

# Phys 512 ODE's

# ODEs

- usual case: we have  $y(x_0)$ , find  $y(x_0+h)$  given  $dy/dx=f(x,y)$
- Harder than integration, because we don't know how to evaluate  $dy/dx$  along the path - that depends on the (unknown) value of  $y$ .
- We can still apply many of the Taylor series tricks we've learned.
- NB - we could have a system of equations just as well as a single equation.
- NB 2 - we can also have higher order equations recast into a system. variables are then  $y, y', y'' \dots$

# 1st order

- We could just take  $y(x+h)=y(x)+h \, dy/dx = y(x)+hf$
- How will error scale with size of  $h$ ?
- This isn't very good...
- Can we do better?

# RK2

- I could take a trial step, then evaluate derivative. Maybe combine with first derivative to do better?
- Taylor expand  $y(x+h)$  to 2nd order.
- Take  $k_1=hf(x,y)$ ,  $k_2=hf(x+\alpha h, y+\beta k_1)=hf(x+\alpha h, y+\beta fh)$
- Let answer be  $y(x+h)=ak_1+bk_2$ .
- Solve for  $a, b, \alpha, \beta$  to make agree with Taylor.
- Answer underconstrained - have freedom to pick. Usual is  $\alpha=\beta=1$ ,  $a=b=1/2$ , or  $\alpha=\beta=1/2$  and  $a=0, b=1$ .

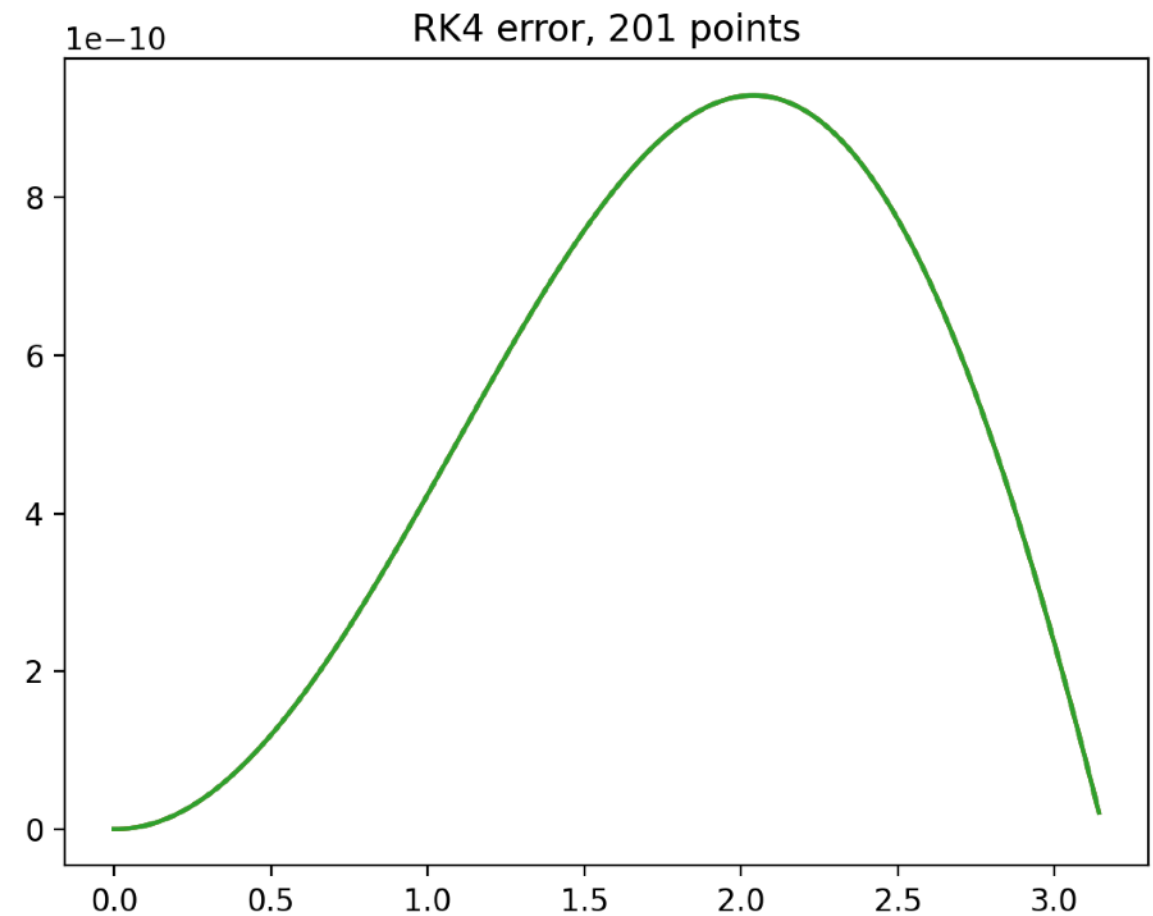
# RK4

- Can extend to 4<sup>th</sup> order. Mathy, but not very illustrative.
- $k_1 = hf(x, y)$   
 $k_2 = hf(x + h/2, y + k_1/2)$   
 $k_3 = hf(x + h/2, y + k_2/2)$   
 $k_4 = hf(x + h, y + k_3)$   
 $y(x + h) = y(x) + (k_1 + 2k_2 + 2k_3 + k_4)/6$
- Accurate to 4th order - ODE equivalent of Simpson's rule.
- Adaptive stepsize *highly* recommended.

```
def f(x,y):
    dydx=np.asarray([y[1],-y[0]])
    return dydx
```

```
def rk4(fun,x,y,h):
    k1=fun(x,y)*h
    k2=h*fun(x+h/2,y+k1/2)
    k3=h*fun(x+h/2,y+k2/2)
    k4=h*fun(x+h,y+k3)
    dy=(k1+2*k2+2*k3+k4)/6
    return y+dy
```

```
npt=201
x=np.linspace(0,np.pi,npt)
y=np.zeros([2,npt])
y[0,0]=1 #starting conditions
y[1,0]=0 #if I start at peak, then first derivative =0
for i in range(npt-1):
    h=x[i+1]-x[i]
    y[:,i+1]=rk4(f,x[i],y[:,i],h)
truth=np.cos(x)
print(np.std(truth-y[0,:]))
```



# Bulirsch-Stoer

- We saw we could take multiple not-very accurate integrals, and combine to get high accuracy by estimating error terms.
- Can do the same with ODEs - Romberg equivalent is called Bulirsch-Stoer.
- If you need high accuracy, have smooth function, try this!

# Stability

- let  $y' = -cy$ . Answer should be  $\exp(-cx)$
- Solve the stupid way:  $y(x+h) = y(x) + hf(x,y) = y(x) - hcy(x)$
- $y(x+h) = y(x)(1-hc)$ .  $y(x+nh) = y(x)(1-hc)^n$ .
- What happens if  $h$  is too large? For  $|1-hc| > 1$ , this *grows* exponentially. Large steps have made our answer *unstable*.



# Stiff Equations

- Very often in systems, one equation (or eigenmode) has a large  $c$ , other has a small  $c$ .
- Solution is large  $c$  converges very quickly, and stays at solution while we wait for small  $c$  to evolve.
- If we aren't careful and take steps for small  $c$ , then large  $c$  becomes unstable, and solution blows up.
- Shrinking time step enough to track large  $c$  may be impracticable.
- Such systems are called *stiff*. Solutions presented here are simplistic, but knowing you have a stiff set is most of the battle.

# U238 Decay

- Radioactive decay common situation.
- U238 takes 4 billion years to go to Th234. Po214 takes 160 microseconds to go to Pb210.
- To solve naively, takes (4 billion years/160 microseconds) =  $10^{21}$  timesteps.

	Half-Life	Time unit	Emitter
Uranium-238	4,468	billion of years	alpha
Thorium-234	24,10	days	beta -
Protactinium-234	6,70	hours	beta -
Uranium-234	245 500	years	alpha
Thorium-230	75380	years	alpha
Radium-226	1 600	years	alpha
Radon-222	3,8235	days	alpha
Polonium-218	3,10	minutes	alpha
Plomb-214	26,8	minutes	beta -
Bismuth-214	19,9	minutes	beta -
Polonium-214	164,3	microseconds	alpha
Plomb-210	22,3	years	beta
Bismuth-210	5,015	years	beta
Polonium-210	138,376	days	alpha
Plomb-206	Stable		

