Thesis Title

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E. B. Legrand



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MASTER OF SCIENCE THESIS

For the degree of Master of Science in Systems and Control at Delft University of Technology

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Delft University of Technology Department of Delft Center for Systems and Control (DCSC)

The undersigned hereby certify that they have read and recommend to the Faculty of Mechanical, Maritime and Materials Engineering for acceptance a thesis entitled

Thesis Title

by

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in partial fulfillment of the requirements for the degree of Master of Science Systems and Control

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Abstract

This is an abstract.

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Preface

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Acknowledgements

I would like to thank my supervisor prof.dr.ir. M.Y. First Reader for his assistance during the writing of this thesis. . .

By the way, it might make sense to combine the Preface and the Acknowledgements. This is just a matter of taste, of course.

Delft, University of Technology May 17, 2022 E. B. Legrand

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Chapter 1

Introduction

Original Liouville ideas:

- Showcase complex behaviour using the van der Pol oscillator
- (Optimal) control of the distributions using the Brockett approach
- (Stochastic) inputs, link with Langevin equations
- Liouville thing (in continuity form, not incompressibility) can be applied to any diff. eq.
- Bayesian inversion of chaotic systems; guess the initial state by sampling after a certain time
- Define as streamtube, continuity equation asserts that streamlines cannot cross; i.e. streamtubes are conserves. To reduce computational complexity, define level sets (curves in 2-D) and check how they deform through the evolution of the phase space fluid; should always contain the same amount of probability troughout the evolution of the system.

Notation check

Object	Roman lower	Roman upper	Greek lower	Greek upper
Standard	abcde	ABCDE	αβγδε	ΓΔΥΩΘ
Vector	abcde	ABCDE	αβγδε	ΓΔΥΩΘ
Tensor	abcde	ABCDE	αβγδε	ΓΔΥΩΘ

Table 1-1: Caption

Math constants: $ie\pi$ Variation: δS

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Musical isomorphism

Flat: X^{\flat} Sharp: ω^{\sharp}

Lie derivative: $\pounds_X H$ Interior product: $X \sqcup \omega$ Lowercase mathcal:

Kinematic momentum: pp

 $E \xrightarrow{\pi} B$

About mathematical notation and sign conventions

For symplectic geometry, the sign convention used by Abraham and Marsden [1] and Cannas da Silva [2] is observed — not the one used by Arnol'd in his *Mathematical methods of classical mechanics*, nonetheless often referred to in this text.

- Matrices, vectors and tensors are bold upper case.
- Differential forms are typically denoted by Greek letters, with their rank as a superscript (cf. Arnol'd).

Chapter 2

A Differential-Geometric Perspective on Economic Engineering

Symplectified Contact Mechanics for Dissipative Systems

The traditional view is that the methods of analytical mechanics, such as the Lagrangian and Hamiltonian formalisms, are only suited for conservative systems. However, several attempts, especially in the previous century, have been made to extend these principles to dissipative systems as well.

3-1 The damped harmonic oscillator

This chapter (and the application in the following chapter) is primarily concerned with the prototypical dissipative mechanical system: the linearly damped harmonic oscillator depicted in fig. 3-1, with the governing second-order differential equation being

$$m\ddot{x} + b\dot{x} + kx = 0. \tag{3-1}$$

The choice for this system is rather perspicuous, since it is arguably the 'easiest' dissipative system that also exhibits second-order dynamics and is linear in all terms. Furthermore, as discussed below, it serves as the test case of the overwhelming majority of research into dissipative Lagrangian and Hamiltonian mechanics [3, 4]. However, the method described in this section can be generalized directly to a general (possibly time-dependent) potential function V = V(x, t).

3-2 Historical perspectives

A traditional, engineering-inclined method to incorporate damping in the framework is to include a Rayleigh damping term in the Lagrangian to emulate linear damping forces, and this works 'mathematically' to derive the correct equations of motion [5]. Although frequently

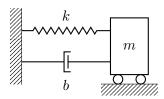


Figure 3-1: Schematic of the mass-spring-damper system.

used for practical problems, this damping term is not really part of the *actual* Lagrangian — rather, it simply makes use of the notion of a generalized force that is not inherently part of the system. As such, this method only 'works' on a superficial level: the pristine differential geometric foundations of mechanics do not leave room for such ad hoc tricks. There is, as a result, also no Hamiltonian counterpart for this method.

