{v}}{\partial t} = 0$ at every position \mathbf{r} .

In this case, the continuity equation becomes $\nabla \cdot (\rho \mathbf{v}) = 0$, so what flows in must flow out. This is most easily implemented by using the integral form, where for any closed surface S we have



$$\int_{S} d\mathbf{S} \cdot (\rho \mathbf{v}) = 0. \tag{6.56}$$

For a steady flow the Bernoulli equation becomes the statement that

$$\frac{\mathbf{v}^2}{2} + \Phi + \frac{\mathcal{P}}{\rho} + U = B \tag{6.57}$$

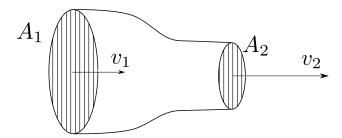
where B is a constant along the paths of fluid elements. Most often we will consider gravity where $\Phi = gz$.

A steady incompressible fluid is one where $\nabla \cdot \mathbf{v} = 0$ and U is constant. Continuity now says $\nabla \rho = 0$ so ρ is constant as well. Moreover, the Bernoulli equation now says

$$\frac{\mathbf{v}^2}{2} + \Phi + \frac{\mathcal{P}}{\rho} = B' \tag{6.58}$$

is also constant.

Example for a horizontal pipe filled with an incompressible fluid (which is approximately true of water at room temperature) of constant density ρ .



Lets consider the two ends to be at approximately at the same height so we can drop the term $\Phi = gz$. At the hatched areas shown we know that the flow must be tangential to the edge of the pipe, so a valid solution is to simply consider the velocities to be uniform and tangential to the enclosing pipe across each of these areas. Using Bernoulli this implies that

$$\frac{\rho v_1^2}{2} + \mathcal{P}_1 = \frac{\rho v_2^2}{2} + \mathcal{P}_2. \tag{6.59}$$

Furthermore, the continuity equation for the enclosed areas shown (some sides being those of the pipe) implies that the flux in at one end must equal the flux out at the other $A_1v_1 = A_2v_2$. Together this gives

$$\mathcal{P}_1 = \mathcal{P}_2 + \frac{1}{2}\rho v_2^2 \left(1 - \left(\frac{A_2}{A_1} \right)^2 \right). \tag{6.60}$$

Since $A_1 > A_2$ we have $v_2 > v_1$, and this implies $\mathcal{P}_1 > \mathcal{P}_2$.

Example let us consider a water tank filled to a height z_1 with a hole at height $z_2 < z_1$ that produces a jet of water.

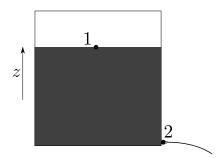
Let us assume $v_1 \approx 0$ (so the tank is much larger than the hole). Then

$$\frac{1}{2}\rho v_2^2 + \mathcal{P}_2 + \rho g z_2 = \mathcal{P}_1 + \rho g z_1. \tag{6.61}$$

Additionally, $\mathcal{P}_2 = \mathcal{P}_1 = \mathcal{P}_{\text{atmosphere}}$, so the pressure terms cancel out, and we can solve for the v_2 velocity to give

$$v_2 = \sqrt{2g(z_1 - z_2)} \,. \tag{6.62}$$

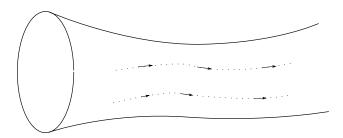
This is the same velocity as that for any mass falling from rest through a height $z_1 - z_2$. Of course a key difference for the jet of water is that this velocity is horizontal rather than vertical.



Lets discuss two common ways to picture flows. One is through *stream lines*, which are lines that are everywhere tangent to the instantaneous velocity, meaning

$$\frac{d\mathbf{x}(s)}{ds} \times \mathbf{v} = 0 \tag{6.63}$$

for some parameter s that determines the distance along the streamline. These lines are drawn at some fixed time and never cross since there is a unique velocity at every point.



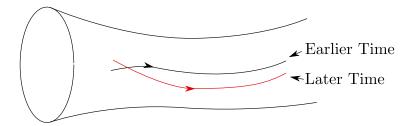
Writing out the cross product we find $0 = d\mathbf{x}/ds \times \mathbf{v} = \hat{z}(v_y \frac{dx}{ds} - v_x \frac{dy}{ds}) + \dots$ implying that we have

$$\frac{dy}{dx} = \frac{v_y}{v_x}, \qquad \frac{dz}{dx} = \frac{v_z}{v_x}, \qquad \left(\text{or } \frac{dz}{dy} = \frac{v_z}{v_y}\right).$$
(6.64)

When we use the equations in this form we would need to be able to switch from s to the variable x to uniquely parameterize the curve.

Another method of picturing the flow is through *flow lines*, which are paths that are followed by fluid elements, meaning

$$\frac{d\mathbf{x}}{dt} = \mathbf{v}(\mathbf{x}, t). \tag{6.65}$$

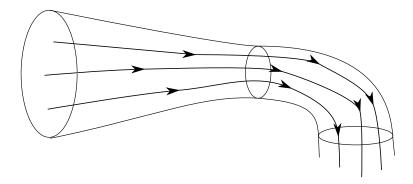


Since flow lines are time dependent, they can in general cross since the path a fluid element takes through a point may differ at a later time.

For a steady flow, the lines are time independent, and stream lines and flow lines are identical. To prove this we consider the flow line equations $dx/dt = v_x$, $dy/dt = v_y$, and note that since the velocities are time independent that we can eliminate time through the ratio $dy/dx = v_y/v_x$ (and similar for other directions), which is the equation for the stream lines. To go the opposite direction we simply pick s = t to parameterize the streamline, and note that $\mathbf{v} \times \mathbf{v} = 0$. Furthermore, for a steady fluid we have Bernoulli's law

$$\frac{\mathbf{v}^2}{2} + \Phi + \frac{\mathcal{P}}{\rho} + U = B \tag{6.66}$$

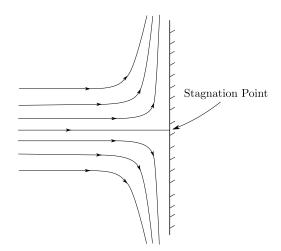
which we now understand is constant along stream lines.



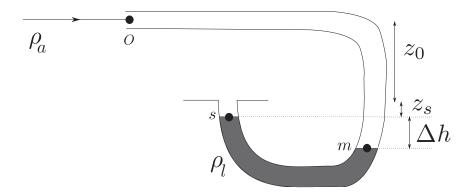
Stagnation Points and a Pitot Tube

Consider an ideal incompressible fluid in a steady flow which flows in an almost uniform manner, and then hits a wall.

Since the velocity must be tangential at the surface, the flow can either curve left or right, and it is often the case that parts of the flow will go in both directions. In this case there is a stagnation point o, where $v_o = 0$. If we consider the streamline that hits the stagnation point then $\mathcal{P}_o = \mathcal{P}_\infty + \frac{1}{2}\rho v_\infty^2$, where v_∞ and \mathcal{P}_∞ are, respectively, the velocity and pressure infinitely far away. Thus the largest pressure in the entire flow occurs at the stagnation point where there is only pressure and no kinetic energy.



Now let us consider a Pitot tube, which is a device used to measure velocity (for example on airplanes).



If the density is ρ_a in the air and ρ_l in the liquid, we can write down the Bernoulli equations obtained by comparing the air and liquid flows at ∞ , the stagnation point o, at the point s near the surface (where the air velocity is the same as at ∞ and liquid is at rest), and at the point m at the top of the liquid inside the column. This gives:

$$\mathcal{P}_{o} = \mathcal{P}_{\infty} + \frac{1}{2}\rho_{a}v_{\infty}^{2}$$

$$\mathcal{P}_{s} = \mathcal{P}_{\infty} + \frac{1}{2}\rho_{a}g(z_{o} + z_{s})$$

$$\mathcal{P}_{m} = \mathcal{P}_{o} + \rho_{a}g(z_{0} + z_{s} + \Delta h)$$

$$= \mathcal{P}_{s} + \rho_{l}q\Delta h$$

Subtracting the 2nd equation from the 1st, subtracting the 4th equation from the 3rd, and then adding these two results cancels all the pressure terms, and leaves

$$\frac{1}{2}\rho_a v_\infty^2 = g(\rho_l - \rho_a)\Delta h. \tag{6.67}$$

This can be rearranged to write

$$v_{\infty}^{2} = 2g \left(\frac{\rho_{l}}{\rho_{a}} - 1\right) \Delta h \approx 2g \frac{\rho_{l}}{\rho_{a}} \Delta h \tag{6.68}$$

allowing us to determine the velocity of the air v_{∞} in terms of the known ratio of densities $\frac{\rho_l}{\rho_a} \gg 1$ and simply the measured height between the liquid on each side, Δh .

