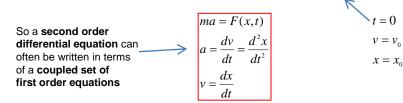
Numeric solution of ordinary differential equations

Many laws of the Physical Sciences are expressed in terms of derivatives. e.g. Newton's Second Law of (non-relativistic) dynamics:

mass x acceleration = vector sum of forces

In one dimension, acceleration is the rate of change of velocity, and velocity is the rate of change of displacement. With appropriate symbols we could write a set of differential equations to describe subsequent motion, given a knowledge of the initial conditions.



Some differential equations can be solved analytically, i.e. in terms of an expression comprising of basic Mathematical functions. However, in general this is not possible. To make further progress we need a numerical method that we can apply using a computational tool such as Excel or MATLAB to evaluate in an *iterative* fashion. Most methods will assume a fixed, small time step (or x step) and 'solve' the equation approximately.

Euler's Method is probably the simplest, but least precise method.

First order equations - i.e. in terms of a single derivative

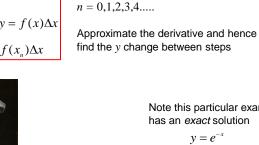
$$\frac{dy}{dx} = f(x); \quad y = y_0 \text{ when } x = x$$

$$x_{n+1} = x_n + \Delta x$$

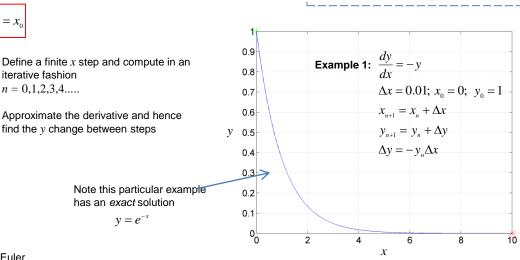
$$y_{n+1} = y_n + \Delta y$$

$$\frac{dy}{dx} \approx \frac{\Delta y}{\Delta x} \therefore \Delta y = f(x)\Delta x$$

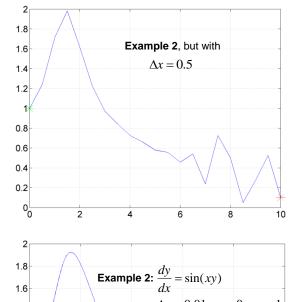
$$\therefore y_{n+1} = y_n + f(x_n)\Delta x$$

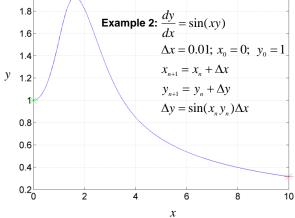


iterative fashion



size of Δx





Leonhard Euler 1707-1783

Mathematics topic handout: Calculus – Numeric solvers of differential equations Dr Andrew French. www.eclecticon.info PAGE 1

Making the

solver

For the Euler method, it can be shown that

errors propagate in direct proportion to the

finite difference

(i.e. the x step)

smaller introduces

less error into the

Euler's Method can be extended for equations involving higher derivatives, by creating a system of coupled first order equations

$$\frac{d^2 y}{dx^2} = f\left(\frac{dy}{dx}, y, x\right); \quad y = y_0 \& \frac{dy}{dx} = m_0 \text{ when } x = x_0$$

$$m = \frac{dy}{dx} \qquad \text{i.e. a coupled set of} \\ \text{first order equations}$$

$$\therefore \frac{dm}{dx} = f(m, y, x)$$

$$x_{n+1} = x_n + \Delta x$$

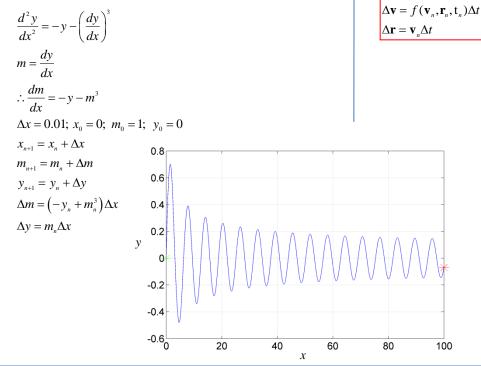
$$m_{n+1} = m_n + \Delta m$$

$$y_{n+1} = y_n + \Delta y$$

$$\Delta m = f(m_n, y_n, x_n) \Delta x$$

$$\Delta y = m_n \Delta x$$

Example using Euler's method:



