

# The Stability of Complex Time Steppers

CIAN J. DUGGAN



Trinity  
College  
Dublin

The University of Dublin

## INTRODUCTION

This poster looks at the stability of numerical methods when applied to problems of differential calculus.

This project aims to add a visual understanding to the concepts of stability. It also aims to explore the use of complex time steps in numerical methods.

## KEY TERMINOLOGY

**Numerical Methods** are iterative algorithms used to approximate solutions to problems that cannot be solved exactly.

These methods are often employed to solve differential equations.

The infinitesimal  $h$  in  $y'(t) = \lim_{h \rightarrow 0} \frac{y(t+h) - y(t)}{h}$  cannot be represented on a computer.

Instead, a finite **time step**  $h$  is used to approximate the derivative.

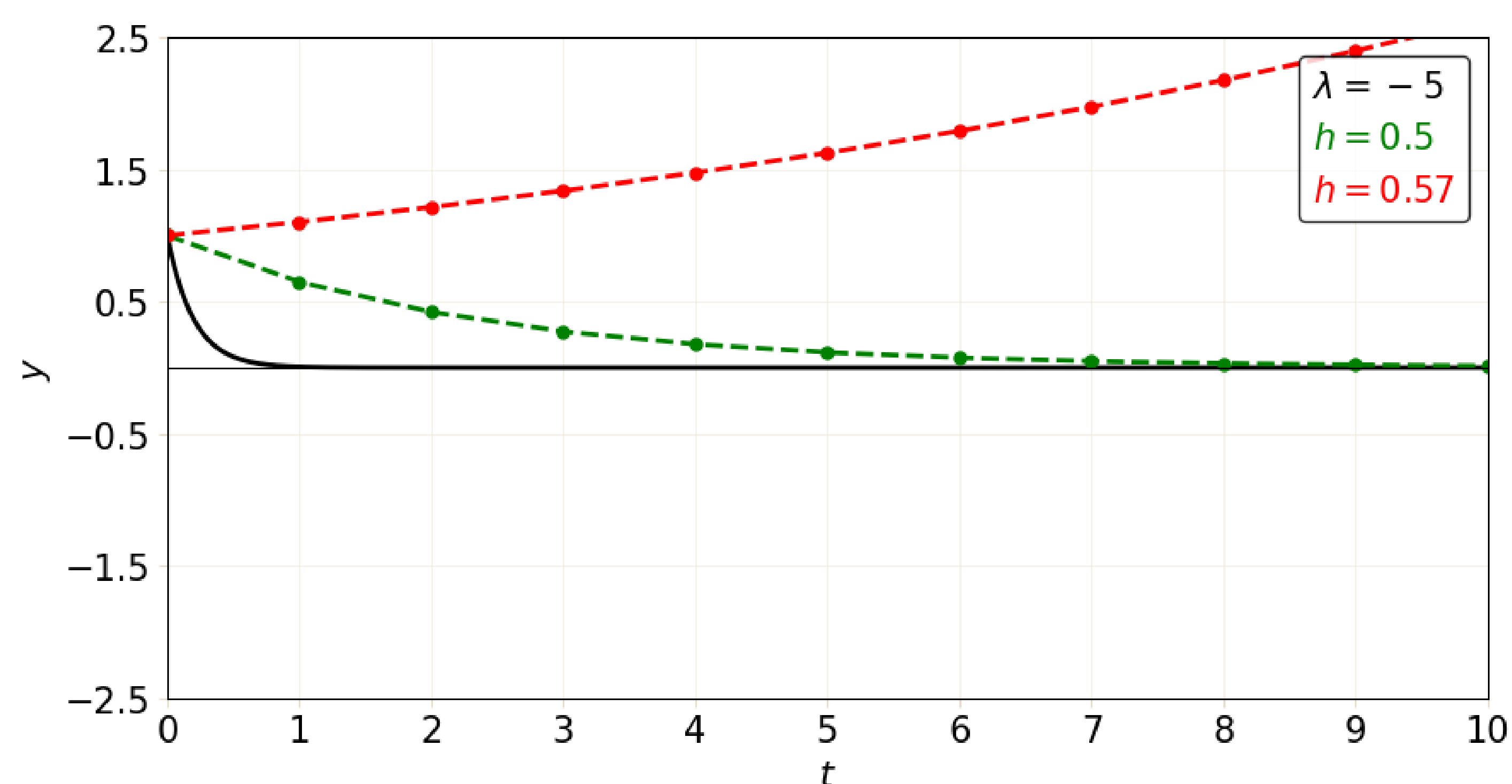
A numerical method is **stable** for a given size of time step if the solution does not diverge from the exact solution.