The historical attempts to do better than the Rayleigh method were primarily motivated by the application of the (dissipative) Hamiltonian formalism in quantum mechanics through discretization. For this application, a sound mathematical structure is of the essence, which calls for a more rigorous approach. A celebrated paper by Dekker [3] provides an excellent summary of many attempts up to 1981. Indeed, the well-studied approach developed by Caldirola [6] and Kanai [7] was intended exactly for this purpose. This method features an explicit time-dependence both in the Lagrangian function

$$L_{\text{CK}}(x, \dot{x}, t) = e^{\gamma t} \left(\frac{1}{2} m \dot{x}^2 - \frac{1}{2} k x^2 \right),$$
 (3-2)

and the corresponding Hamiltonian function:

$$H_{\rm CK}(x, \mathfrak{p}, t) = \frac{\mathfrak{p}^2}{2m} e^{-\gamma t} + \frac{1}{2} k x^2 e^{\gamma t}.$$
 (3-3)

In the Hamiltonian equation, \mathfrak{p} refers¹ to a special 'canonical momentum' (which is to be made clear later), defined by $\mathfrak{p} = p e^{\gamma t} = m \dot{x} e^{\gamma t}$. From either eq. (3-2) or eq. (3-3), the equations of motion are readily derived (for the Hamiltonian case with respect to \mathfrak{p} after which the transformation to p can be effected). Indeed, after taking the appropriate derivatives, one obtains:

$$\frac{\mathrm{d}}{\mathrm{d}t} \left(\frac{\partial L_{\mathrm{CK}}}{\partial \dot{x}} \right) - \frac{\partial L_{\mathrm{CK}}}{\partial x} = 0$$

$$\Rightarrow \mathrm{e}^{\gamma t} (m\ddot{x} + m\gamma \dot{x} + kx) = 0$$

for the Lagrangian case. Hamilton's equations amount to: [8]

$$\dot{x} = \frac{\partial H_{\text{CK}}}{\partial \mathfrak{p}} = \frac{\mathfrak{p}}{m} e^{-\gamma t} = \frac{p}{m},$$
$$\dot{\mathfrak{p}} = -\frac{\partial H_{\text{CK}}}{\partial x} = -kxe^{\gamma t}.$$

The relation between the time derivatives of the momenta \dot{p} and $\dot{\mathfrak{p}}$ is slightly more involved since one must invoke the product rule as a result of their time-dependencent relation:

$$\dot{\mathfrak{p}} = \mathrm{e}^{\gamma t} (\dot{p} + \gamma p).$$

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¹Not to be confused with the conventional notation for (Lie) algebras.

Substition yields the correct equation for p, though the equation is again multiplied by $e^{\gamma t}$. Because the latter is very well-behaved (that is, it has no zeros), it can be divided without any problems.

Geometric perspective The derivation stated above corresponds to the *suspension* of the time-dependent Hamiltonian vector field: ²

$$X_{H_{\text{CK}}} = -kx \mathrm{e}^{\gamma t} \frac{\partial}{\partial \mathfrak{p}} + \frac{\mathfrak{p}}{m} \mathrm{e}^{-\gamma t} \frac{\partial}{\partial x} + \frac{\partial}{\partial t}$$

dictated by the symplectic 2-form (in terms of the canonical momentum)

$$\omega = \mathrm{d}x \wedge \mathrm{d}\mathfrak{p}$$
,

such that $X_{H_{\text{CK}}}|_{t} \perp \omega = d(H_{C}K)$!!! not yet correct

Suspension includes the $\frac{\partial}{\partial t}$ coordinate in the vector field, see Foundations p. 373

Some peculiarities of the Caldirola-Kanai method deserve to be pointed out, since they will play a vital role in what is to come in the remainder of this chapter. [...]

