

# 6 Body Solar System\_IRKNGL LUZEA

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## 1 6-Body Solar System IRKNGL

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### 1.1 Loading packages and functions

```
[1]: using LinearAlgebra
      using Plots
      using OrdinaryDiffEq
      using JLD2, FileIO
      using Parameters, NBInclude
      #using IRKGaussLegendre
      using LaTeXStrings
      using BenchmarkTools
      #using RecursiveArrayTools
```

```
[2]: PATH_DATA="./Data/"

      PATH_SRC="../../src_simd/"
      include(string(PATH_SRC,"IRKGL_SIMD.jl"))
      using .IRKGL_SIMD

      PATH_SRC="../../src_seq/"
      include(string(PATH_SRC,"IRKGL_SEQ.jl"))
      using .IRKGL_SEQ

      PATH_SRC="../../src/"
      include(string(PATH_SRC,"MyBenchmarksTools.jl"))
```

```
[2]: launch_method_tests (generic function with 1 method)
```

```
[3]: PATH_SRC2="../../src_seq2/"
include(string(PATH_SRC2, "IRKGL_SEQ.jl"))
#include(string(PATH_SRC2, "IRKGL_Seq_Solver.jl"))
#include(string(PATH_SRC2, "IRKGL_Step_Functions.jl"))

using .IRKGL_SEQ2
```

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## 1.2 Definition of the N-body problem

In Nbody.jl below, the following functions are defined: NbodyEnergy(u,Gm), NbodyODE!(du,u,Gm,t), and NbodyODE1!(du,u,Gm,t), where

$$u = \begin{pmatrix} q_1 & v_1 \\ \vdots & \vdots \\ q_N & v_N \end{pmatrix} \in \mathbb{R}^{2 \times 3 \times N}, \quad Gm = (Gm_1, \dots, Gm_N) \in \mathbb{R}^N.$$

The energy, as a function of the positions  $q_1, \dots, q_N \in \mathbb{R}^3$  and the velocities  $v_1, \dots, v_N \in \mathbb{R}^3$  of the  $N$  bodies is:

$$\text{Energy}(q_1, \dots, q_N, v_1, \dots, v_N) = \frac{1}{2} \sum_{i=1}^N m_i \|v_i\|^2 - G \sum_{1 \leq i < j \leq N} \frac{m_i m_j}{\|q_i - q_j\|}.$$

The ODE system of the N-body problem, as a function of the positions  $q_1, \dots, q_N \in \mathbb{R}^3$  and the velocities  $v_1, \dots, v_N \in \mathbb{R}^3$  of the  $N$  bodies is:

$$\begin{aligned} \frac{d}{dt} q_i &= v_i, \\ \frac{d}{dt} v_i &= G \sum_{j \neq i} \frac{m_j}{\|q_j - q_i\|^3} (q_j - q_i). \end{aligned}$$

This system of ODEs can be written in compact form as

$$\frac{du}{dt} = f(t, u, Gm)$$

```
[4]: PATH_ODES="../../ODEProblems/"

include(string(PATH_ODES, "Initial6Body.jl"))
include(string(PATH_ODES, "Nbody.jl"))
include(string(PATH_ODES, "Nbody2nd.jl"))
include(string(PATH_ODES, "NbodyDyn.jl"));
```

### 1.2.1 Initial value problem: 6-body problem (outer solar system)

We consider  $N = 6$  bodies of the outer solar system: the Sun, Jupiter, Saturn, Uranus, Neptune and Pluto. The initial values  $u_{00}$  are taken from DE430, Julian day (TDB) 2440400.5 (June 28, 1969).

```
[5]: nruns=1

u0, Gm, bodylist = Initial6Body(Float64)
u0_B, Gm_B, bodylist = Initial6Body(BigFloat)

q0=u0[:, :, 1]
v0=u0[:, :, 2]
dim=length(size(u0))

N = length(Gm)

show(bodylist)
EO=NbodyEnergy(u0,Gm)
```

```
["Sun" "Jupiter" "Saturn" "Uranus" "Neptune" "Pluto"]
```

```
[5]: -9.522620605966966e-12
```

```
[6]: # Comprobación: centro de masas del sistema solar exterior en el origen de
      ↪ coordenadas con velocidad nula

q0_ = zeros(3)
v0_ = zeros(3)
Gm0_ = 0.
for i in 1:6
    Gmi = Gm[i]
    qi = u0[:, i, 1]
    vi = u0[:, i, 2]
    q0_ += Gmi*qi
    v0_ += Gmi*vi
    Gm0_ += Gmi
end
(q0_/Gm0_, v0_/Gm0_)
```

```
[6]: ([1.417048569677335e-18, 9.119357686204585e-20, 1.5082339679877506e-19],
      [1.0942040830907497e-22, -2.709864045560887e-22, 1.6237832716049163e-22])
```

```
[7]: t0=0.
      tF=1e9 # days

tspan= (t0,tF)
tspan_B=(BigFloat(t0),BigFloat(tF))

dim=length(size(u0))
prob = ODEProblem(NbodyODE!, u0,tspan , Gm)
probDyn= DynamicalODEProblem(NbodyODE!,NbodyODEv!,q0,v0,tspan,Gm)
prob2nd = SecondOrderODEProblem(NbodyODE2nd!,v0,q0,(t0,tF),Gm)
```

```
prob_B = ODEProblem(NbodyODE!, u0_B, tspan_B, Gm_B);
```

