

```

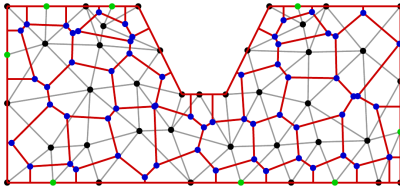
- begin
-   ENV["LANG"]="C"
-   using Pkg
-   Pkg.activate(mktempdir())
-   using Revise
-   Pkg.add("Revise")
-   Pkg.add(["PyPlot", "PlutoUI", "ExtendableGrids", "GridVisualize", "VoronoiFVM"])
-   using PlutoUI, PyPlot, ExtendableGrids, VoronoiFVM, GridVisualize
-   PyPlot.svg(true)
- end;

```

Finite volumes: transient problems

Construction of control volumes

- Start with a triangulation of a polygonal domain (intervals in 1D, triangles in 2D, tetrahedra in 3D).
- Join triangle circumcenters by lines → create Voronoi cells which can serve as control volumes, akin to representative elementary volumes (REV) used to derive conservation laws.



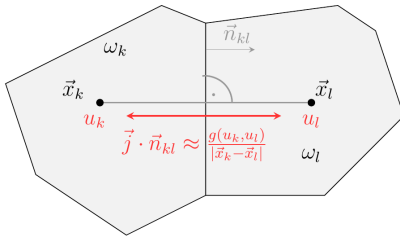
- Black + green: triangle nodes
- Gray: triangle edges
- Blue: triangle circumcenters
- Red: Boundaries of Voronoi cells

Condition on triangulation

- There is a 1:1 incidence between triangulation nodes and Voronoi cells. Moreover, the angle between the interface between two Voronoi cells and the edge between their corresponding nodes is $\frac{\pi}{2}$.
- Requires (in 2D) that sums of angles opposite to triangle edges are less than π and that angles opposite to boundary edges are less than $\frac{\pi}{2}$.
- "boundary conforming Delaunay property". It has different equivalent definitions and analogues in 3D.
- Construction:
 - "by hand" (or script) from tensor product meshes
 - Mesh generators: Triangle, TetGen
 - Julia packages: Triangulate.jl, TetGen.jl; SimplexGridFactory.jl

The discretization approach

- Use Voronoi cells as REVs aka control volumes aka finite volume cells.



- Given a continuity equation $\nabla \cdot \vec{j} = 0$ in a domain Ω , integrate this over a control volume ω_k with associated node \vec{x}_k and apply Gauss theorem:

$$\begin{aligned} 0 &= \int_{\omega_k} \nabla \cdot \vec{j} \, d\omega = \int_{\partial\omega_k} \vec{j} \cdot \vec{n} \, ds \\ &= \sum_{l \in N_k} \int_{\omega_k \cap \omega_l} \vec{j} \cdot \vec{n} \, ds + \int_{\partial\omega_k \cap \partial\Omega} \vec{j} \cdot \vec{n} \, ds \\ &\approx \sum_{l \in N_k} \frac{\sigma_{kl}}{h_{kl}} g(u_k, u_l) + \gamma_k b(u_k) \end{aligned}$$

- Here, N_k is the set of neighbor control volumes, $\sigma_{kl} = |\omega_k \cap \omega_l|$, $h_{kl} = |\vec{x}_k - \vec{x}_l|$, $\gamma_k = |\partial\omega_k \cap \partial\Omega|$, where $|\cdot|$ denotes the measure (length resp. area) of a geometrical entity.

Flux functions

For instance, for the diffusion flux $\vec{j} = -D\nabla u$, we use $g(u_k, u_l) = D(u_k - u_l)$.

For a convective diffusion flux $\vec{j} = -D\nabla u + u\vec{v}$, one can chose the upwind flux

$$g(u_k, u_l) = D(u_k - u_l) + v_{kl} \begin{cases} u_k, & v_{kl} > 0 \\ u_l, & v_{kl} \leq 0, \end{cases}$$

where $v_{kl} = \frac{h_{kl}}{\sigma_{kl}} \int_{\omega_k \cap \omega_l} \vec{v} \cdot \vec{n}_{kl} \, ds$ Fluxes also can depend nonlinearly on u .

