# oscode: fast solutions of oscillatory ODEs in cosmology

Fruzsina J. Agocs \* 1,2

<sup>1</sup>Kavli Institute for Cosmology, Cambridge <sup>2</sup>Astrophysics Group, Cavendish Laboratory, Cambridge

oscode is a Python/C++ package for the fast solution of oscillatory ordinary differential equations. It handles equations of the form  $\ddot{x}(t) + 2\gamma(t)\dot{x}(t) + \omega^2(t)x(t) = 0$ . The time-dependence of the frequency  $\omega$  and the damping term  $\gamma$  can be explicit or implicit; below are examples of both. The algorithm is significantly more efficient than conventional (Runge-Kutta-based) solvers found in numerical libraries, thanks to reduced number of steps needed to traverse highly oscillatory regions.

### **Airy equation**

**Burst equation** 



**Figure 1:** Steps and accuracy of **oscode** solving  $\ddot{x} + tx = 0$ .

• Left: different coloured dots show the internal steps oscode takes in `RK' and `WKB' mode, the continuous line being the exact solution

**Schrödinger equation** 

- **Right:** relative accuracy of **oscode**'s approximate solution
- Compare to scenario where oscode only takes `RK' steps

**Figure 2:** Solving  $\ddot{x} + \frac{n^2 - 1}{(1 + t^2)^2}x = 0$ , an equation exhibiting a burst of oscillations.

- Equation has a parameter  $n, n \sim no.$  of oscillations in total
- Left: oscode's internal steps for n = 40
- Middle: Relative error in oscode's solution, and error in `pure RK' solution for comparison
- **Right:** number of oscillations crossed per step for  $n = 10^5$

### **Primordial power spectra**



**Figure 3:** Eigenstates of the Schrödinger equation  $(\Psi'' + 2m(E - V(x))\Psi = 0)$  in a harmonic potential well.

- Above is an example where the eigenenergies are known, but can compute them numerically:
- Start integration from well outside potential well on either side and meet at x = 0.5
- Compute  $\Psi'/\Psi$  at the meeting point for both solutions; the difference is minimal if the energy is an eigenvalue
- Calculated energy eigenvalues in a harmonic well with quartic anharmonicity up to the  $10000^{\text{th}}$  energy level to 1 in  $10^7$  accuracy

## Algorithm

- **RK (Runge-Kutta) step:** approximates the solution at a later time t + h as a Taylor expansion around t.
- WKB (Wentzel-Kramers-Brillouin) step: uses an analytic approximation which gets better the slower the frequency of oscillations change, allowing **oscode** to traverse highly oscillatory regions in few steps.
- **oscode** switches between RK/WKB mode, choosing whichever allows for the largest step within a given error tolerance.
- The stepsize is updated to ensure the local error stays within the given tolerance. Derivatives of the frequency and damping term are required for both the RK and WKB steps, these are computed numerically as finite differences.
- WKB steps also require integrals, which are computed using Gauss-Lobatto integration.
- The RK steps and finite difference formulae were designed to minimise function evaluations.

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**Figure 4:** Scalar primordial power spectra in closed universes with varying start-of-inflation curvature  $\Omega_k^i$ .

- Solve the Mukhanov–Sasaki (MS) equation for each wavenumber k, frequency is roughly proportional to k
- MS equation is an example of an oscillator where  $\omega$  and  $\gamma$  are not explicit functions of time, are only available numerically
- oscode deals with this by performing interpolation on a grid of  $\omega$  and  $\gamma$  values
- Can test models where the perturbations have to be computed while they are deep inside the Hubble horizon and go through many oscillations before freeze-out.

For full paper, code and an animated version of Fig. 4, take a photo of this QR code or visit



https://github.com/fruzsinaagocs/oscode

\* e-mail: fa325@cam.ac.uk