6.4 Potential Flow

When a flow is everywhere both irrotational and incompressible it is known as *potential* flow. Such flows may be steady or not steady. Since $\Omega = \nabla \times \mathbf{v} = 0$ the velocity field is conservative. This means there exists a velocity potential

$$\phi(\mathbf{r}) = \int_{\mathbf{r}_0}^{\mathbf{r}} \mathbf{v}(\mathbf{r}', t) \cdot d\mathbf{r}'$$
 (6.69)

which depends only on the endpoints of the integration, and not the path taken between them, such that

$$\mathbf{v} = \nabla \phi \tag{6.70}$$

Since $\nabla \cdot \mathbf{v} = 0$ for an incompressible fluid, then ϕ must solve Laplace's equation

$$\nabla^2 \phi = 0. ag{6.71}$$

Solving for the scalar ϕ (with suitable boundary conditions), then immediately gives \mathbf{v} .

We can then use the Euler equation to immediately get the pressure. For $\Omega=0$ one form of the Euler equation was

$$\frac{\partial \mathbf{v}}{\partial t} + \nabla \left(\frac{\mathbf{v}^2}{2} + \Phi \right) + \frac{\nabla \mathcal{P}}{\rho} = 0.$$
 (6.72)

If ρ is constant, then

$$\nabla \left(\frac{\partial \phi}{\partial t} + \frac{\mathbf{v}^2}{2} + \Phi + \frac{\mathcal{P}}{\rho} \right) = 0, \tag{6.73}$$

SO

$$\frac{\partial \phi}{\partial t} + \frac{\mathbf{v}^2}{2} + \Phi + \frac{\mathcal{P}}{\rho} = b(t) \tag{6.74}$$

for some function b. For each t we can pick the zero of ϕ so that b(t) is constant. (This is equivalent to shifting $\phi \to \phi + \int^t b(t') dt'$, where adding this constant that is independent of \mathbf{x} gives a solution that is equally valid.) The remaining constant b be fixed by a boundary condition on the pressure. Thus the full pressure as a function of \mathbf{x} and t is determined by

$$\frac{\mathcal{P}}{\rho} = -\frac{\partial \phi}{\partial t} - \frac{\mathbf{v}^2}{2} - \Phi + b \,, \tag{6.75}$$

where in principal the first three terms on the right hand side carry both spatial and time dependence. Often we are interested in a steady flow, in which case the term $\partial \phi / \partial t = 0$.

Example if $\phi = v_0 x$, then $\mathbf{v} = v_0 \hat{x}$, which is a specific case of a uniform flow.

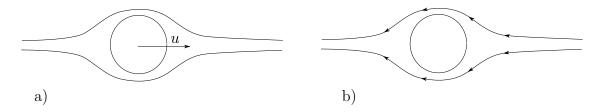
Example if $\phi = A \ln(r)$ in 2 dimensions, then for all r > 0

$$\mathbf{v} = \frac{\partial \phi}{\partial r}\hat{r} + \frac{1}{r}\frac{\partial \phi}{\partial \theta}\hat{\theta} = \frac{A}{r}\hat{r},\tag{6.76}$$

which is a point source. An analogous point source in 3 dimensions would arise from $\phi = -\frac{A}{r}$, which gives $\mathbf{v} = \frac{A}{r^2}\hat{r}$ for all r > 0. (At r = 0 there would be a delta function source, so we do not satisfy Laplace's equation at this point.)

Example Consider $\phi = \frac{\Gamma}{2\pi}\theta$ in polar coordinates in the 2 dimensional plane. Then $v_r = \frac{\partial \phi}{\partial r} = 0$ and $v_{\theta} = \frac{1}{r} \frac{\partial \phi}{\partial \theta} = \frac{\Gamma}{2\pi r}$ for all r > 0. This corresponds to a potential vortex about the point r = 0.

Example Consider a sphere of radius R moving with constant velocity \mathbf{u} through an incompressible ideal fluid and find its velocity by solving for its velocity potential. Equivalently we can consider the problem of finding the velocity when a sphere is held at rest and the fluid flows in from far away with uniform velocity $-\mathbf{u}$ at infinity. The first situation is shown in figure a) and the second as figure b).



We use coordinates centered on the sphere, and define the axes so that $\mathbf{u} = u\hat{x}$. The problem is spherically symmetric other than the directionality from \mathbf{u} , so its natural to expect $\mathbf{v} \propto \mathbf{u}$. Since $\nabla \phi = \mathbf{v}$ this means that we expect $\phi \propto \mathbf{u}$. (Effectively the boundary condition is linear in \mathbf{u} and the equation $\nabla^2 \phi = 0$ is linear. We could also explicitly demonstrate the proportionality $\mathbf{v} \propto \mathbf{u}$ using dimensional analysis, as we will discuss in more detail later in this chapter.)

It is actually easier to consider the sphere being at rest with the fluid moving past it as in b), so lets start with this case. Since $\nabla^2 \phi = 0$ is linear, we can solve using superposition. The velocity potential

$$\phi(\mathbf{x}) = -ux + \phi'(\mathbf{x}) \tag{6.77}$$

has a term -ux giving the correct uniform flow far away from the sphere. Therefore, with r the distance from the center of the sphere, we have

$$\lim_{r \to \infty} \phi'(\mathbf{x}) = 0. \tag{6.78}$$

Another way to see this is that taking the gradient of Eq. (6.77) gives

$$\mathbf{v} = -u\hat{x} + \mathbf{v}',\tag{6.79}$$

which is simply the translation between the velocity field \mathbf{v} for b) and the velocity field \mathbf{v}' for a). For the situation a) we would anticipate Eq. (6.78) as the correct boundary condition, since the fluid is at rest at infinity when it is infinitely far away from the disturbance caused by dragging the sphere. We can look for a solution for ϕ' .

As $\nabla^2 \phi' = 0$ with $\lim_{r \to \infty} \phi' = 0$, one option could be $\phi' = \frac{1}{r}$ in 3 dimensions, but this would give a point source solution with velocity moving radially outward from our sphere and hence make it impossible to satisfy the appropriate boundary condition on the sphere (its also not $\propto \mathbf{u}$). Instead, let us consider a dipole source

$$\phi' = A\mathbf{u} \cdot \nabla \left(\frac{1}{r}\right) \tag{6.80}$$

for some constant A. This proposal is linear in \mathbf{u} , and satisfies $\nabla^2 \phi' = 0$ for r > 0 since the derivatives commute: $\nabla^2 \phi' = A(\mathbf{u} \cdot \nabla) \nabla^2 (1/r) = 0$. It remains to compute ϕ' and \mathbf{v} explicitly and check that we can satisfy the boundary conditions (and proper dimensions) with this solution. Since

$$\nabla\left(\frac{1}{r}\right) = -\frac{\mathbf{r}}{r^3},\tag{6.81}$$

we have

$$\phi' = -\frac{A\mathbf{u} \cdot \mathbf{r}}{r^3}.\tag{6.82}$$

Moreover,

$$\nabla(\mathbf{u} \cdot \mathbf{r}) = \mathbf{u}, \qquad \nabla r^{-n} = -nr^{-(n+2)}\mathbf{r}, \tag{6.83}$$

so the solution for \mathbf{v}' is

$$\mathbf{v}' = \nabla \phi' = \frac{A}{r^3} \left(-\mathbf{u} + \frac{3(\mathbf{u} \cdot \mathbf{r})\mathbf{r}}{r^2} \right)$$
 (6.84)

and we then also have obtained $\mathbf{v} = -\mathbf{u} + \mathbf{v}'$. The boundary condition on the surface of the sphere (which is r = R for case b) where the sphere is at rest) is $\mathbf{v} \cdot \hat{r} = 0$. This means

$$0 = -\mathbf{u} \cdot \hat{r} - \frac{A\mathbf{u} \cdot \hat{r}}{R^3} + \frac{3A\mathbf{u} \cdot \hat{r}}{R^3},\tag{6.85}$$

which has the solution $A = \frac{R^3}{2}$. Thus,

$$\mathbf{v} = -\mathbf{u} + \frac{R^3}{2r^3} \left(\frac{3(\mathbf{u} \cdot \mathbf{r})\mathbf{r}}{r^2} - \mathbf{u} \right). \tag{6.86}$$

This solution has the right dimensions and satisfies the boundary conditions on the sphere and at infinity.

For steady flow, we can then use Bernoulli's equation to get the pressure on the sphere, constant $= (\mathcal{P} + \frac{1}{2}\rho v^2)_{r=R} = (\mathcal{P} + \frac{1}{2}\rho v^2)_{r=\infty}$. Squaring our result for the velocity on the sphere, setting $\mathbf{u} \cdot \mathbf{r} = ur \cos \theta$, and simplifying we find

$$\mathcal{P} = \mathcal{P}_{\infty} + \frac{\rho \mathbf{u}^2}{8} \left(9\cos^2 \theta - 5 \right). \tag{6.87}$$

This result for the pressure says that it is the same on the front and back of the sphere, since its unchanged by taking $\theta \to \pi/2 - \theta$. This is quite counterintuitive, since we expect a force on the sphere in b) that would try to push it downstream. This actually results from our approximation that the fluid is ideal (viscosity can not be neglected when trying to answer questions near surfaces).

Another possibility is that our approximation of potential flow is suspect. To explore this, lets ask how common is potential flow? Consider

$$\frac{d}{dt}\left(\nabla \times \mathbf{v}\right) = \nabla \times \frac{d\mathbf{v}}{dt} = \nabla \times \left(\frac{\partial \mathbf{v}}{\partial t} + \mathbf{v} \cdot \nabla \mathbf{v}\right),\tag{6.88}$$

which expands into

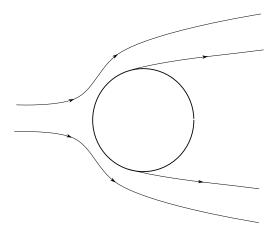
$$\frac{\partial}{\partial t} \left(\nabla \times \mathbf{v} \right) + \nabla \times \left(\nabla \left(\frac{\mathbf{v}^2}{2} \right) - \mathbf{v} \times (\nabla \times \mathbf{v}) \right) = \frac{\partial \mathbf{\Omega}}{\partial t} - \nabla \times (\mathbf{v} \times \mathbf{\Omega}) = 0 \tag{6.89}$$

where in the last step we used our result in Eq. (6.36) that we derived from the Euler equation. Thus, $\nabla \times \mathbf{v}$ is conserved along flow lines. If we consider a steady flow which starts out as uniform infinitely far away, then there is no vorticity at infinity, and

$$\nabla \times \mathbf{v} = 0 \tag{6.90}$$

on every stream line, and remains that way for the entire flow. Thus we have a potential flow with no vorticity.

If we wanted to get around the counterintuitive behavior we have found in our ideal fluid solution, but stick with the ideal fluid framework, then we would have to allow for the existence of discontinuous solutions. For an ideal fluid flowing past a sphere, we could propose that stream lines exist that start tangential to the spherical surface, and hence satisfy the boundary conditions, and can have $\nabla \times \mathbf{v} \neq 0$ since they are not connected to infinity. Behind the sphere we could then imagine having fluid at rest.



The correct treatment of boundary layers near objects and of wakes, does not require discontinuous solutions of this sort, but instead simply requires the inclusion of viscosity, which we will turn to shortly, after treating one final important example from ideal fluids.

6.5 Sound Waves

In this section we will explore an example where $\nabla \cdot \mathbf{v} \neq 0$ plays an important role. To set things up, consider a compressible fluid at rest with pressure \mathcal{P}_0 and density ρ_0 in equilibrium with an external force density \mathbf{f}_0 . If \mathcal{P}_0 and ρ_0 are constant and uniform, then

$$\frac{\nabla \mathcal{P}_0}{\rho_0} = \frac{\mathbf{f}_0}{\rho_0} \tag{6.91}$$

from the Euler equation. Now lets add disturbances \mathcal{P}' and ρ' to this system

$$\mathcal{P} = \mathcal{P}_0 + \mathcal{P}',\tag{6.92}$$

$$\rho = \rho_0 + \rho' \,, \tag{6.93}$$

with $\mathcal{P}' \ll \mathcal{P}_0$ and $\rho' \ll \rho_0$. These disturbances will induce a velocity field as well, $\mathbf{v}(\mathbf{r},t)$, which we will also assume is small, so that perturbation theory applies. We will therefore drop terms that are second order or higher in any of $\{\mathcal{P}', \rho', \mathbf{v}\}$. Using perturbation theory on the Euler equation, that is

$$0 = \rho \frac{\partial \mathbf{v}}{\partial t} + \rho (\mathbf{v} \cdot \nabla) \mathbf{v} + \nabla p - \mathbf{f}_0 = \nabla \mathcal{P}_0 - \mathbf{f}_0 + \rho_0 \frac{\partial \mathbf{v}}{\partial t} + \nabla \mathcal{P}' + \dots$$
 (6.94)

where we have dropped terms $\rho \mathbf{v} \cdot \nabla \mathbf{v} = \mathcal{O}(\rho_0 \mathbf{v}^2)$ and $\mathcal{O}(\rho' \mathbf{v})$. Using $\mathbf{f}_0 = \nabla \mathcal{P}_0$ then gives

$$\frac{\partial \mathbf{v}}{\partial t} = -\frac{\nabla \mathcal{P}'}{\rho_0} \tag{6.95}$$

as the Euler equation to 1st order in perturbations. To the same order, continuity says

$$\frac{\partial \rho'}{\partial t} = -\nabla \cdot (\rho \mathbf{v}) = -\rho_0 \nabla \cdot \mathbf{v}. \tag{6.96}$$

where we have again dropped second order terms. Finally the appropriate thermodynamic equation of state is

$$\rho' = \frac{\rho_0}{B} \mathcal{P}' \tag{6.97}$$

where B is a constant known as the bulk modulus. The bulk modulus describes a substance's resistance to compression, and this formula arises from $B = \rho \frac{\partial \mathcal{P}}{\partial \rho} \approx \rho_0 \frac{\mathcal{P}'}{\rho'}$. Using this result we can eliminate density ρ' to get a second equation involving only the pressure and velocity disturbances,

$$\frac{\partial \mathcal{P}'}{\partial t} = \frac{B}{\rho_0} \frac{\partial \rho'}{\partial t} = -B\nabla \cdot \mathbf{v} \tag{6.98}$$

Combining Eq. (6.95) and Eq. (6.98) we can derive a differential equation for the pressure disturbance

$$\frac{\partial^2 \mathcal{P}'}{\partial t^2} = -B\nabla \cdot \frac{\partial \mathbf{v}}{\partial t} = \frac{B}{\rho_0} \nabla^2 \mathcal{P}', \tag{6.99}$$

which can be written more simply as

$$\frac{\partial^2 \mathcal{P}'}{\partial t^2} - c_{\rm S}^2 \nabla^2 \mathcal{P}' = 0 \tag{6.100}$$

which is a wave equation for \mathcal{P}' , whose solutions move at a velocity $c_{\rm S} = \sqrt{\frac{B}{\rho_0}}$ which is known as the *speed of sound*. Due to the simple proportionality from the equation of state we also immediately know that

$$\frac{\partial^2 \rho'}{\partial t^2} - c_{\rm S}^2 \nabla^2 \rho' = 0 \tag{6.101}$$

so the same wave equation also holds for the density.