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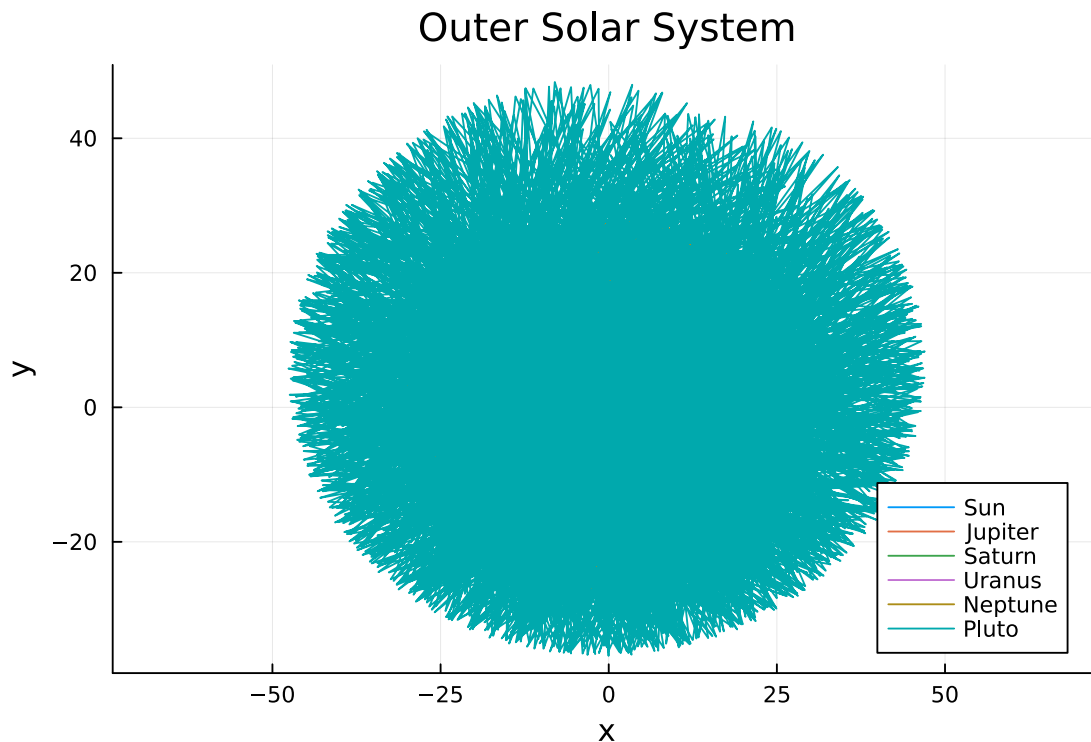
## 1.3 Integrations

### 1.3.1 IRKNGl integratioa-0 (exact solution)

```
[8]: m=5000
dt0=100.
alg=IRKNGl_simd(s=8, initial_interp=1, m=m,myoutputs=true)
saveat=m*dt0
sol0, iters0, steps0=solve(prob,alg,dt=dt0, saveat=saveat,adaptive=false);
```

```
[9]: pl = plot(title="Outer Solar System",
               xlabel="x", ylabel="y", aspect_ratio=1)

for j = 1:6
    x = [u[1,j,1] for u in sol0.u]
    y = [u[2,j,1] for u in sol0.u]
    pl = plot!(x,y, label=bodylist[j])
end
display(pl)
```