The choice of step size is crucial; too large, and the solution will diverge, too small, and the solution will cost computational resources.

Shown below is **Runge-Kutta 4** applied to the Exponential Decay Problem.

The exact solution is known, and shown in black.

Stable and unstable choices of  $h$  are shown in green and red respectively.



## THE EXPONENTIAL DECAY PROBLEM

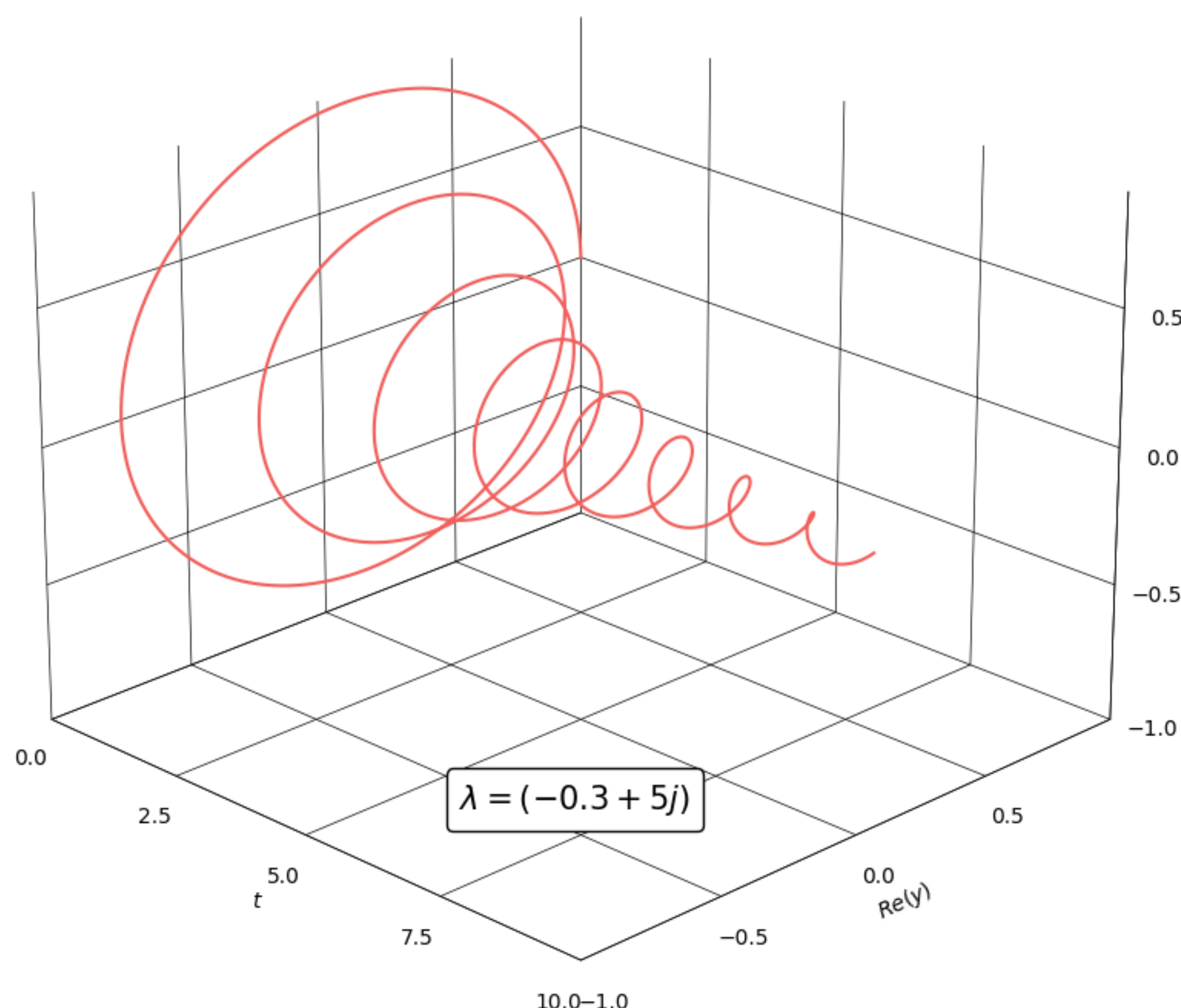
Consider a quantity  $y$  that decays relative to its current value.