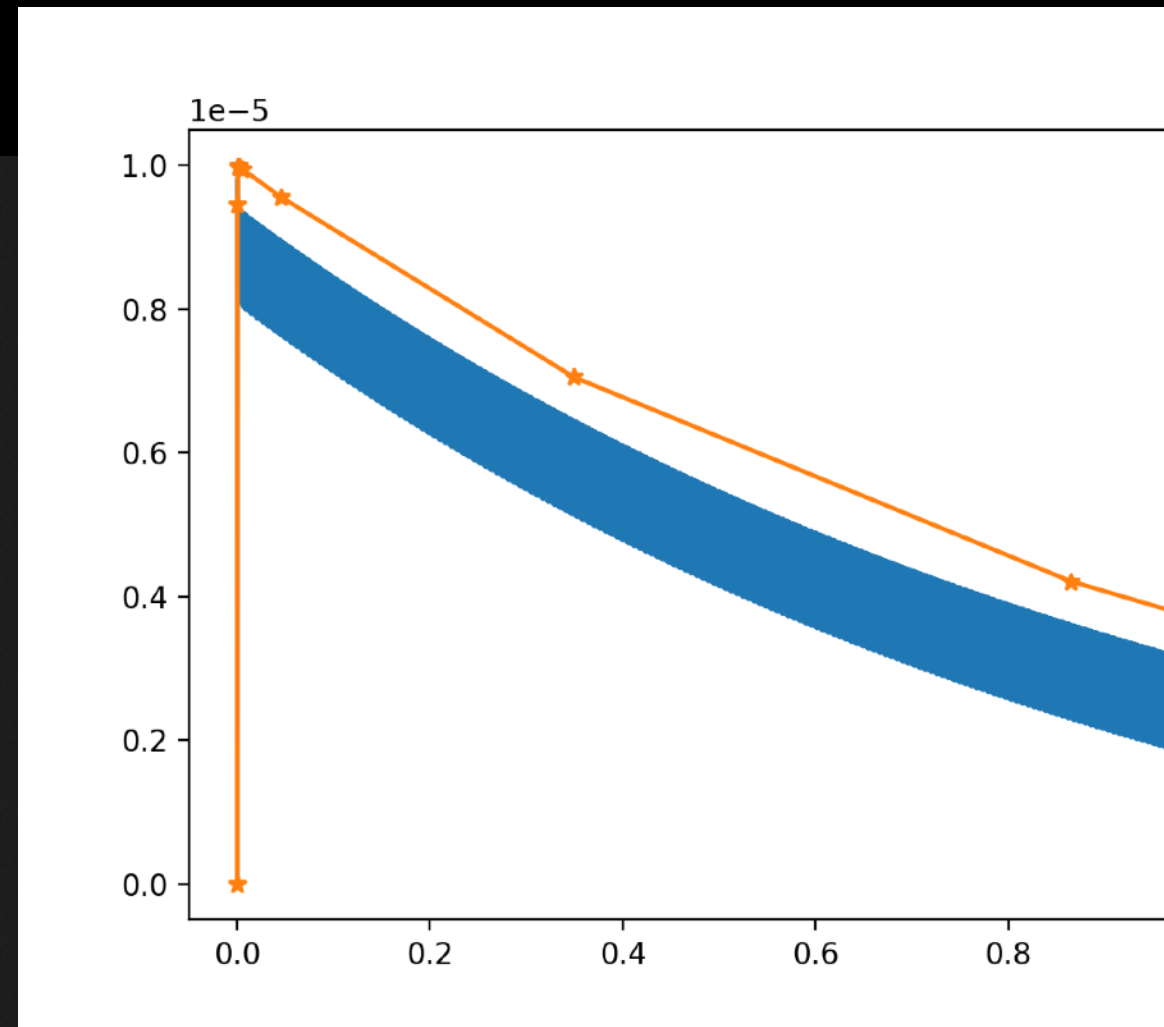
# Implicit

- Common solution - use derivative at end of interval.
- In constant coefficient case  $y_{n+1}=y_n-hcy_{n+1}$ .
- Solve:  $y_{n+1}(1+hc)=y_n$ ,  $y_{n+1}=y_n/(1+hc)$ . (reminder - old way was  $y_{n+1}=y_n(1-hc)$ . Agrees to 1<sup>st</sup> order but not 2<sup>nd</sup>. )
- I can now make  $h$  very large and remain stable. I may not be accurate, but I can crank up  $h$ .
- This is just an introduction, much better ways to handle stiff equations exist, but at least you know to look for them...

# Scipy ODE Driver

```
import numpy as np
from scipy import integrate
def fun(x,y, half_life=[1,1e-5]):
    #let's do a 2-state radioactive decay
    dydx=np.zeros(len(half_life)+1)
    dydx[0]=-y[0]/half_life[0]
    dydx[1]=y[0]/half_life[0]-y[1]/half_life[1]
    dydx[2]=y[1]/half_life[1]
    return dydx

y0=np.asarray([1,0,0])
x0=0
x1=1
ans_rk4=integrate.solve_ivp(fun,[x0,x1],y0)
ans_stiff=integrate.solve_ivp(fun,[x0,x1],y0,method='Radau')
print('took ',ans_rk4.nfev,' evaluations to solve with RK4.')
print('took ',ans_stiff.nfev,' evaluations to solve implicitly')
print('final values were ',ans_rk4.y[0,-1],' and ',ans_stiff.y[0,-1],'
```



```
[>>> exec(open('stiff.py').read())
took 212618 evaluations and 3.127746820449829 seconds to solve with RK4.
took 72 evaluations and 0.0035657882690429688 seconds to solve implicitly
final values were 0.3678794411714445 and 0.3678803705878816 with truth 0.367879441171442
[>>> plt.clf();plt.plot(ans_rk4.t,ans_rk4.y[1,:]);plt.plot(ans_stiff.t,ans_stiff.y[1,:], '*-')
```