Software API and implementation

$$\partial_t s(u) + \nabla \cdot \vec{j}(u) + r(u) = f$$

The entities describing the discrete system can be subdivided into two categories:

- geometrical data: $|\omega_k|, \gamma_k, \sigma_{kl}, h_{kl}$ together with the connectivity information of the triangles
- physical data: the number m and the functions s, g, r, f describing the particular problem, where g is a flux function approximating \vec{j} .

This structure allows to describe the problem to be solved by data derived from the discretization grid and by the functions describing the physics, giving rise to a software API.

The solution of the nonlinear systems of equations can be performed by Newton's method combined with various direct and iterative linear solvers.

The generic programming capabilities of Julia allow for an implementation of the method which results in an API which consists in the implementation of functions s, g, r, f without the need to write code for their derivatives.

Examples

General settings

Initial value problem with homogeneous Neumann boundary conditions

$$\Omega = (0, 1)^d, \quad d = 1, 2$$

$$T = [0, t_{end}]$$

evolution (generic function with 1 method)

```

. # Function describing evolution of system with initial value inival
. # using the Implicit Euler method
. function evolution(inival, # initial value
.     sys, # finite volume system
.     grid, # simplex grid
.     timestep, # initial time step
.     tend, # end time
.     dtgrowth # time step growth factor
. )
.     time=0.0
.     # record time and solution
.     times=[time]
.     solutions=[copy(inival)]
.
.     solution=copy(inival)
.     while time<tend
.         time=time+timestep
.         solve!(solution,inival,sys,timestep) # solve implicit Euler time step
.         inival.=solution # copy solution to inival
.         push!(times,time)
.         push!(solutions,copy(solution))
.         timestep=dtgrowth # increase timestep by factor when approaching stationary
.     state
.     end
.     # return result and grid
.     (times=times,solutions=solutions,grid=grid)
. end
. 
```

fpeak (generic function with 2 methods)

```

. # Define function for initial value $u_0$ with two methods - for 1D and 2D problems
. begin
.     fpeak(x)=exp(-100*(x-0.25)^2)
.     fpeak(x,y)=exp(-100*((x-0.25)^2+(y-0.25)^2))
. end
. 
```

create_grid (generic function with 1 method)

```

. # Create discretization grid in 1D or 2D with approximately n nodes
. function create_grid(n,dim)
.     nx=n
.     if dim==2
.         nx=ceil(sqrt(n))
.     end
.     X=collect(0:1.0/nx:1)
.     if dim==1
.         grid=simplexgrid(X)
.     else
.         grid=simplexgrid(X,X)
.     end
. end
. 
```

Diffusion problem

$$\partial_t u - \nabla \cdot D \nabla u = 0 \text{ in } \Omega$$

$$D \nabla u \cdot \vec{n} = 0 \text{ on } \partial \Omega$$

$$u|_{t=0} = u_0$$

diffusion (generic function with 1 method)

```

. function diffusion(;n=100,dim=1,timestep=1.0e-4,tend=1, D=1.0, dtgrowth=1.1)
.     grid=create_grid(n,dim)
.
.     ## Diffusion flux between neighboring control volumes
.     function flux!(f,u,edge)
.         uk=viewK(edge,u)
.         ul=viewL(edge,u)
.         f[1]=D*(uk[1]-ul[1])
.     end
.
.     ## Storage term (under time derivative)
.     function storage!(f,u,node)
.         f[1]=u[1]
.     end
.
.     ## Create a physics structure
. 
```

```

    physics=VoronoiFVM.Physics(
        flux=flux!,
        storage=storage!)
    sys=VoronoiFVM.DenseSystem(grid,physics)
    enable_species!(sys,1,[1])
    ## Create a solution array
    inival=unknowns(sys)
    ## Broadcast the initial value
    inival[1,:].=map(fpeak,grid)
    evolution(inival,sys,grid,tstep,tend,dtgrowth)
end

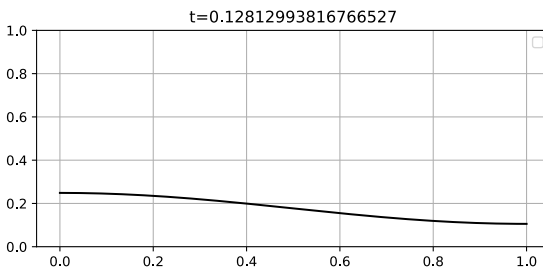
```

```

result_diffusion=diffusion(dim=1,n=1000);

```

time=



Reaction-diffusion problem

Diffusion + physical process which 'eats' species

$$\partial_t u - \nabla \cdot D \nabla u + Ru = 0 \text{ in } \Omega$$

$$D \nabla u \cdot \vec{n} = 0 \text{ on } \partial \Omega$$

$$u|_{t=0} = u_0$$

reaction_diffusion (generic function with 1 method)

```

function reaction_diffusion(;
    n=1000,
    dim=1,
    tstep=1.0e-4,
    tend=1,
    D=1.0,
    R=10.0,
    dtgrowth=1.1)
    grid=create_grid(n,dim)
    ## Diffusion flux between neighboring control volumes
    function flux!(f,u,edge)
        uk=viewK(edge,u)
        ul=viewL(edge,u)
        f[1]=D*(uk[1]-ul[1])
    end
    ## Storage term (under time derivative)
    function storage!(f,u,node)
        f[1]=u[1]
    end
    ## Reaction term
    function reaction!(f,u,node)
        f[1]=R*u[1]
    end
    ## Create a physics structure
    physics=VoronoiFVM.Physics(
        flux=flux!,
        reaction=reaction!,
        storage=storage!)
end

```

```

- sys=VoronoiFVM.DenseSystem(grid,physics)
- enable_species!(sys,1,[1])
- ## Create a solution array
- inival=unknowns(sys)
-
- ## Broadcast the initial value
- inival[1,:].=map(fpeak,grid)
-
- evolution(inival,sys,grid,tstep,tend,dtgrowth)
- end

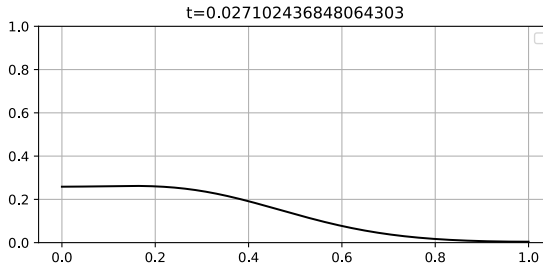
```

```

- result_reaction_diffusion=reaction_diffusion(dim=1,n=1000,R=10);

```

time=



Convection-Diffusion problem

$$\partial_t u - \nabla \cdot (D \nabla u - u \vec{v}) = 0 \text{ in } \Omega$$

$$(D \nabla u - u \vec{v}) \cdot \vec{n} = 0 \text{ on } \partial \Omega$$

$$u|_{t=0} = u_0$$

```

convection_diffusion (generic function with 1 method)
- function convection_diffusion(;
-     n=20,
-     dim=1,
-     timestep=1.0e-4,
-     tend=1,
-     D=0.01,
-     vx=10.0,
-     vy=10.0,
-     dtgrowth=1.1,
-     scheme="expfit")
- grid=create_grid(n,dim)
- ## copy vx, vy into vector
- if dim==1
-     V=[vx]
- else
-     V=[vx,vy]
- end
- ## Bernoulli function
-
- B(x)=x/(exp(x)-1)
-
- function flux_expfit!(f,u,edge)
-     uk=viewK(edge,u)
-     ul=viewL(edge,u)
-     vh=project(edge,V) ## Calculate projection v * (x_L-x_K)
-     f[1]=D*(B(-vh/D)*uk[1]- B(vh/D)*ul[1])
- end
-
- function flux_centered!(f,u,edge)
-     uk=viewK(edge,u)
-     ul=viewL(edge,u)
-     vh=project(edge,V)
-     f[1]=D*(uk[1]-ul[1])+ vh*0.5*(uk[1]+ul[1])
- end
-
- function flux_upwind!(f,u,edge)
-     uk=viewK(edge,u)
-     ul=viewL(edge,u)
-     vh=project(edge,V)
-     f[1]=D*(uk[1]-ul[1])+ (vh>0.0 ? vh*uk[1] : vh*ul[1])
- end