It remains to derive a differential equation for the velocity. Taking the curl of Eq. (6.95) yields

$$\frac{\partial}{\partial t} \left(\nabla \times \mathbf{v} \right) = -\frac{1}{\rho_0} \nabla \times \nabla \mathcal{P}' = 0. \tag{6.102}$$

This means $\nabla \times \mathbf{v}$ does not explicitly depend on time, so we can conveniently take $\nabla \times \mathbf{v} = 0$ initially everywhere, and hence for all times. Using Eq. (6.95) and Eq. (6.98) now gives

$$\frac{\partial^{2} \mathbf{v}}{\partial t^{2}} = -\frac{1}{\rho_{0}} \nabla \left(\frac{\partial \mathcal{P}'}{\partial t} \right) = \frac{B}{\rho_{0}} \nabla (\nabla \cdot \mathbf{v})
= \frac{B}{\rho_{0}} \left(\nabla^{2} \mathbf{v} - \nabla \times (\nabla \times \mathbf{v}) \right)
= \frac{B}{\rho_{0}} \nabla^{2} \mathbf{v} ,$$
(6.103)

since $(\nabla \times \mathbf{v}) = 0$. This yields

$$\frac{\partial^2 \mathbf{v}}{\partial t^2} - c_{\rm S}^2 \nabla^2 \mathbf{v} = 0 \tag{6.104}$$

which means that the velocity of the fluid disturbance also satisfies the same wave equation. The solutions are thus *sound waves* in pressure, density, and velocity, with speed $c_{\rm S}$. An example of a solution is a plane wave (here written for pressure) which looks like

$$\mathcal{P}' = \mathcal{P}'(\mathbf{r} \cdot \hat{n} - c_{S}t) \tag{6.105}$$

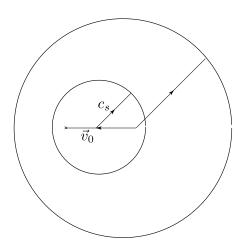
when traveling in the direction of \hat{n} .

Note that $\nabla \cdot \mathbf{v} \neq 0$ was important for this derivation. It is reasonable to ask if there is a way that we can determine when the approximation $\nabla \cdot \mathbf{v} = 0$ may be justified. For a flow with characteristic velocity v_0 this can be done by defining the *Mach number*

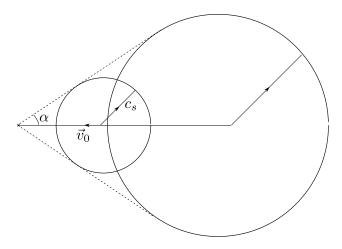
$$M = \frac{v_0}{c_{\rm S}},$$
 (6.106)

since the scaling of terms involving $\nabla \cdot \mathbf{v}$ will be determined by this ratio. If $M \ll 1$ then we can treat flow as approximately incompressible, with $\nabla \cdot \mathbf{v} = 0$.

Considering flows with large values of M leads to the concept of shock waves. Consider a flow with initial velocity v_0 in which there is a disturbance. If M < 1 then the flow is said to be subsonic, and the perturbation spreads everywhere, because the speed of the perturbation is larger than that of the flow.



On the other hand if M > 1, then the disturbance is swept downstream to the right by the flow, and actually propagates downstream within a cone of angle γ defined by $\sin(\gamma) = c_s/v_0 = \frac{1}{M}$, as shown below.



If we consider a supersonic plane, then we should view this picture the other way around, where the fluid is static and the disturbance (plane) moves through it, traveling to the left at faster than the speed of sound. This causes a sonic boom, which is the air pressure wave given by the dashed lines trailing the plane, which moves at speed $c_{\rm S}$. (Another example is thunder, where the rapid increase in temperature of plasma of ions causes rapid air expansion, making a shockwave.)

6.6 Viscous Fluid Equations

Internal friction occurs in a fluid when neighboring fluid particles move with different velocities, which means that $\frac{\partial v_i}{\partial x_j} \neq 0$. Viscous stresses oppose such relative motion. Adding these friction terms changes some of our fluid equations. In particular, the continuity equation remains unchanged, while the Euler equations along with the conservation laws for momentum, energy, and entropy must be modified.

To consider this friction we will work to first order in the $\frac{\partial v_i}{\partial x_j}$ partial derivatives, treating these as the most important terms. These derivatives can be arranged into

$$\sigma_{ij} = \eta \left(\frac{\partial v_i}{\partial x_j} + \frac{\partial v_j}{\partial x_i} - \frac{2}{3} \delta_{ij} \nabla \cdot \mathbf{v} \right) + \zeta \delta_{ij} \nabla \cdot \mathbf{v}$$
 (6.107)

to define the viscous stress tensor with elements σ_{ij} . The constant coefficients of the two terms are the shear viscosity η and the bulk viscosity ζ , where $\eta > 0 \& \zeta > 0$. The form of the viscous stress tensor σ_{ij} is dictated by the fact that it must vanish for constant \mathbf{v} and for uniform rotation $\mathbf{v} = \boldsymbol{\omega} \times \mathbf{r}$ where there is no friction. Writing out $\mathbf{v} = \omega_x(y\hat{z} - z\hat{y}) + \dots$ we see that $\nabla \cdot \mathbf{v} = 0$ and $\frac{\partial v_z}{\partial y} + \frac{\partial v_y}{\partial z} = 0$, etc., for the uniform rotation, dictating the symmetric form of the terms in σ_{ij} . The remaining organizational choice is to let η multiply a traceless tensor.

Momentum conservation still comes from

$$\frac{\partial}{\partial t} \left(\rho \mathbf{v} \right) + \nabla \cdot \mathbb{T} = \mathbf{f} \tag{6.108}$$

where now we include a friction term in the stress tensor to account for the viscous transfer of momentum. Thus

$$T_{ij} = \mathcal{P}\delta_{ij} + \rho v_i v_j - \sigma_{ij} \tag{6.109}$$

is the new total stress tensor.

A simple rule for incorporating σ_{ij} is simply to replace $\mathcal{P}\delta_{ij} \to \mathcal{P}\delta_{ij} - \sigma_{ij}$. With this we can add friction to the Euler equation. In particular we have

$$(\nabla \mathcal{P})_i = \frac{\partial}{\partial x_k} \delta_{ki} \mathcal{P} \to \frac{\partial}{\partial x_k} (\delta_{ki} \mathcal{P} - \sigma_{ki}), \qquad (6.110)$$

where we can compute that

$$\frac{\partial}{\partial x_k} \sigma_{ki} = \eta \left(\frac{\partial^2 v_i}{\partial x_k \partial x_k} + \frac{\partial}{\partial x_i} \nabla \cdot \mathbf{v} - \frac{2}{3} \frac{\partial}{\partial x_i} \nabla \cdot \mathbf{v} \right) + \zeta \frac{\partial}{\partial x_i} \nabla \cdot \mathbf{v}
= \eta \nabla^2 v_i + \left(\zeta + \frac{\eta}{3} \right) \frac{\partial}{\partial x_i} \nabla \cdot \mathbf{v} .$$
(6.111)

Plugging this into the Euler equation yields

$$\frac{\partial \mathbf{v}}{\partial t} + (\mathbf{v} \cdot \nabla)\mathbf{v} + \frac{\nabla \mathcal{P}}{\rho} - \frac{\eta}{\rho}\nabla^2\mathbf{v} - \frac{1}{\rho}\left(\zeta + \frac{\eta}{3}\right)\nabla(\nabla \cdot \mathbf{v}) = \frac{\mathbf{f}}{\rho}$$
(6.112)

which is the Navier-Stokes equation.