3-2-1 Contact manifolds

!!! Our Lagrangian is CK-Lagrangian, but derivative wrt $\dot{q} =>$ obtain correct momentum for Hamiltonian formulation. In Hamiltonian formulation, choose p_0 and q_0 accordingly

'Contact' Hamiltonian (naive, from CK -> not complete!)

$$\hat{H} = \frac{p^2}{2m} + V(q)$$

Look at CK, call $p_0 = e^{\gamma t}$ and $p = p_1/p_0$ (for now explicit definition). Symplectify, i.e.

$$H = p_0 \left(\frac{1}{2m} \left(\frac{p_1}{p_0} \right)^2 + V(q) \right)$$

such that indeed,

$$H(p_0, p_1, q) = p_0 \hat{H}(p, q) = p_0 \hat{H}(p_1/p_0, q)$$

and we can use the contact relations given. Clearly, they are equivalent to the CK-method. HOWEVER:

observe that, $q_0 = \int L \, \mathrm{d}t$, and that $\frac{\partial \hat{H}}{\partial q_0} = \frac{\dot{p_0}}{p_0}$. Hence, the contact Hamiltonian can be modified to incorporate this, and presto, there is Bravetti's Hamiltonian. Hence, these methods are precisely equivalent through symplectization. TODO: Lagrangian counterpart, recover from our method what is three to save

²A time-dependent vector field on a manifold M is a mapping $X: N \times \mathbb{R} \to TM$ such that for each $t \in \mathbb{R}$, the restriction X_t of X to $N \times \{t\}$ is a vector field on M. [?] An additional construction of importance, called the suspension of the vector field, (...)

Split-Quaternions as Dynamical **Systems**

! orthogonal refers to 'regular' orthogonal, Lorentz-orthogonal makes the distinction.

Motivation: u seems to be 'aligned' with major direction of the elliptic trajectory in the Lorentz-orthogonal subspace, generated by the action of its cross-product. Show this formally by making use of the eigenvectors.

The basis vectors $\{e_2, e_3\}$, where e_2 is the orthogonal projection of the vector $e_1 = \hat{u}$ on its Lorentz-orthogonal subspace, and $e_3 \triangleq e_1 \times_L e_2$, form the real and imaginary parts of two of the eigenvectors of the matrix $U_{\times_{L}}$.

Because the basis vectors e_2 and e_3 are also orthogonal in the Euclidean sense, the

Proof. Let $\hat{u} = u_1 \hat{i} + u_2 \hat{j} + u_3 \hat{k}$. A normal vector to the Lorentz-orthogonal subspace is $\hat{\boldsymbol{n}} = u_1 \hat{\boldsymbol{i}} - u_2 \hat{\boldsymbol{j}} - u_3 \hat{\boldsymbol{k}}$. Then, the basis vectors are

$$e_{2} = \hat{\boldsymbol{u}} - \frac{\langle \hat{\boldsymbol{u}}, \hat{\boldsymbol{n}} \rangle}{\langle \hat{\boldsymbol{n}}, \hat{\boldsymbol{n}} \rangle} \hat{\boldsymbol{n}}$$

$$e_{3} = \hat{\boldsymbol{u}} \times_{L} e_{2} = -\frac{\langle \hat{\boldsymbol{u}}, \hat{\boldsymbol{n}} \rangle}{\langle \hat{\boldsymbol{n}}, \hat{\boldsymbol{n}} \rangle} (\hat{\boldsymbol{u}} \times_{L} \hat{\boldsymbol{n}}),$$
(4-1)

because the Lorentz-cross product distributes over addition and $\hat{\boldsymbol{u}} \times_{\mathbf{L}} \hat{\boldsymbol{u}} = \mathbf{o}$. The proposition above claims that $e_2 + ie_3$ is an eigenvector of the matrix U_{\times_L} . Hence, it must be the case that $U_{\times_{\mathbf{L}}}(\mathbf{e}_2 + \mathrm{i}\mathbf{e}_3) = \lambda(\mathbf{e}_2 + \mathrm{i}\mathbf{e}_3)$, where λ is then an eigenvalue of the matrix. This can be verified by replacing the action of U_{\times_L} with the cross product. Plugging in the definition and exploiting the linearity of the Lorentz cross-product, one obtains:

$$\begin{split} \hat{\boldsymbol{u}} \times_{L} \left(\boldsymbol{e}_{2} + \mathrm{i}\boldsymbol{e}_{3} \right) &= \hat{\boldsymbol{u}} \times_{L} \boldsymbol{e}_{2} + \mathrm{i} (\hat{\boldsymbol{u}} \times_{L} \boldsymbol{e}_{3}) \\ &= \boldsymbol{e}_{3} + (\hat{\boldsymbol{u}} \times_{L} \boldsymbol{e}_{3}) \mathrm{i} \\ &= \boldsymbol{e}_{3} + (\hat{\boldsymbol{u}} \times_{L} (\hat{\boldsymbol{u}} \times_{L} \boldsymbol{e}_{2})) \mathrm{i} \\ &= \boldsymbol{e}_{3} - \frac{\langle \hat{\boldsymbol{u}}, \hat{\boldsymbol{n}} \rangle}{\langle \hat{\boldsymbol{n}}, \hat{\boldsymbol{n}} \rangle} (\hat{\boldsymbol{u}} \times_{L} (\hat{\boldsymbol{u}} \times_{L} \hat{\boldsymbol{n}})) \mathrm{i}. \end{split}$$

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The triple cross-product expansion, or 'Lagrange formula', relates the regular cross product to the corresponding dot product:

$$a \times (b \times c) = b \langle c, a \rangle - c \langle a, b \rangle.$$

This well-known identity generalizes (easily verified) to the Lorentzian counterpart of the cross- and inner products:

$$\boldsymbol{a} \times_{\mathrm{L}} (\boldsymbol{b} \times_{\mathrm{L}} \boldsymbol{c}) = \boldsymbol{b} \langle \boldsymbol{c}, \, \boldsymbol{a} \rangle_{\mathrm{L}} - \boldsymbol{c} \langle \boldsymbol{a}, \, \boldsymbol{b} \rangle_{\mathrm{L}}.$$

Using the Lagrange formula, the above expression becomes

$$e_{3} - \frac{\langle \hat{\boldsymbol{u}}, \hat{\boldsymbol{n}} \rangle}{\langle \hat{\boldsymbol{n}}, \hat{\boldsymbol{n}} \rangle} (\hat{\boldsymbol{u}} \langle \hat{\boldsymbol{u}}, \hat{\boldsymbol{n}} \rangle_{L} - \hat{\boldsymbol{n}} \langle \hat{\boldsymbol{u}}, \hat{\boldsymbol{u}} \rangle_{L}) i$$

$$= e_{3} - \left(\hat{\boldsymbol{u}} \frac{\langle \hat{\boldsymbol{u}}, \hat{\boldsymbol{n}} \rangle_{L} \langle \hat{\boldsymbol{u}}, \hat{\boldsymbol{n}} \rangle}{\langle \hat{\boldsymbol{n}}, \hat{\boldsymbol{n}} \rangle} - \hat{\boldsymbol{n}} \frac{\langle \hat{\boldsymbol{u}}, \hat{\boldsymbol{n}} \rangle}{\langle \hat{\boldsymbol{n}}, \hat{\boldsymbol{n}} \rangle} \right) i$$

$$= e_{3} - \left(\hat{\boldsymbol{u}} - \hat{\boldsymbol{n}} \frac{\langle \hat{\boldsymbol{u}}, \hat{\boldsymbol{n}} \rangle}{\langle \hat{\boldsymbol{n}}, \hat{\boldsymbol{n}} \rangle} \right) i$$

$$= e_{3} - e_{2} i.$$

The latter is the scalar multiple of the vector e_2+e_3 by -i - hence, this is indeed an eigenvector of the corresponding matrix.

Because e_2 and e_3 are also orthogonal in the normal sense, they are aligned with the major axes of the elliptic trajectories generated by the cross product. Hence, they can be used to find a basis of the invariant subspace which makes the trajectories identical to those in the phase plane.

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Chapter 5

Conclusion

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14 BIBLIOGRAPHY

Glossary

List of Acronyms

Mathematical notation

 \boldsymbol{v} A (tangent) vector

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