We can readily extend the method to two, three or higher dimensional problems which are parameterized in terms of a single variable by using vectors.

x v **r** =

 $\mathbf{v} = \frac{d\mathbf{r}}{dt} = \begin{pmatrix} \frac{dx}{dt} \\ \frac{dy}{dt} \end{pmatrix}$

 $\therefore \frac{d\mathbf{v}}{dt} = f(\mathbf{v}, \mathbf{r}, \mathbf{t})$

 $t_{n+1} = t_n + \Delta t$ $\mathbf{v}_{n+1} = \mathbf{v}_n + \Delta \mathbf{v}$

 $\mathbf{r}_{n+1} = \mathbf{r}_n + \Delta \mathbf{r}$

$$\mathbf{r} = \begin{pmatrix} x \\ y \end{pmatrix}$$

$$\frac{d^{2}\mathbf{r}}{dt^{2}} = f\left(\frac{d\mathbf{r}}{dt}, \mathbf{r}, \mathbf{t}\right); \quad \mathbf{r} = \begin{pmatrix} x_{0} \\ y_{0} \end{pmatrix} \& \frac{d\mathbf{r}}{dt} = \begin{pmatrix} u_{s} \\ u_{y} \end{pmatrix} \text{ when } t = 0$$

$$\mathbf{v} = \frac{d\mathbf{r}}{dt} = \begin{pmatrix} \frac{dx}{dt} \\ \frac{dy}{dt} \end{pmatrix}$$

$$\therefore \frac{d\mathbf{v}}{dt} = f(\mathbf{v}, \mathbf{r}, \mathbf{t})$$

$$t_{s+1} = r_{s} + \Delta \mathbf{t}$$

$$\mathbf{v}_{s+1} = \mathbf{v}_{s} + \Delta \mathbf{v}$$

$$\mathbf{r}_{s+1} = \mathbf{r}_{s} + \Delta \mathbf{r}$$

$$\Delta \mathbf{v} = f(\mathbf{v}_{s}, \mathbf{r}_{s}, \mathbf{t}_{s}) \Delta t$$

$$\Delta \mathbf{r} = \mathbf{v}_{s} \Delta t$$

$$\mathbf{r} = \mathbf{v}_{s} \Delta t$$

2

3

х

4

5

0

1

There are various methods which can offer greater precision given a fixed step size. The **Runge-Kutta** method below will extend to higher derivatives and vector equations in exactly the same way as described for the Euler method. The **Verlet** method is really specific to kinematic problems, but will naturally extend to two or three dimensions using vectors.

Verlet Method – assume 'constant acceleration motion' between time steps

$$v = \frac{dx}{dt}$$

$$\frac{dv}{dt} = a(x, v, t); \quad x = x_0 \& v = u \quad \text{when } t = 0$$

$$t_{n+1} = t_n + \Delta t$$

$$v_{n+1} = v_n + \Delta v$$

$$x_{n+1} = x_n + \Delta x$$

$$\Delta x = v_n \Delta t + \frac{1}{2} a(x_n, v_n, t_n) (\Delta t)^2$$

$$V = v_n + a(x_n, v_n, t_n) \Delta t$$

$$\Delta v = \frac{1}{2} \{a(x_n, t_n) + a(x_{n+1}, V, t_{n+1})\} \Delta t$$

This is a **second order** method i.e. errors are proportional to $(\Delta x)^2$

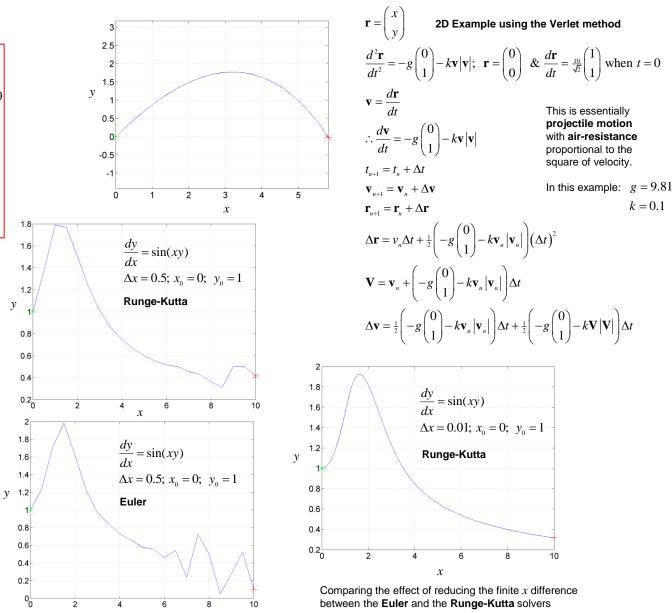
Note if acceleration is functionally dependent on velocity we have to use a first order Euler method to evaluate the second acceleration term.