A common case we will study is when $\mathbf{f} = 0 \& \nabla \cdot \mathbf{v} = 0$, which reduces the Navier-Stokes equation to

$$\frac{\partial \mathbf{v}}{\partial t} + \mathbf{v} \cdot \nabla \mathbf{v} = -\frac{\nabla \mathcal{P}}{\rho} + \nu \nabla^2 \mathbf{v}$$
 (6.113)

where

$$\nu \equiv \frac{\eta}{\rho} \tag{6.114}$$

is the *kinematic viscosity*, and the bulk viscosity term has dropped out. The dimensions of the kinematic viscosity are $[\nu] = m^2/s$, which is simpler than $[\eta] = kg/(ms)$.

Other useful equations can be derived for the situation where $\nabla \cdot \mathbf{v} = 0 \& \rho$ is constant. Taking the divergence of the Navier-Stokes equation removes both the $\partial \mathbf{v}/\partial t$ term, and the $\nu \nabla^2 \mathbf{v}$ term, leaving $\rho \nabla \cdot [(\mathbf{v} \cdot \nabla)\mathbf{v}] = -\nabla^2 \mathcal{P}$. Writing this out in components we find

$$\nabla^2 \mathcal{P} = -\rho \frac{\partial}{\partial x_j} v_i \frac{\partial}{\partial x_i} v_j = -\rho \frac{\partial v_i}{\partial x_j} \frac{\partial v_j}{\partial x_i}, \qquad (6.115)$$

since $\partial v_j/\partial x_j = \nabla \cdot \mathbf{v} = 0$. This equation can be used to compute the pressure if the velocity is determined, since it simply acts like a source term. Taking the curl of the Navier-Stokes equation, and recalling that $\Omega = \nabla \times \mathbf{v}$ we find

$$\frac{\partial \mathbf{\Omega}}{\partial t} - \nabla \times (\mathbf{v} \times \mathbf{\Omega}) = \nu \nabla^2 \mathbf{\Omega} \,, \tag{6.116}$$

where the algebra to arrive at the terms on the LHS was consider earlier in our discussion of the Euler equation, and the new pieces is the term on the RHS.

In the presence of viscosity the boundary conditions change from what we had previously. Molecular forces between the viscous fluid & any surface mean that the fluid adheres to the surface, and hence that the velocity of the fluid and surface must be the same for both the tangential and longitudinal components. Therefore the boundary condition for a moving wall is $\mathbf{v} = \mathbf{v}_{\text{wall}}$, which also covers the case $\mathbf{v} = 0$ for a wall at rest.

Another important concept is the force exerted by the fluid on a surface. This has a contribution both from the pressure as well as from the friction. The force per unit area \mathcal{F}_i is given by

$$\mathcal{F}_i = -n_j(\mathcal{P}\delta_{ji} - \sigma_{ji}) = -\mathcal{P}n_i + \sigma_{ij}n_j, \qquad (6.117)$$

where \hat{n} is the normal vector pointing out of the surface, and the first term is the pressure acting along this normal vector, while the second is the friction that has tangential components.

Starting with the Navier-Stokes equation we can also derive a modified form for energy conservation. Rather than carrying out this derivation explicitly, we will just examine the final result in integral form, which is a bit more intuitive:

$$\frac{\partial}{\partial t} \int_{V} \left(\frac{1}{2} \rho \mathbf{v}^{2} \right) dV = -\oint_{\partial V} \left(\rho \left(\frac{\mathbf{v}^{2}}{2} + \frac{\mathcal{P}}{\rho} \right) v_{i} - v_{j} \sigma_{ji} \right) d\mathcal{S}_{i} - \int_{V} \sigma_{ij} \frac{\partial v_{i}}{\partial x_{j}} dV. \tag{6.118}$$

Here the term on the LHS is the change of the kinetic energy in the volume V with time. The first integral on the RHS is the energy flux through the closed bounding surface ∂V , and the second integral is the decrease in energy that is caused by dissipation. To see this even more clearly we can consider integrating over the whole fluid with $\mathbf{v} = 0$ at ∞ (or on ∂V). This removes the flux term and leaves

$$\frac{\partial E}{\partial t} = -\int_{V} \eta \left(\frac{\partial v_{i}}{\partial x_{j}} + \frac{\partial v_{j}}{\partial x_{i}} \right) \frac{\partial v_{i}}{\partial x_{j}} dV = -\frac{\eta}{2} \int_{V} \left(\frac{\partial v_{i}}{\partial x_{j}} + \frac{\partial v_{j}}{\partial x_{i}} \right)^{2} dV < 0.$$
 (6.119)

where we can check the second equality by squaring and manipulating the summed over dummy indices i and j. Thus we see that friction causes energy to dissipate just as we would expect (and this also justifies our sign choice of $\eta > 0$).

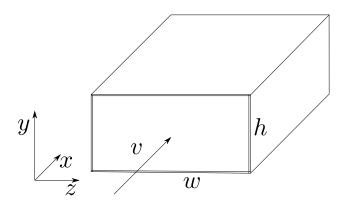
Entropy conservation is modified at temperature T to $\rho T\dot{S} = \sigma_{ij}\frac{\partial v_i}{\partial x_j}$, where the left-hand side of the equation is the heat gain per unit volume, and the right-hand side of the equation is the energy dissipated by viscosity. (If we allow thermal conduction (from temperature gradients) in the fluid, then there is another term on the right-hand side that appears as $\nabla \cdot (\kappa \nabla T)$ for conductivity κ .)

6.7 Viscous Flows in Pipes and Reynolds Number

We start our investigation of fluid flow with viscosity, by studying steady flows in pipes, typically caused by a pressure gradient. This type of steady viscous flow is also called Poiseuille flow, after Jean Poiseuille who first studied them experimentally. We take the flow to be incompressible, $\nabla \cdot \mathbf{v} = 0$, and hence the continuity equation implies that ρ is constant (just as it did in the ideal fluid case). Thus the Navier-Stokes equation for such flows reduces to

$$\mathbf{v} \cdot \nabla \mathbf{v} = -\frac{\nabla \mathcal{P}}{\rho} + \nu \nabla^2 \mathbf{v} \tag{6.120}$$

Example Lets start by considering flow in a long rectangular duct, aligned with the x-axis, with height h and width w. We also take it be a thin duct with $h \ll w$, and with a pressure gradient along the pipe in the x-direction.



This can be approximated by infinite parallel plates, taking $w \to \infty$ and holding h fixed. Here

$$\mathbf{v} = v_x(y)\,\hat{x},\tag{6.121}$$

since there is no dependence on x or z by the translational symmetry in the plane, and we equally well can not develop velocity components v_y or v_z due to this symmetry. This implies that the term $\mathbf{v} \cdot \nabla \mathbf{v} = v_x \frac{\partial \mathbf{v}}{\partial x} = 0$ Taking the inner product of Eq.(6.120) with \hat{y} and \hat{z} then removes all the terms that depends on the velocity (which is only in the \hat{x} direction), giving

$$\frac{\partial \mathcal{P}}{\partial y} = \frac{\partial \mathcal{P}}{\partial z} = 0. \tag{6.122}$$

Thus the pressure $\mathcal{P} = \mathcal{P}(x)$ and can have a gradient only in the \hat{x} direction. Taking the inner product of Eq.(6.120) with \hat{x} gives

$$\frac{\partial \mathcal{P}}{\partial x} = \rho \nu \nabla^2 v_x = \eta \frac{\partial^2 v_x}{\partial y^2} = k \,, \tag{6.123}$$

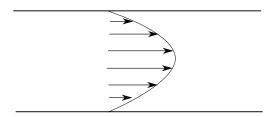
where we have introduced a constant k. Since $\frac{\partial \mathcal{P}}{\partial x}$ only depends on x, while $\eta \frac{\partial^2 v_x}{\partial y^2}$ only depends on y, they must both be equal to a constant. Let us say $\frac{\partial \mathcal{P}}{\partial x} = k < 0$, so that the pressure drops as we move in the \hat{x} direction. (In the Navier-Stokes equation, this pressure drop balances the viscous stress term.) Then integrating the equation for $v_x(y)$ gives

$$v_x(y) = \frac{k}{2\eta} y^2 + ay + b, \qquad (6.124)$$

where we have introduced two integration constants a and b. To solve for a and b we impose the boundary conditions that the velocity must vanish at y = 0 and y = h, giving

$$v_x(y) = \frac{(-k)}{2\eta} y(h - y). \tag{6.125}$$

Recalling that k < 0 we see that $v_x(y) > 0$, flowing down the pipe. (If we had reversed k the flow would be in the opposite direction.) The velocity field we have derived flows in the pipe with a parabolic profile with its maximum in the center:

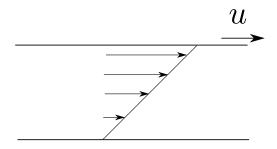


Lets also calculate the friction force per unit area that the fluid exerts on the pipe wall. The bottom plate at y = 0 has a unit vector $\hat{n} = \hat{y}$, so from Eq. (6.117) the force along \hat{x} is

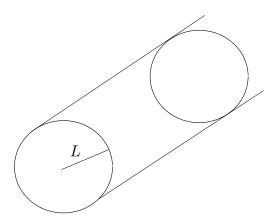
$$\mathcal{F}_x = \sigma_{yx} = \eta \frac{\partial v_x}{\partial y}|_{y=0} = -\frac{hk}{2} > 0. \tag{6.126}$$

Intuitively this is the direction we expect, the fluid tries to drag the pipe along with it.

Example Lets now consider the same example of fluid between infinite parallel plates, but now with no pressure gradient. Instead we take the top plate to move with velocity $\mathbf{u} = u\hat{x}$. Here $\mathcal{P} = \mathcal{P}_0$ is constant and uniform with k = 0, so the second derivative of $v_x(y)$ is zero, $\nabla^2 v_x = 0$, and the solution for $v_x(y)$ can at most be linear. The solution for this case is $v_x(y) = \frac{uy}{h}$, which satisfies the boundary conditions $v_x(0) = 0$ and $v_x(h) = u$. Thus the fluids velocity field is linear for this case:



Example Next consider a long cylindrical pipe of radius L, oriented along \hat{x} , again with a pressure gradient along the pipe. We will approximate the pipe as being infinitely long so there is a translational symmetry along x.



Due to the translational symmetry we know that $\mathbf{v} = \mathbf{v}(y, z)$. To fully exploit the consequences of the symmetry it is useful to use cylindrical coordinates (x, r, θ) so we can also easily impose the rotational symmetry about \hat{x} to conclude $\mathbf{v} = \mathbf{v}(r)$. The fact that there is r dependence makes sense since we know that \mathbf{v} must vanish at the edge of the pipe, r = L, but we do not want it to vanish everywhere. Continuity and symmetry also imply that the velocity is only in the \hat{x} direction, so in fact

$$\mathbf{v} = v_x(r)\hat{x} \,. \tag{6.127}$$

For example, consider an annulus shaped closed surface formed by the region between two cylinders cocentric with the pipe. The flow into and out of this surface must be balanced by continuity. The flow on the ends of the surface automatically balance each other since \mathbf{v} is independent of x. But since \mathbf{v} is r dependent, the only way the flow through the circular sides can balance each other is if there is no flow in the \hat{r} direction.

With this setup we can again confirm that $(\mathbf{v} \cdot \nabla)\mathbf{v} = v_x(r)\frac{\partial}{\partial x}v_x(r)\hat{x} = 0$, leaving $\nabla \mathcal{P} = \eta \nabla^2 \mathbf{v}$ from the Navier-Stokes equation. Taking the inner product with \hat{y} and \hat{z} we see that $\frac{\partial \mathcal{P}}{\partial y} = \frac{\partial \mathcal{P}}{\partial z} = 0$, so $\mathcal{P} = \mathcal{P}(x)$. Taking the inner product with \hat{x} gives

$$\frac{\partial \mathcal{P}(x)}{\partial x} = \eta \nabla^2 v_x(r) = k \tag{6.128}$$

where since $\frac{\partial \mathcal{P}}{\partial x}$ depends only on x while $\eta \nabla^2 v_x(r)$ is independent of x, the two must be equal to a constant k. Again we choose k < 0 to have a pressure gradient that pushes the fluid down the pipe in the \hat{x} direction (k > 0 would simply reverse the flow). For the velocity this gives

$$\frac{1}{r}\frac{\partial}{\partial r}\left(r\frac{\partial v_x}{\partial r}\right) = \frac{k}{\eta},\tag{6.129}$$

and integrating this gives

$$v_x(r) = \frac{k}{4\eta}r^2 + a\ln\left(\frac{r}{r_0}\right) + b \tag{6.130}$$

for some constants a and b. (The constant r_0 is introduced to make the equation dimensionfully correct, but is not independent, since any change to r_0 can be compensated by a change to b). Since $v_x(r)$ has to be finite at r=0 we must have a=0 (if the geometry excluded the region at the middle). The condition $v_x=0$ at r=L fixes b so that

$$v_x(r) = \frac{(-k)}{4\eta} \left(L^2 - r^2 \right) . \tag{6.131}$$

Lets calculate the discharge rate of fluid in such a pipe, as mass per unit time. This is given by

discharge rate =
$$\rho \int_0^L 2\pi r v_x(r) dr = -\frac{\pi k L^4}{8\nu} > 0$$
. (6.132)

Note that this rate is quite sensitive to the radius, being proportional to L^4 . This is why you don't want narrow blood vessels.

Reynolds Number and Similarity

How do we quantify the importance of viscosity? Let us consider a flow and identify a characteristic length L (examples of which might be the radius of a pipe or the size of a sphere inserted in a flow) and a characteristic speed u (examples of which might be the fluid velocity in the middle of a pipe or the fluid velocity upstream from flow around a sphere). The quantity

$$R = \frac{uL}{\nu} = \frac{\rho uL}{\eta} \tag{6.133}$$

is the dimensionless Reynolds number. A large R means viscosity is not as important compared to the characteristic u and L of the system, while small R means that viscosity plays an important role. It is important to note that viscosity is always important near a surface where it modifies the boundary condition on \mathbf{v} . (The above examples of Poiseuille flow have $R \lesssim 2000$.)

The introduction of R leads nicely into a discussion of the utility of exploiting dimensional analysis in fluid dynamics problems. This is best explained by way of examples.

Example Consider a steady incompressible flow in a system that is characterized by a single length scale L and initial velocity u. What could the possible functional form be for the velocity and pressure in such a system? Considering the coordinates and fluid field variables we can only have

$$\frac{\mathbf{v}}{u} = \mathbf{h} \left(\frac{\mathbf{r}}{L}, R \right), \qquad \frac{\mathcal{P}}{\rho u^2} = g \left(\frac{\mathbf{r}}{L}, R \right). \tag{6.134}$$

Here **h** and g are a dimensionless vector and scalar function respectively. The ratio \mathbf{r}/L is dimensionless, as is R, and the dimensions of \mathbf{u} and \mathcal{P} are compensated by u and ρu^2 respectively. Note that if we consider flows that have the same R, then the properties of those flows are related by changes to the scales associated with \mathbf{v} , \mathbf{r} , or \mathcal{P} . Such flows are called *similar*.

Example Consider a viscous flow past a sphere of radius a with initial velocity given by $\lim_{x\to-\infty} \mathbf{v} = u\hat{x}$. Here the Reynolds number is $R = \frac{u}{a}\nu$. If we double ν and double u then R is unchanged. Due to the relations in Eq. (6.134) we thus can predict that we will have the exact same form for the solutions with \mathbf{v} twice as large as before, and \mathcal{P} being four times as large as before.

Note that in general other dimensionless ratios, like the ratio of two length scales, or the Mach number $M=\frac{u}{c_{\rm S}}$ could also appear. (For $M\ll 1$ we treat the fluid as incompressible and neglect M for the dimensional analysis.) To determine how many independent dimensionless ratios can be formed for the analysis of a general dimensional analysis problem, one counts the number of variables and subtracts the number of unrelated types of dimensions that appear in these variables. For most fluid problems this will mean subtracting three for kg, meters, and seconds.