The quantity at time  $t$  is  $y(t)$ .

The rate of decay is  $\lambda \in \mathbb{C}$  with  $\text{Re}(\lambda) < 0$ .

The simplest case of this problem is described by the differential equation:  $y'(t) = \lambda y(t)$  with  $y(0) = 1$

This has the exact solution  $y(t) = e^{\lambda t}$  giving the name *Exponential Decay*.



## STABILITY

We can define the **Stability Function**  $s(\lambda, h)$  for any numerical method by writing the algorithm in the form  $y_{j+1} = s(\lambda, h)y_j$ .

A method is stable for any pair in  $S = \{(\lambda, h) : |s(\lambda, h)| < 1\} \in \mathbb{C}$ , the **Stability Region**.

Applying **Runge-Kutta 4**:  $y_{j+1} = (1 + \lambda h + \frac{(\lambda h)^2}{2} + \frac{(\lambda h)^3}{6} + \frac{(\lambda h)^4}{24})y_j$

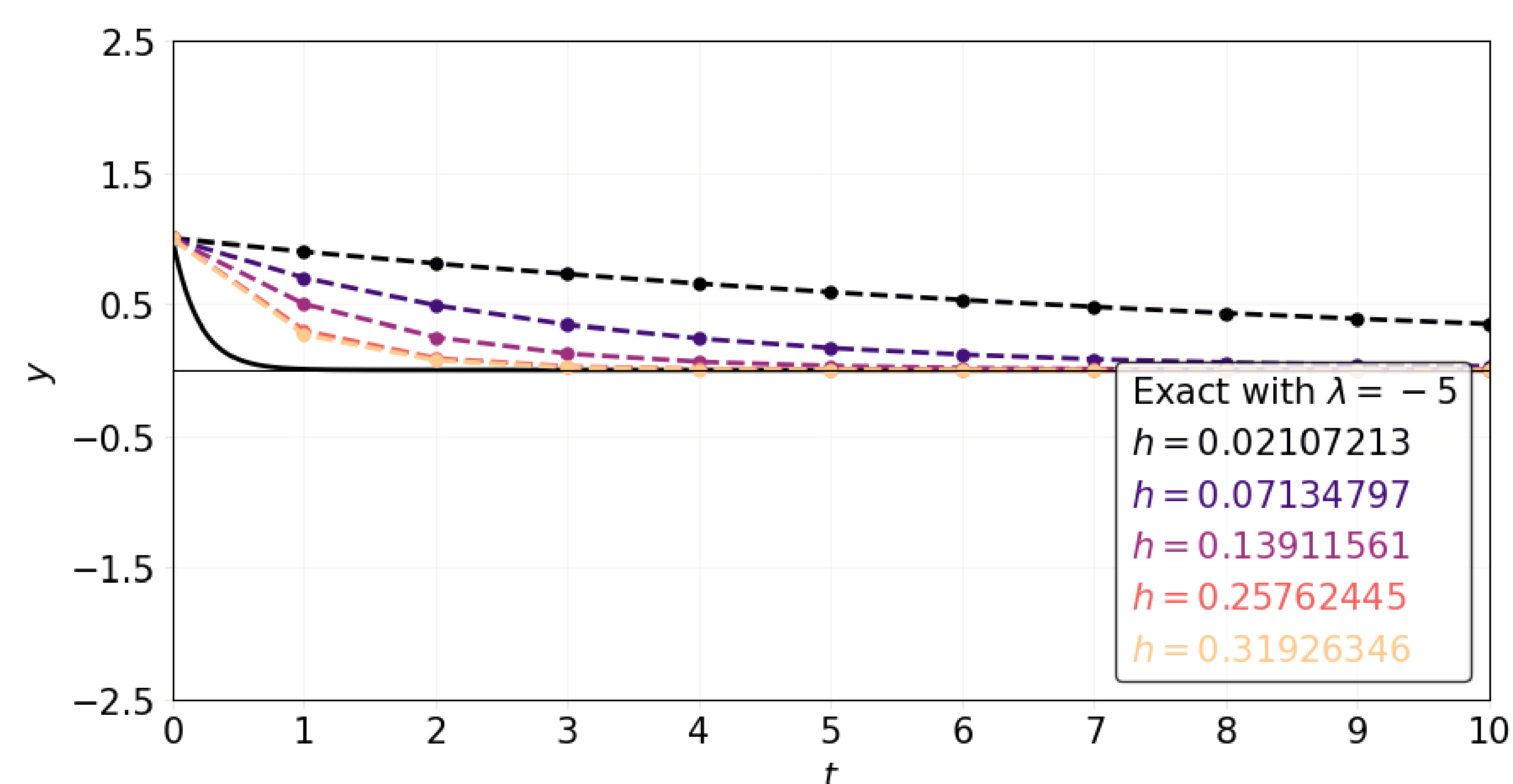
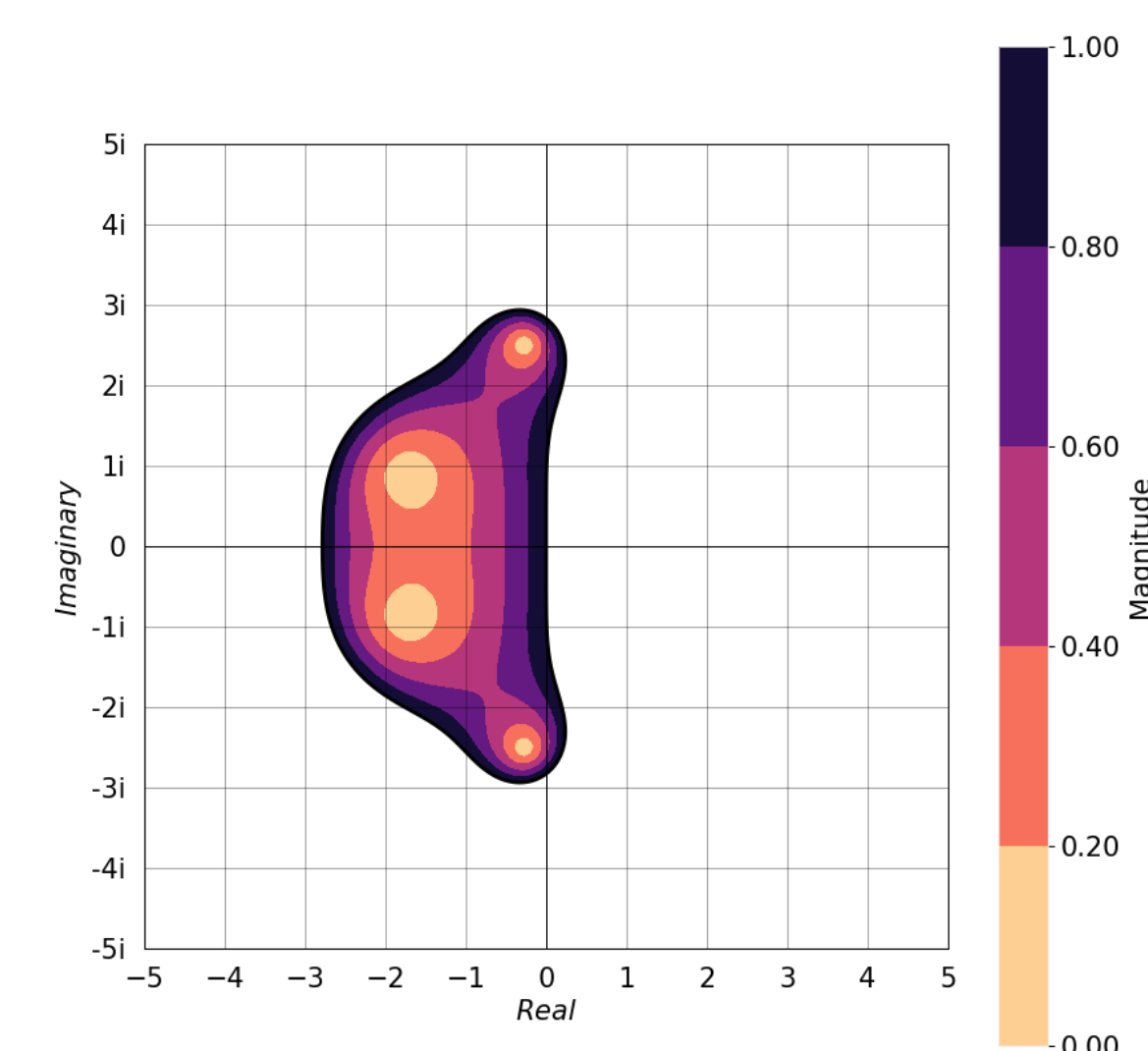
The stability function is