### 1.3.2 IRKNGL

```
[10]: m=1000
      dt=500.
      alg=IRKNGL_Seq(s=8, initial_interp=1, m=m, myoutputs=true)
      alg2=IRKNGL_simd(s=8, initial_interp=1, m=m, myoutputs=true)

      sol1, iters1, steps1=solve(prob, alg, dt=dt, adaptive=false);
      sol2, iters2, steps2=solve(prob, alg2, dt=dt, adaptive=false);

[11]: solve(prob, alg, dt=dt, save_everystep=false, adaptive=false);
      cpu_IRKGLseq=0.
      for i in 1:nruns
          cpu_IRKGLseq+=@elapsed solve(prob, alg, dt=dt, save_everystep=false,
          ↪ adaptive=false);
      end
      cpu_IRKGLseq=cpu_IRKGLseq/nruns

      solve(prob, alg2, dt=dt, save_everystep=false, adaptive=false);
      cpu_IRKGLsimd=0.
      for i in 1:nruns
          cpu_IRKGLsimd+=@elapsed solve(prob, alg2, dt=dt, save_everystep=false,
          ↪ adaptive=false);
      end
      cpu_IRKGLsimd=cpu_IRKGLsimd/nruns;
```

### 1.3.3 Sofronio

```
[12]: m=2000
      dt3=250.0
      method = SofSpa10

      m=4000
      dt3=250.0/2
      method = KahanLi8

      saveat=m*dt3
      sol3=solve(prob2nd, method(), dt=dt3, saveat=saveat, adaptive=false);
      T=typeof(u0)
      solu3=Array{T}(undef, length(sol3.u));
      for k in 1:length(sol3.u)
          solu3[k]=zero(u0)
          solu3[k][:, :, 1].=sol3.u[k].x[2]
          solu3[k][:, :, 2].=sol3.u[k].x[1]
      end
```

```
[13]: solve(prob2nd, method(), dt=dt3, save_everystep=false, adaptive=false);
      cpu_SofSpa10=0.
      for i in 1:nruns
          cpu_SofSpa10+=@elapsed solve(prob2nd, method(), dt=dt3,
          ↪save_everystep=false, adaptive=false)
      end
      cpu_SofSpa10=cpu_SofSpa10/nruns;
```

### 1.3.4 DPRKN12

```
[14]: m=2500
      dt4=200.
      saveat=m*dt4
      sol4=solve(prob2nd, DPRKN12(), dt=dt4, saveat=saveat,adaptive=false);
      T=typeof(u0)
      solu4=Array{T}(undef,length(sol4.u));
      for k in 1:length(sol4.u)
          solu4[k]=zero(u0)
          solu4[k][:,1].=:sol4.u[k].x[2]
          solu4[k][:,2].=:sol4.u[k].x[1]
      end
```

```
[15]: solve(prob2nd, DPRKN12(), dt=dt4, save_everystep=false, adaptive=false);
      cpu_DPRKN12=0.
      for i in 1:nruns
          cpu_DPRKN12+=@elapsed solve(prob2nd, DPRKN12(), dt=dt4,
          ↪save_everystep=false, adaptive=false)
      end
      cpu_DPRKN12=cpu_DPRKN12/nruns;
```

### 1.3.5 Some checks

```
[16]: sols=[sol0, sol1, sol2, sol3, sol4]
      for sol in sols
          println("length=",length(sol.t))
      end
```

```
length=2001
length=2001
length=2001
length=2001
length=2001
```

```
[17]: k=20
      for sol in sols[1:3]
          println("norm=", norm(sol0.u[1]-sol.u[1]), ", sol.t[k]=",sol.t[k], ", sol.
          ↪t[end]=", sol.t[end])
```

```
end
```

```
norm=0.0, sol.t[k]=9.5e6, sol.t[end]=1.0e9  
norm=0.0, sol.t[k]=9.5e6, sol.t[end]=1.0e9  
norm=0.0, sol.t[k]=9.5e6, sol.t[end]=1.0e9
```