6.8 Viscous Flow Past a Sphere (Stokes Flow)

Lets consider a steady, incompressible, viscous fluid with small R flowing past a sphere of radius a. The fluid has velocity \mathbf{u} when it is far away from the sphere. This is the same problem we treated for an ideal fluid with potential flow, but now we want to consider the impact of viscosity, and resolve the puzzle we found in our solution for an ideal fluid. To make the problem solvable we work in the limit $R \ll 1$.

Here the Navier-Stokes equation is

$$\mathbf{v} \cdot \nabla \mathbf{v} = -\frac{\nabla \mathcal{P}}{\rho} + \frac{\eta}{\rho} \nabla^2 \mathbf{v} \,. \tag{6.135}$$

Lets determine the relative importance of the two velocity terms in the $R \ll 1$ limit. Using a to characterize spatial variations, and u to characterize velocity we find

$$\mathbf{v} \cdot \nabla \mathbf{v} = \mathcal{O}\left(\frac{u^2}{a}\right), \qquad \frac{\eta}{\rho} \nabla^2 \mathbf{v} = \mathcal{O}\left(\frac{\eta u}{\rho a^2}\right) = \mathcal{O}\left(\frac{u^2}{aR}\right).$$
 (6.136)

Therefore the viscosity term, which is enhaced by a factor of 1/R, dominates. Neglecting the $\mathbf{v} \cdot \nabla \mathbf{v}$ term, the Navier-Stokes equation reduces to

$$\nabla \mathcal{P} = \eta \nabla^2 \mathbf{v} \,. \tag{6.137}$$

Note that when written in terms of the shear viscosity η , that the density ρ has dropped out of this equation, and hence the constant ρ will not play a part in the solution. Taking the

divergence of this equation we find $\nabla^2 \mathcal{P} = \eta \nabla^2 \nabla \cdot \mathbf{v} = 0$, so the pressure satisfies Laplace's equation. Using dimensional analysis for the pressure we expect a solution of the form

$$\mathcal{P} = \frac{\eta \mathbf{u} \cdot \mathbf{r}}{a^2} g\left(\frac{\mathbf{r}}{a}\right),\tag{6.138}$$

where the dimensional analysis requirement of having only a single factor of the velocity \mathbf{u} and a scalar result, leads to including the factor of $\mathbf{u} \cdot \mathbf{r}$. Note that we have not included R as a possible argument since we are expanding with $R \ll 1$. Due to the fact that \mathcal{P} must satisfy Laplace's equation and is proportional to $\mathbf{u} \cdot \mathbf{r}$, we can immediately recognize the dipole solution (which we met earlier for the potential ϕ in the ideal fluid case):

$$\mathcal{P} = \mathcal{P}_0 + k \frac{\eta a \mathbf{u} \cdot \mathbf{r}}{r^3} \,. \tag{6.139}$$

Here we have included an overall constant pressure \mathcal{P}_0 to satisfy any boundary condition on the pressure at $r \to \infty$, and a dimensionless constant k which is still to be determined. Note that with $\mathbf{u} \cdot \mathbf{r} = ur \cos \theta$ that the pressure is no longer symmetric on the front and back of the sphere (with respect to the incoming flow), thus qualitatively differing from the ideal fluid case.

Next we note that the vorticity of the fluid satisfies

$$\nabla \times \mathbf{\Omega} = \nabla \times (\nabla \times \mathbf{v}) = \nabla(\nabla \cdot \mathbf{v}) - \nabla^2 \mathbf{v} = -\nabla^2 \mathbf{v} = -\frac{1}{\eta} \nabla \mathcal{P}$$
$$= -\frac{ka}{r^3} \left(\mathbf{u} - \frac{3(\mathbf{u} \cdot \mathbf{r})\mathbf{r}}{r^2} \right). \tag{6.140}$$

Enforcing that Ω vanishes as $r \to \infty$, the solution for this equation is

$$\Omega = ka \frac{\mathbf{u} \times \mathbf{r}}{r^3} \,. \tag{6.141}$$

Thus we see that unlike the ideal fluid case, there is now a non-zero vorticity in the fluid here.

Next we turn to determining the velocity, which can only depend on $\mathbf{v} = \mathbf{v}(\mathbf{r}, \mathbf{u}, a)$. Again we do not include R as an argument since we are expanding for $R \ll 1$. By dimensional analysis the velocity must be linear in \mathbf{u} so the most general possible solution takes the form

$$\mathbf{v} = \mathbf{u} f\left(\frac{r}{a}\right) + \frac{\mathbf{r}(\mathbf{u} \cdot \mathbf{r})}{a^2} g\left(\frac{r}{a}\right), \tag{6.142}$$

where the functions f and g are dimensionless. The gradient of one of these dimensionless functions gives $\nabla f(r/a) = (\mathbf{r}/(ar))f'(r/a)$. Computing the divergence of the velocity with this form we have

$$0 = \nabla \cdot \mathbf{v} = \frac{\mathbf{u} \cdot \mathbf{r}}{ar} f'\left(\frac{r}{a}\right) + \frac{(\mathbf{u} \cdot \mathbf{r})r}{a^3} g'\left(\frac{r}{a}\right) + \frac{4(\mathbf{u} \cdot \mathbf{r})}{a^2} g\left(\frac{r}{a}\right). \tag{6.143}$$

Therefore we find that the scalar functions must satisfy

$$f'\left(\frac{r}{a}\right) = -\frac{r^2}{a^2}g'\left(\frac{r}{a}\right) - \frac{4r}{a}g\left(\frac{r}{a}\right). \tag{6.144}$$

Next we equate the Laplacian of the velocity and gradient of the pressure, which can be simplified to give

$$\nabla^{2}\mathbf{v} = \left(\nabla^{2}f + \frac{2g}{a^{2}}\right)\mathbf{u} + \left(\nabla^{2}g + \frac{4g'}{ar}\right)\frac{(\mathbf{u} \cdot \mathbf{r})\mathbf{r}}{a^{2}}$$

$$= \frac{1}{\eta}\nabla\mathcal{P} = \frac{ka}{r^{3}}\mathbf{u} + \left(-\frac{3ka^{3}}{r^{5}}\right)\frac{(\mathbf{u} \cdot \mathbf{r})\mathbf{r}}{a^{2}}.$$
(6.145)

Note that here $\nabla^2 g(r/a) = (1/r^2)(d/dr)r^2(d/dr)g(r/a)$. Equating the coefficients of the two structures we find

$$\nabla^2 g + \frac{4g'}{ar} = \frac{g''}{a^2} + \frac{6g'}{ar} = -\frac{3ka^3}{r^5},$$

$$\frac{f''}{a^2} + \frac{2f'}{ar} + \frac{2g}{a^2} = \frac{ka}{r^3}.$$
(6.146)

To solve the equation for g we try a polynomial solution of the form $g(x) = C_n x^n$ giving

$$C_n[n(n-1)+6n]x^{n-2} = -3kx^{-5}.$$
 (6.147)

Here n = -3 is a particular solution to the full inhomogeneous equation with $C_{-3} = k/2$. Also n = 0 and n = -5 are homogeneous solutions where the LHS vanishes, and the corresponding coefficients C_0 and C_5 must be fixed by boundary conditions. Looking back at our starting point in Eq. (6.142) we see that the full set of boundary conditions are

$$\mathbf{v} = \mathbf{u} \text{ at } r = \infty : \qquad \lim_{r \to \infty} r^2 g(r/a) \to 0, \qquad \lim_{r \to \infty} f(r/a) \to 1,$$

$$\mathbf{v} = 0 \text{ at } r = a : \qquad g(1) = 0, \qquad f(1) = 0, \qquad (6.148)$$

This fixes $C_0 = 0$ and $C_{-5} = -k/2$ so that

$$g(r/a) = \frac{ka^3}{2r^3} \left(1 - \frac{a^2}{r^2} \right). \tag{6.149}$$

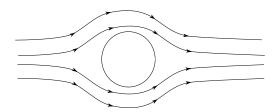
Using Eq. (6.144) and integrating once we find that

$$f(r/a) = 1 + \frac{ka}{2r} + \frac{ka^3}{6r^3}, (6.150)$$

where we have set the integration constant to one to satisfy the boundary condition at $r = \infty$. The final boundary condition, f(1) = 1 then requires us to take the constant k = -3/2. Note that this fixes the constant k that appeared in the vorticity Ω and in the pressure \mathcal{P} . All together we have that the final solution for the velocity is

$$\mathbf{v} = \mathbf{u} \left(1 - \frac{3a}{4r} - \frac{a^3}{4r^3} \right) - \frac{3a^3}{4r^3} \frac{\mathbf{r}(\mathbf{u} \cdot \mathbf{r})}{a^2} \left(1 - \frac{a^2}{r^2} \right). \tag{6.151}$$

The flow looks like:



Next we turn to determining the drag force on the sphere. In general the drag force on an object in the direction j is given by an integral of the force per unit area over the surface,

$$\mathbf{F}_{\mathrm{D}j} = \int_{\partial V} d\mathcal{S}_i \left(\mathcal{P} \delta_{ij} - \sigma_{ij} \right). \tag{6.152}$$

Lets take the inflowing velocity to be in the \hat{x} direction, $\mathbf{v}(r \to \infty) = u\hat{x}$. Then $\mathbf{F}_{\mathrm{D}} \cdot \hat{x}$ will be the drag force on our sphere in the direction of the bulk fluid flow. With spherical coordinates (r, θ, ϕ) where θ is the polar angle, we have $d\mathbf{S} = \hat{r}a^2d\cos\theta d\phi$ with r = a, as well as $\hat{r} \cdot \hat{x} = \cos\theta$ and $\hat{\theta} \cdot \hat{x} = -\sin\theta$. Thus

$$\hat{x} \cdot \mathbf{F_D} = a^2 \int d\cos\theta \, d\phi \left(-\mathcal{P}\cos\theta + \sigma_{rr}\cos\theta - \sigma_{r\theta}\sin\theta \right). \tag{6.153}$$

Computing the needed components on the sphere r = a we find

$$\sigma_{rr} = 2\eta \frac{\partial v_r}{\partial r}\Big|_{r=a} = 0, \qquad \sigma_{r\theta} = \eta \left(\frac{1}{r} \frac{\partial v_r}{\partial \theta} + \frac{\partial v_{\theta}}{\partial r} - \frac{v_{\theta}}{r}\right)_{r=a} = -\left(\frac{3\eta}{2a}\right) u \sin \theta,$$

$$\mathcal{P} = \mathcal{P}_0 - \left(\frac{3\eta}{2a}\right) u \cos \theta. \qquad (6.154)$$

Thus the drag force on the sphere is

$$\hat{x} \cdot \mathbf{F_D} = \left(\frac{3\eta u}{2a}\right) \int d\cos\theta \, d\phi \, a^2 = (6\pi\eta a)u \,, \tag{6.155}$$

which is the famous Stoke's formula for the viscous (friction) drag force on a sphere. (Note that we could have obtained the factor of ηau by dimensional analysis.)

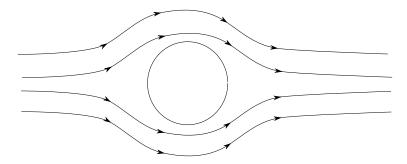
In addition to drag forces like this, that point in the direction of the fluid flow, objects may also experience *lift forces* that are tangential to direction of the fluid flow. Such forces

occur for wing-shaped objects and are important for many physical phenomena, including lift on airplanes.

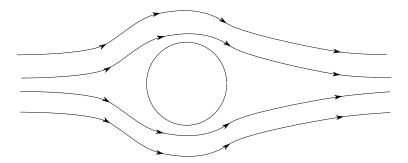
Dynamic Vortices and Turbulence

For our flow about the fixed sphere, lets consider what happens as we increase R. From our analysis above it is clear that at some point the non-linear $\mathbf{v} \cdot \nabla \mathbf{v}$ term we dropped will become important. The $\partial \mathbf{v}/\partial t$ will also become important, with flows that are more dynamical, changing with time. Lets consider how the flow appears for various values of R:

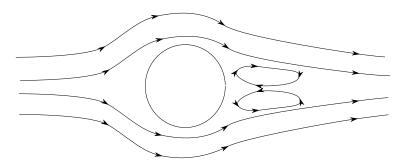
• For $R \ll 1$, the flow is symmetric and is (somewhat counter intuitively) qualitatively like the case of $\eta \to 0$. This is also called "Stokes flow" or a "laminar flow".



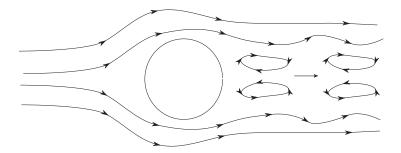
• For $R \approx 1$, the flow is still like Stokes flow, but the stream lines are no longer as symmetric, with a more clear wake developing behind the sphere.



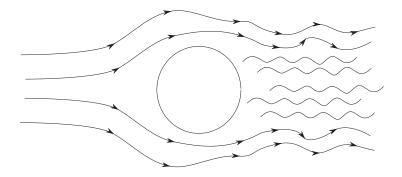
• For $10 \lesssim R \lesssim 100$, detached vortices called *eddies* form behind the sphere, though the flow is still steady. Note that directly behind the sphere between the vortices that the fluid is now flowing in the opposite direction to the asymptotic inflow \mathbf{u} . As R increases, the flow becomes looses its steady nature, with the time dependence emerging by having through oscillations of the vortices.



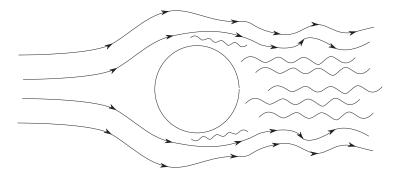
• For $150 \lesssim R \lesssim 1000$, vortices start to be cyclically shed and drift downstream in a wake behind the sphere. This time dependent solution appears like it has interaction between the eddies, where one pair pushes the next downstream.



• For $10^3 \lesssim R \lesssim 2 \times 10^5$, the wake becomes highly irregular, exhibiting a phenomena known as turbulence which we will discuss in more detail below. Here there are unsteady, interacting vortices at all length scales.



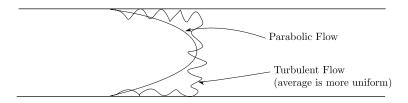
• For $R \gtrsim 2 \times 10^5$, the turbulent wake narrows and the boundary layer around the sphere is no longer laminar, also becoming turbulent.



Turbulence is characterized by a flow that: i) is highly irregular (technically chaotic) in \mathbf{x} and/or t, ii) has nonzero vorticity, iii) has much greater dissipation of energy compared to more uniform laminar viscous flows, and iv) has eddies and vortices forming over many length scales with energy that in a three dimensional flow cascades from the largest eddies down to the smallest eddies where it dissipates into heat due to viscous friction. Turbulent mixing is a very effective mechanism of transport for energy, momentum, temperature, and so on. Examples of turbulence include many familiar phenomena: the circulation of air in the atmosphere, the water flow from a faucet which becomes white at a certain point when the flow rate is increased, water in rapids, dust kicked up by wind, water beside a ship moving in an otherwise smooth lake, clear air turbulence causing a drop in lift for airplanes, and so on.

In the last value of R discussed for flow around our shere, a turbulent boundary layer appeared. This causes an abrupt drop in drag in the flow over objects, and is a very useful phenomena. In particular, by introducing imperfections we can cause this turbulent boundary layer to form at smaller values R, meaning smaller velocities. This is why golf balls have dimples and baseballs and tennis balls have visible seams.

We also get turbulence in flow through pipes at large R. The viscous flow in pipes we previously considered were laminar flow at smaller R values and had velocity distributions that were parabolic, while in contrast a turbulent flow will be non-uniform at small scales, but when averaged causes a more uniform flow down the pipe at larger length scales.



In general the chaotic and irregular nature of turbulence makes it difficult to treat with analytic methods, and a complete description of turbulence remains an unsolved problem.