Runge-Kutta method

$$\frac{dy}{dx} = f(y, x); \quad y = y_0 \text{ when } x = x_0$$

$$\begin{aligned} x_{n+1} &= x_n + \Delta x \\ y_{n+1} &= y_n + \Delta y \\ k_1 &= f(y_n, x_n) \\ k_2 &= f\left(y_n + \frac{1}{2}\Delta x k_1, x_n + \frac{1}{2}\Delta x\right) \\ k_3 &= f\left(y_n + \frac{1}{2}\Delta x k_2, x_n + \frac{1}{2}\Delta x\right) \\ k_4 &= f\left(y_n + \Delta x k_3, x_n + \Delta x\right) \\ \Delta y &= \frac{1}{6}\Delta x \left(k_1 + 2k_2 + 2k_3 + k_4\right) \end{aligned}$$

This is a **fourth order** method i.e. errors are proportional to $(\Delta x)^4$



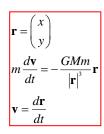


Carl Runge 1856-1927



Martin Kutta 1867-1944

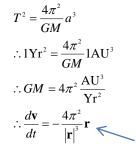
Solar system simulation using the Verlet method



Newton's law of Universal Gravitation. M is the mass of the Sun, which is assumed to be so massive compared to the other planets that the interaction of their gravity on it is negligible. i.e. it is fixed in position.

The gravitational interaction *between* the planets is also ignored.

Using Kepler's Third law



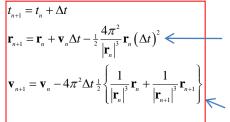
T is the orbital period a is the semi-major axis of the orbit (in general, bound orbits are *ellipses*)

Yr is an Earth year AU is an astronomical unit i.e. the average Earth-Sun separation

If lengths are measured in AU and times in Years. Note planet masses cancel and therefore are not required.

Can't readily **Runge-Kutta** since the left had side of the differential equation is functionally dependent on *neither* velocity or time

Solver using the Verlet method



i.e. constant acceleration motion between time steps

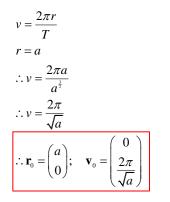
> Use the average acceleration between the time steps to work out the new velocity

Initial conditions for circular orbits.

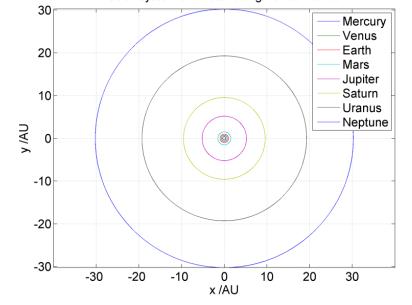
In (Yr, AU) units, Kepler III is:



Circular Motion:



 $M_{\odot} = 1.99 \times 10^{30} \text{ kg}$ $G = 6.67 \times 10^{-11} \text{ Nm}^2 \text{ kg}^{-2}$ $AU = 1.49597871 \times 10^{11} \text{ m}$ $24 \times 3600 \text{ s} = 1 \text{ day}$ $M_{\odot} = 332,837 m_{\oplus}$ $m_{\oplus} = 5.972 \times 10^{24} \text{ kg}$



	Planet	<i>T I</i> years	<i>r </i> AU	<i>m I</i> Earth masses	Rotation period /days	Orbital eccentricity
1	Mercury	0.241	0.387	0.055	58.646	0.21
	Venus	0.615	0.723	0.815	243.018	0.01
	Earth	1.000	1.000	1.000	1.000	0.02
	Mars	1.881	1.523	0.107	1.026	0.09
	Jupiter	11.861	5.202	317.85	0.413	0.05
	Saturn	29.628	9.576	95.159	0.444	0.06
	Uranus	84.747	19.293	14.5	0.718	0.05
	Neptune	166.344	30.246	17.204	0.671	0.01
	Pluto	248.348	39.509	0.003	6.387	0.25

Solar system simulator using Verlet method