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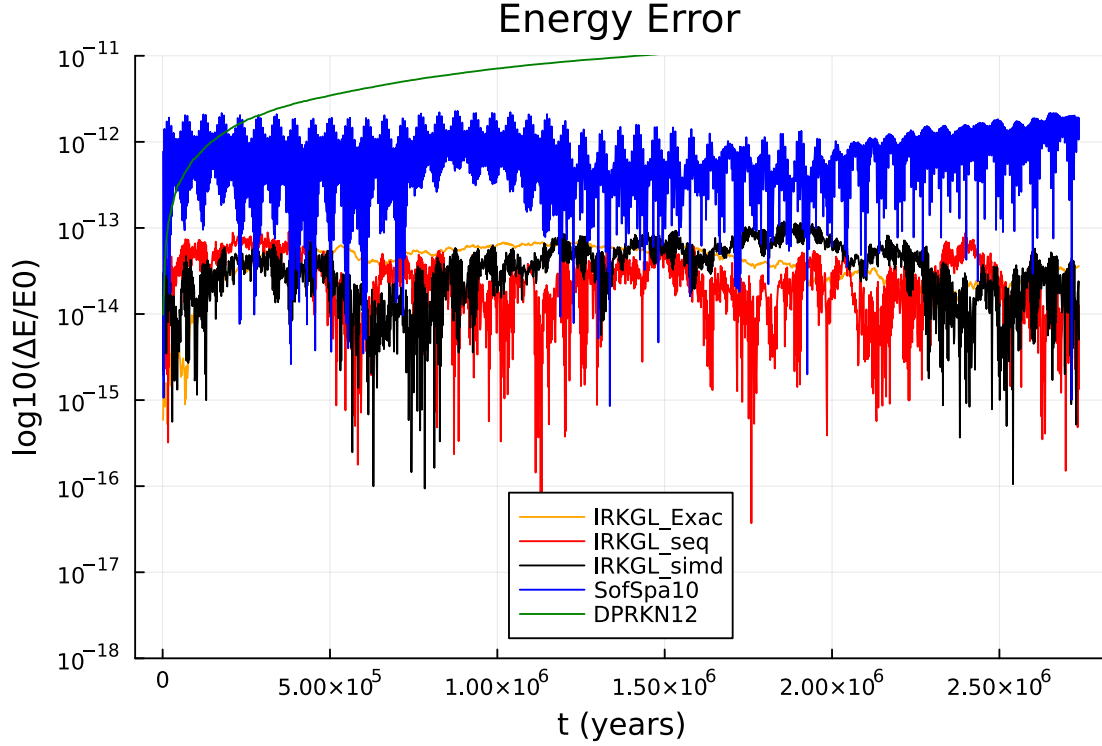
## 1.4 Errors in energy

```
[18]: println("CPU IRKNGl_seq=$cpu_IRKGLseq, CPU IRKNGl_simd=$cpu_IRKGLsimd, CPU_  
      ↪IRKNGl_seq/CPU IRKNGl_simd=$(cpu_IRKGLseq/cpu_IRKGLsimd)")  
println("CPU SofSpa10=$cpu_SofSpa10, CPU DPRKN12=$cpu_DPRKN12" )
```

```
CPU IRKNGl_seq=52.813707364, CPU IRKNGl_simd=9.954895036, CPU IRKNGl_seq/CPU  
IRKNGl_simd=5.305300274187642  
CPU SofSpa10=14.770874216, CPU DPRKN12=10.259620822
```

```
[19]: E0=NbodyEnergy(u0_B, Gm_B)  
year=365.25  
yrange=(1e-18,1e-11)  
ΔE0 = map(x->NbodyEnergy(BigFloat.(x),Gm_B), sol0.u)./E0.-1  
ΔE1 = map(x->NbodyEnergy(BigFloat.(x),Gm_B), sol1.u)./E0.-1  
ΔE2 = map(x->NbodyEnergy(BigFloat.(x),Gm_B), sol2.u)./E0.-1  
ΔE3 = map(x->NbodyEnergy(BigFloat.(x),Gm_B), solu3)./E0.-1  
ΔE4 = map(x->NbodyEnergy(BigFloat.(x),Gm_B), solu4)./E0.-1  
plot(title="Energy Error",xlabel="t (years)", ylabel="log10(ΔE/E0)",  
      yscale=:log10, ylims=yrange, legend=:bottom)  
plot!(sol0.t./year,abs.(ΔE0),label="IRKGL_Exac", color="orange")  
plot!(sol1.t./year,abs.(ΔE1),label="IRKGL_seq", color="red")  
plot!(sol2.t./year,abs.(ΔE2),label="IRKGL_simd", color="black")  
plot!(sol3.t./year,abs.(ΔE3),label="SofSpa10", color="blue")  
plot!(sol4.t./year,abs.(ΔE4),label="DPRKN12", color="green")
```