$$s(\lambda, h) = \sum_{n=0}^4 \frac{(\lambda h)^n}{n!}$$

$S$  is shown on the right.

$S$  is divided into subregions, each corresponding to a different range of  $|s(\lambda, h)|$  values.

The methods with these step sizes are shown below, for  $\lambda \in \mathbb{R}^-$ :



## THE COMPLEX DOMAIN

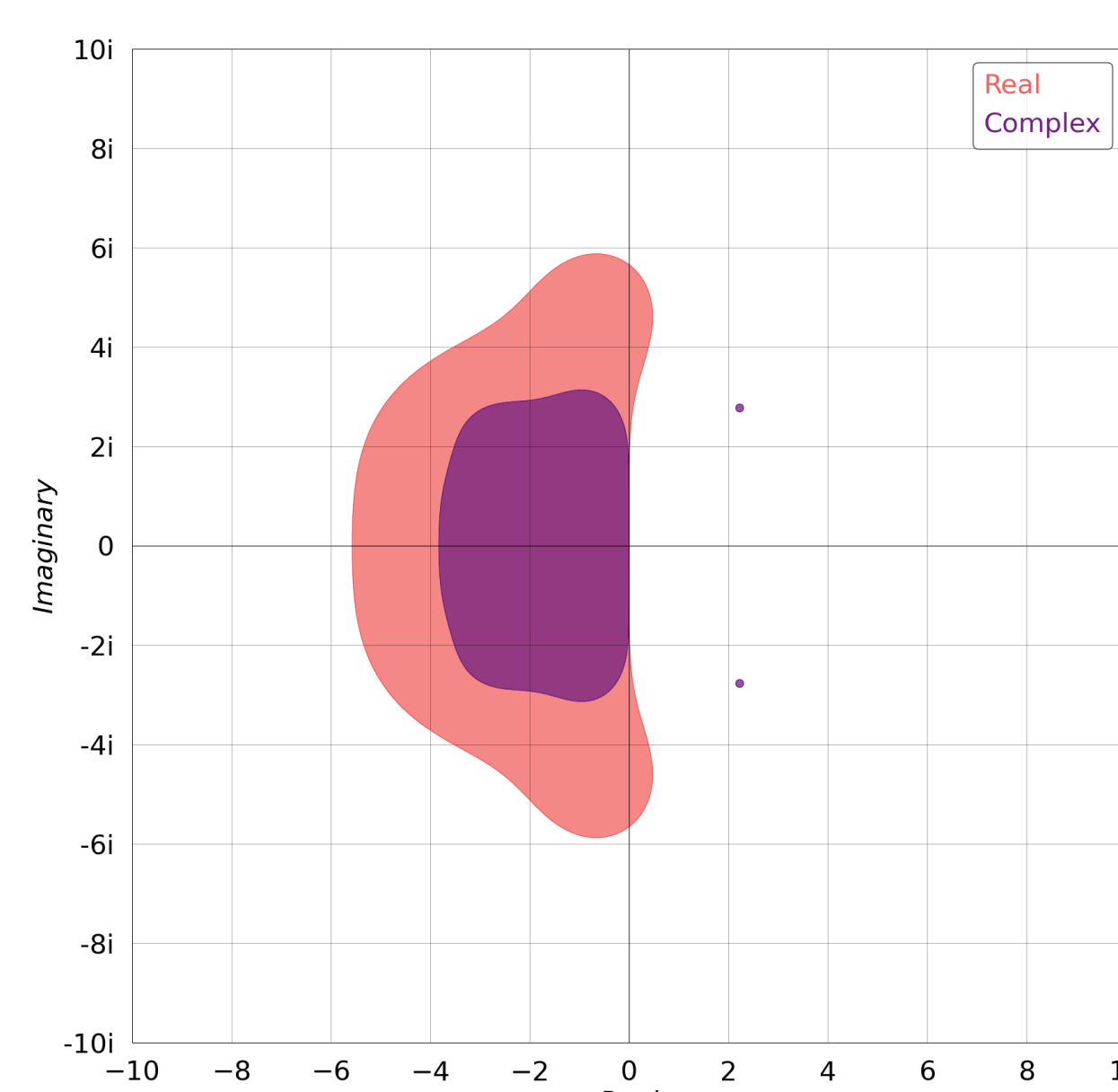
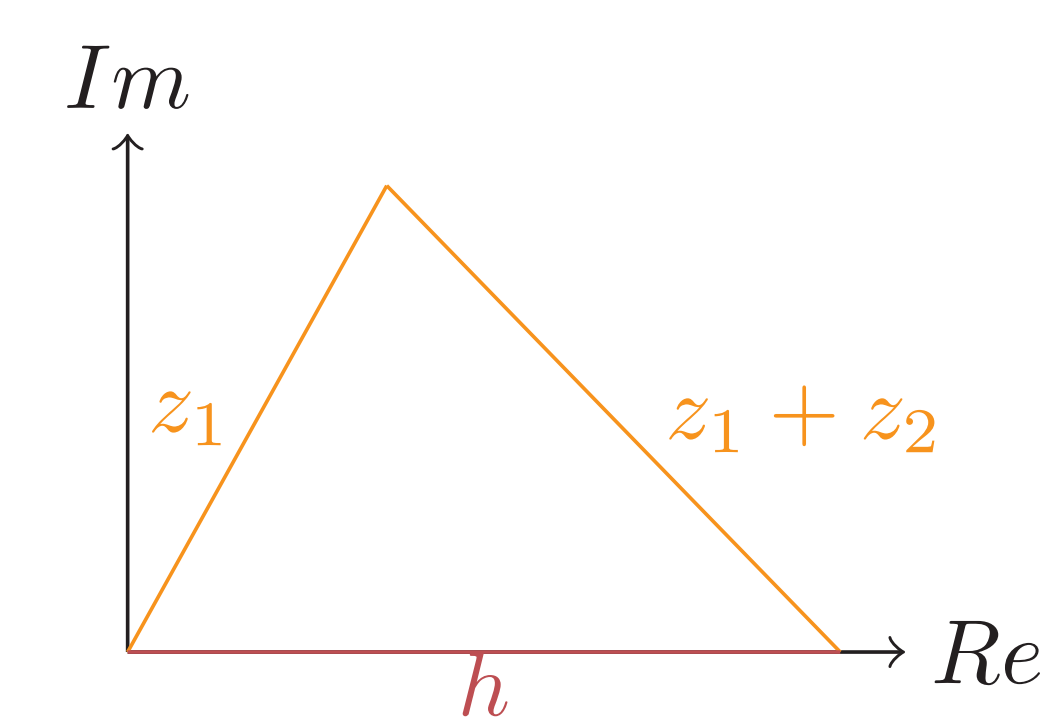
It has often proven useful to extend the analysis of a problem to the complex domain. We will have a look at the stability of numerical methods when the chosen time-step is complex.

We want to take an overall step of size  $h \in \mathbb{R}$  comprised of two steps  $z_1, z_2 \in \mathbb{C}$  such that  $z_1 + z_2 = h$ .

We compare to the case where we take 2 steps of size  $\frac{h}{2} \in \mathbb{R}$ .

The following diagram shows the case

$z_1 = \frac{h}{2}(1 + i)$  and  $z_2 = \bar{z}_1$



George, Yung and Mangan[1] show that the error is reduced when using complex time steps.

For Runge-Kutta 4, we obtain two different stability functions

$s_{\mathbb{R}}(\lambda, h)$  and  $s_{\mathbb{C}}(\lambda, h)$

They give the stability regions shown on the left.

While the error may decrease, the stability region decreases with it.

## REFERENCES

- [1] Jithin D. George, Samuel Y. Jung, and Niall M. Mangan, *Walking into the complex plane to 'order' better time integrators*, 2021. <https://arxiv.org/abs/2110.04402>
- [2] The GitHub repository for this project, <https://github.com/CianJDuggan/Capstone-Project>