```

```

.
.
.   flux! =flux_upwind!
.   if scheme=="expfit"
.       flux! =flux_expfit!
.   elseif scheme=="centered"
.       flux! =flux_centered!
.   end
.
.
.   ## Storage term (under time derivative)
.   function storage!(f,u,node)
.       f[1]=u[1]
.   end
.
.
.   ## Create a physics structure
.   physics=VoronoiFVM.Physics(
.       flux=flux!,
.       storage=storage!)
.
.   sys=VoronoiFVM.DenseSystem(grid,physics)
.   enable_species!(sys,1,[1])
.   ## Assume homogeneous Neumann boundary conditions, so do nothing
.
.   ## Create a solution array
.   inival=unknowns(sys)
.
.   ## Broadcast the initial value
.   inival[1,:].=map(fpeak,grid)
.   evolution(inival,sys,grid,tstep,tend,dtgrowth)
. end
.

```

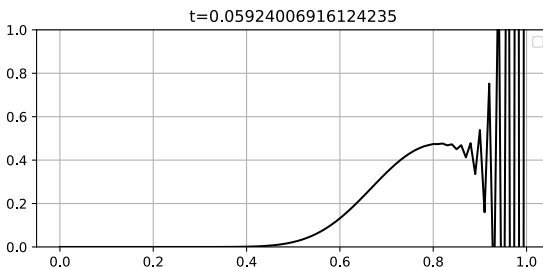
scheme:

```

. result_convection_diffusion=convection_diffusion(n=100,
. dim=1,
. scheme=scheme[1],
. vx=10,vy=10);

```

time=



Brusselator system

Two species interacting via a reaction:

$$\begin{aligned} \partial_t u_1 - \nabla \cdot (D_1 \nabla u_1) + (B+1)u_1 - A - u_1^2 u_2 &= 0 \\ \partial_t u_2 - \nabla \cdot (D_2 \nabla u_2) + u_1^2 u_2 - B u_1 &= 0 \end{aligned}$$

```

brusselator (generic function with 1 method)
. function brusselator(;n=100,dim=1,A=4.0,B=6.0,D1=0.01,D2=0.1,perturbation=0.1,
. tstep=0.05, tend=150,dtgrowth=1.05)
.
.   grid=create_grid(n,dim)
.   function storage!(f,u,node)
.       f.=u
.   end
.
.
.   function bruss_diffusion!(f,u,edge)
.       u=unknowns(edge,_u)
.       f[1]=D1*(u[1,1]-u[1,2])
.       f[2]=D2*(u[2,1]-u[2,2])
.   end

```

```

- # Reaction:
- function bruss_reaction!(f,u,node)
-     f[1]= (B+1.0)*u[1]-A-u[1]^2*u[2]
-     f[2]= u[1]^2*u[2]-B*u[1]
- end
- # Create system
- bruss_physics=VoronoiFVM.Physics(flux=bruss_diffusion!, storage=storage!,
-                                 num_species=2, reaction=bruss_reaction!)
- brusselator_system=VoronoiFVM.DenseSystem(grid,bruss_physics)
- enable_species!(brusselator_system,1,[1])
- enable_species!(brusselator_system,2,[1])
-
- inival=unknowns(brusselator_system)
- for i=1:num_nodes(grid)
-     inival[1,i]=1.0+perturbation*randn()
-     inival[2,i]=1.0+perturbation*randn()
- end
-
- evolution(inival,brusselator_system,grid,tstep,tend,dtgrowth)
- end

```

```

- result_brusselator=brusselator(n=500,dim=1);

```

time=

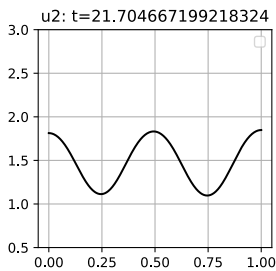
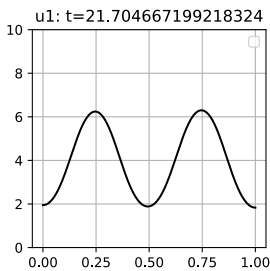


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```

Status `~/tmp/jl_op3LMf/Project.toml`
 [cfc395e8] ExtendableGrids v0.7.4
 [5eed8a63] GridVisualize v0.1.3
 [7f904dfe] PlutoUI v0.7.2
 [d330b81b] PyPlot v2.9.0
 [295af30f] Revise v3.1.12
 [82b139dc] VoronoiFVM v0.10.5

```