[19]:



```
[20]: Jupiter_period = 11.9*365
      tF/Jupiter_period
```

```
[20]: 230229.0779325429
```

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## 1.5 Errors in position

```
[21]: qe1=Array{Array{Float64,1}}(undef,N)
      qe2=Array{Array{Float64,1}}(undef,N)
      qe3=Array{Array{Float64,1}}(undef,N)
      qe4=Array{Array{Float64,1}}(undef,N)
      for i in 1:N
          qe1[i]=map((u0,u1)-> log10(norm(u0[:,i,1]-u1[:,i,1])), sol0.u[2:end],sol1.
          ↪u[2:end])
          qe2[i]=map((u0,u1)-> log10(norm(u0[:,i,1]-u1[:,i,1])), sol0.u[2:end],sol2.
          ↪u[2:end])
          qe3[i]=map((u0,u1)-> log10(norm(u0[:,i,1]-u1[:,i,1])), sol0.u[2:
          ↪end],solu3[2:end])
          qe4[i]=map((u0,u1)-> log10(norm(u0[:,i,1]-u1[:,i,1])), sol0.u[2:
          ↪end],solu4[2:end])
      end
```



```
[22]: MinE=-18
MaxE=-4

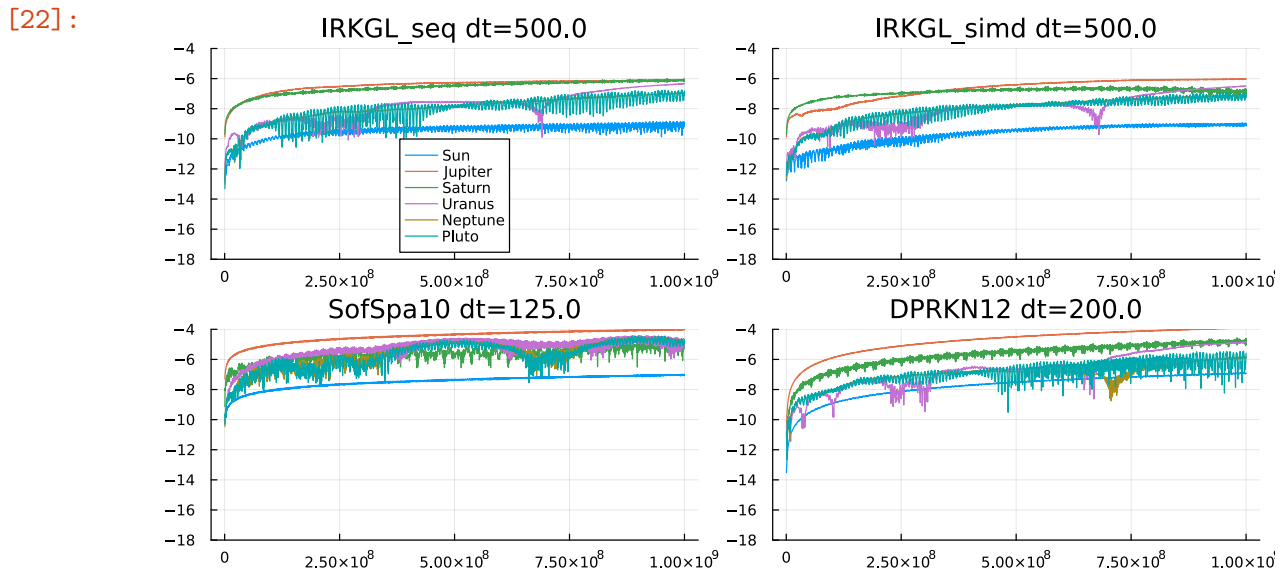
pe1=plot(sol0.t[2:end], qe1,
         title="IRKGL_seq dt=$dt",
         ylims=(MinE,MaxE),
         label=bodylist,legend=:bottom)

pe2=plot(sol0.t[2:end], qe2,
         title="IRKGL_simd dt=$dt",
         ylims=(MinE,MaxE),
         legend=false)

pe3=plot(sol0.t[2:end], qe3,
         title="SofSpa10 dt=$dt3",
         ylims=(MinE,MaxE),
         legend=false)

pe4=plot(sol0.t[2:end], qe4,
         title="DPRKN12 dt=$dt4",
         ylims=(MinE,MaxE),
         legend=false)

plot(pe1, pe2, pe3, pe4,
     layout=(2,2), size=(900,450),plot_title="")
```



[ ]:

[ ]: