

Immersed Boundary Tutorial

Part 1

Aaron Barrett

July 22, 2025

- What to expect:

- What not to expect:

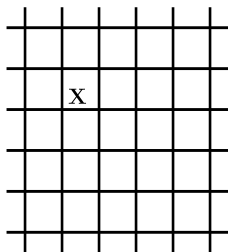
- What to expect:
 - Introduction to IB
 - Introduction to IBAMR
 - Run IBAMR examples
 - Visualize IBAMR output
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 - Basic understanding of design, example layout, input file
 - Basis to start exploring and considering implementation
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- What not to expect:
 - Expertise in IBAMR
 - Expertise in C++
 - Expertise in VisIt

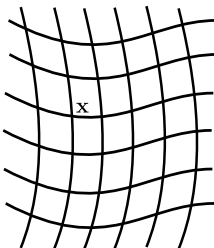
Lagrangian vs Eulerian Variables

Reference Configuration



$\Omega(0)$

Current Configuration



$\Omega(t)$

$\chi(\mathbf{X}, t)$

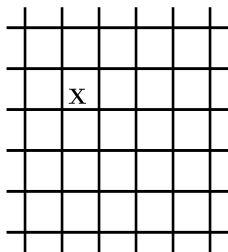
- Fluid parcels are associated with a label $\mathbf{X}$.
- Velocity in a Lagrangian frame

- Spatial coordinates $\mathbf{x}$ are fixed.
- Eulerian velocity: $\mathbf{u}(\mathbf{x}, t)$.

$$\frac{\partial \chi(\mathbf{X}, t)}{\partial t}$$

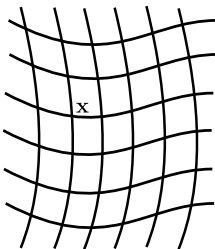
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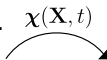


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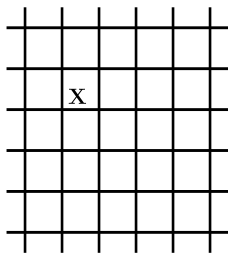
$$\frac{\partial \chi(\mathbf{X}, t)}{\partial t}$$

- Spatial coordinates $\mathbf{x}$ are fixed.
- Eulerian velocity: $\mathbf{u}(\mathbf{x}, t)$.
- We have that

$$\begin{aligned} \frac{\partial \chi(\mathbf{X}, t)}{\partial t} &= \mathbf{u}(\chi(\mathbf{X}, t), t) \\ &= \int_{\Omega} \mathbf{u}(\mathbf{x}, t) \delta(\mathbf{x} - \chi(\mathbf{X}, t)) d\mathbf{x} \end{aligned}$$

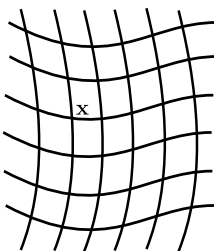
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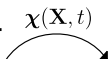


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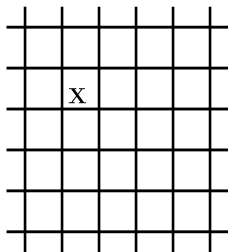
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- Rates of change of a scalar variable
 $f(\mathbf{x}, t) = f(\chi(\mathbf{X}, t), t) = F(\mathbf{X}, t)$:

$$\frac{d}{dt} f(\chi(\mathbf{X}, t), t) = \frac{\partial f}{\partial t} + \frac{\partial \chi}{\partial t} \frac{\partial f}{\partial \chi}$$

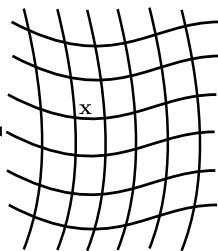
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$\chi(\mathbf{X}, t)$

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$$\frac{d}{dt} f(\chi(\mathbf{X}, t), t) = \left(\frac{\partial}{\partial t} + \mathbf{u} \cdot \nabla \right) f$$

- We can define the acceleration of the fluid in two ways:
 - Lagrangian:

$$\frac{\partial^2 \chi}{\partial t^2}$$

- Eulerian:

$$\frac{D\mathbf{u}}{Dt} = \frac{\partial \mathbf{u}}{\partial t} + \mathbf{u} \cdot \nabla \mathbf{u}$$

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- The deformation gradient $\mathbb{F} = \frac{\partial \chi}{\partial \mathbf{X}}$ maps line elements $d\mathbf{X}$ in the reference configuration to line elements $d\mathbf{x}$ in the current configuration:

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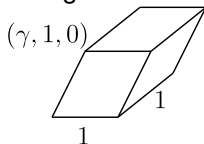
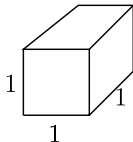
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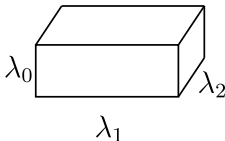
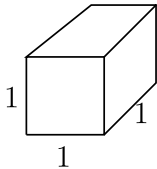
$$d\mathbf{x} = \mathbb{F} \cdot d\mathbf{X}$$

- Consider a cube being *sheared*.



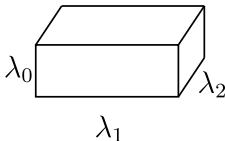
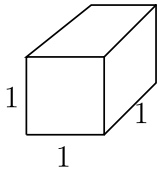
$$\mathbb{F} = \begin{pmatrix} 1 & 0 & 0 \\ \gamma & 1 & 0 \\ 0 & 0 & 1 \end{pmatrix}$$

- Consider a cube being stretched (extensional deformation).



$$\mathbb{F} = \begin{pmatrix} \lambda_0 & 0 & 0 \\ 0 & \lambda_1 & 0 \\ 0 & 0 & \lambda_2 \end{pmatrix}$$

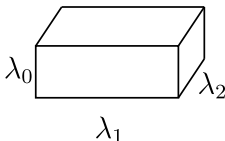
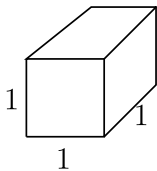
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- For $\mathbb{F}$ to represent an *incompressible deformation*, we need $\lambda_0 \lambda_1 \lambda_2 = 1$.
- The Jacobian $J = \det \mathbb{F}$ measures volume changes:

$$\frac{dJ}{dt} = (\nabla \cdot \mathbf{u})J$$

- Incompressible materials satisfy $J = 1$ or $\nabla \cdot \mathbf{u} = 0$.

- Stress measures the response of material to strain.
- Traction on a surface with normal $\mathbf{n}$ is

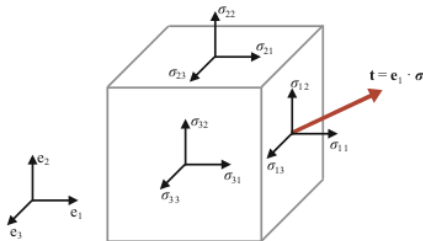
$$\mathbf{t} = \boldsymbol{\sigma} \cdot \mathbf{n}$$

- Net force on surface S

$$\mathbf{F} = \int_{\partial S} \boldsymbol{\sigma} \cdot \mathbf{n} dS$$

- Conservation of angular momentum:

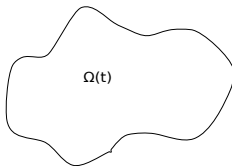
$$\boldsymbol{\sigma} = \boldsymbol{\sigma}^T$$



Conservation of Mass

- Consider a volume of fluid $\Omega(t)$ with density $\rho(\mathbf{x}, t)$. Total mass:

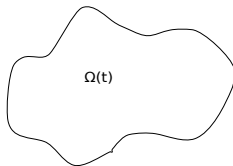
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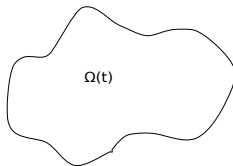


- Conservation of mass:

$$\begin{aligned} 0 &= \frac{d}{dt} \int_{\Omega(t)} \rho(\mathbf{x}, t) dV_{\mathbf{x}} = \frac{d}{dt} \int_{\Omega(0)} \rho(\chi(\mathbf{X}, t), t) J dV_{\mathbf{X}} \\ &= \int_{\Omega(0)} \left(\frac{D\rho(\chi(\mathbf{X}, t), t)}{Dt} + \rho(\chi(\mathbf{X}, t), t) (\nabla \cdot \mathbf{u}) \right) J dV_{\mathbf{X}} \\ &= \int_{\Omega(t)} \left(\frac{D\rho(\mathbf{x}, t)}{Dt} + \rho(\mathbf{x}, t) (\nabla \cdot \mathbf{u}) \right) dV_{\mathbf{x}} \end{aligned}$$

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- This holds for all material volumes:

$$\frac{D\rho}{Dt} + \rho(\nabla \cdot \mathbf{u}) = 0 \quad \text{or} \quad \frac{D\rho}{Dt} = 0$$

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$$\mathbf{p}(t) = \int_{\Omega(t)} \rho(\mathbf{x}, t) \mathbf{u}(\mathbf{x}, t) dV_{\mathbf{x}}$$

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- Forces on fluid:
 - External body forces $\mathbf{f}$, e.g. gravity.
 - Surface forces $\mathbf{t} = \mathbf{n} \cdot \boldsymbol{\sigma}$ e.g. viscous or elastic forces.
 - $\boldsymbol{\sigma}$ is the Cauchy stress tensor.

- Rate of change of momentum:

$$\frac{d}{dt} \mathbf{p}(t) = \int_{\Omega(t)} \rho \frac{D\mathbf{u}}{Dt} dV_{\mathbf{x}}$$

- Forces on fluid:

$$\int_{\Omega(t)} \mathbf{f} dV_{\mathbf{x}} + \int_{\partial\Omega(t)} \mathbf{t} dS_{\mathbf{x}}$$

- Newton's second law:

$$\int_{\Omega(t)} \rho \frac{D\mathbf{u}}{Dt} dV_{\mathbf{x}} = \int_{\Omega(t)} \mathbf{f} dV_{\mathbf{x}} + \int_{\partial\Omega(t)} \mathbf{t} dS_{\mathbf{x}}$$

or

$$\rho \frac{D\mathbf{u}}{Dt} = \mathbf{f} + \nabla \cdot \boldsymbol{\sigma}$$

- Different fluids of interest reduce to specifying a constitutive equation for $\boldsymbol{\sigma}$.
- We first decompose the stress into *hydrostatic pressure* p and the *deviatoric stress* $\boldsymbol{\tau}$.

$$\boldsymbol{\sigma} = -p\mathbb{I} + \boldsymbol{\tau}$$

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- Velocity gradient:

$$\frac{\partial}{\partial t} \mathbb{F} = \frac{\partial}{\partial t} \frac{\partial \mathbf{x}}{\partial \mathbf{X}} = \frac{\partial \mathbf{u}}{\partial \mathbf{x}} \frac{\partial \mathbf{x}}{\partial \mathbf{X}} = \nabla \mathbf{u} \cdot \mathbb{F}$$

- Extensional and shear flows:

$$\nabla \mathbf{u} = \begin{pmatrix} \dot{\epsilon}_1 & 0 & 0 \\ 0 & \dot{\epsilon}_2 & 0 \\ 0 & 0 & \dot{\epsilon}_3 \end{pmatrix}, \quad \nabla \mathbf{u} = \begin{pmatrix} 0 & 0 & 0 \\ \dot{\gamma} & 0 & 0 \\ 0 & 0 & 0 \end{pmatrix}$$

- Rate of strain:

$$\dot{\boldsymbol{\epsilon}} = (\nabla \mathbf{u} + (\nabla \mathbf{u})^T)$$

- Vorticity tensor:

$$\boldsymbol{\omega} = (\nabla \mathbf{u} - (\nabla \mathbf{u})^T)$$

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- Navier-Stokes equations:

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- Vorticity tensor:

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$$\rho \left(\frac{\partial \mathbf{u}}{\partial t} + \mathbf{u} \cdot \nabla \mathbf{u} \right) = \nabla \cdot \boldsymbol{\sigma}$$

$$\boldsymbol{\sigma} = -p\mathbb{I} + \mu \dot{\boldsymbol{\epsilon}}, \quad \nabla \cdot \mathbf{u} = 0$$

or that

$$\rho \left(\frac{\partial \mathbf{u}}{\partial t} + \mathbf{u} \cdot \nabla \mathbf{u} \right) = -\nabla p + \mu \Delta \mathbf{u}$$

- Consider a thin, massless boundary immersed in the fluid.
- Track the interface in Lagrangian coordinates:

- Velocity:

$$\frac{\partial \chi(\mathbf{X}, t)}{\partial t} = \int_{\Omega} \mathbf{u}(\mathbf{x}, t) \delta(\mathbf{x} - \chi(\mathbf{X}, t)) d\mathbf{x}$$

- Elastic energy:

$$E[\chi] \rightarrow \text{Depends on deformation!}$$

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$$\delta E = \int -\mathbf{F} \delta \chi d\mathbf{X}$$

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- Ex: Let q parameterize $\chi(q, t)$. Suppose

$$E = \int \frac{k}{2} \left(\left| \frac{\partial \chi}{\partial q} \right| - 1 \right)^2 dq,$$

Energy is minimized when tangent vector is magnitude 1

- Forces are given by derivative of energy and q parameterizes $\chi(q, t)$,

$$\delta E = \int -\mathbf{F} \delta \chi dq, \quad E = \int \frac{k}{2} \left(\left| \frac{\partial \chi}{\partial q} \right| - 1 \right)^2 dq$$

- Then

$$\begin{aligned} \delta E &= \int \delta \left[\left(\left| \frac{\partial \chi}{\partial q} \right| - 1 \right)^2 \right] dq \\ &= \int k \left(\left| \frac{\partial \chi}{\partial q} \right| - 1 \right) \frac{\partial \chi / \partial q}{|\partial \chi / \partial q|} \frac{\partial}{\partial q} \delta \chi dq \\ &= \int -\frac{\partial}{\partial q} \left(k \left(\left| \frac{\partial \chi}{\partial q} \right| - 1 \right) \frac{\partial \chi / \partial q}{|\partial \chi / \partial q|} \right) \delta \chi dq \end{aligned}$$

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- This gives us that

$$F = \frac{\partial}{\partial q} \left(k \left(\left| \frac{\partial \chi}{\partial q} \right| - 1 \right) \frac{\partial \chi / \partial q}{|\partial \chi / \partial q|} \right)$$

- Consider a *thin massless interface*.
 - Structure moves according to local velocity

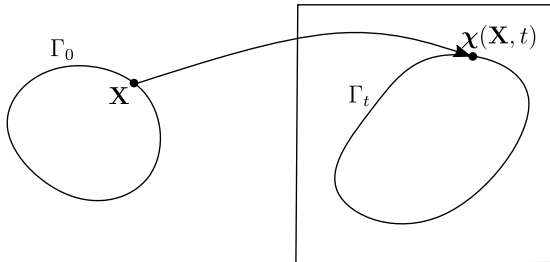
$$\frac{\partial \chi}{\partial t} = \int_{\Omega} \mathbf{u}(\mathbf{x}, t) \delta(\mathbf{x} - \chi(\mathbf{X}, t)) d\mathbf{x}$$

- Lagrangian force density is

$$\mathbf{F}(\mathbf{X}, t) = \frac{\delta E}{\delta \chi}$$

- Eulerian force density is

$$\mathbf{f}(\mathbf{x}, t) = \int_{\Gamma} \mathbf{F}(\mathbf{X}, t) \delta(\mathbf{x} - \chi(\mathbf{X}, t)) d\mathbf{X}$$



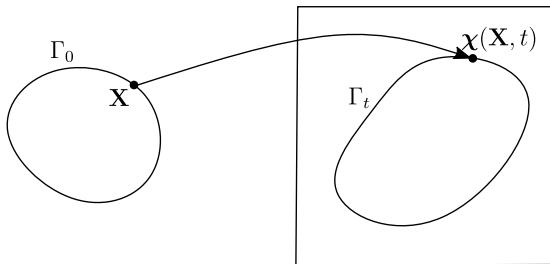
- The structure is *immersed* in a fluid

$$\frac{\partial \mathbf{u}}{\partial t} + \mathbf{u} \cdot \nabla \mathbf{u} = \nabla \cdot \boldsymbol{\sigma} + \mathbf{f}$$

$$\nabla \cdot \mathbf{u} = 0$$

$$\frac{\partial \boldsymbol{\chi}}{\partial t} = \int_{\Omega} \mathbf{u}(\mathbf{x}, t) \delta(\mathbf{x} - \boldsymbol{\chi}(\mathbf{X}, t)) d\mathbf{x}$$

$$\mathbf{f}(\mathbf{x}, t) = \int_{\Gamma} \mathbf{F}(\mathbf{X}, t) \delta(\mathbf{x} - \boldsymbol{\chi}(\mathbf{X}, t)) d\mathbf{X}$$



- IB equations:

$$\frac{\partial \mathbf{u}}{\partial t} + \mathbf{u} \cdot \nabla \mathbf{u} = \nabla \cdot \boldsymbol{\sigma} + \mathbf{f}$$

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- Solution strategy:
 - *Explicit structure, implicit fluid*
 - Move structure by interpolating velocity
 - Compute forces from structure deformation
 - Spread forces to Eulerian grid
 - Solve Navier-Stokes equations

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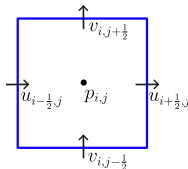
$$\mathbf{f}(\mathbf{x}, t) = \int_{\Gamma} \mathbf{F}(\mathbf{X}, t) \delta(\mathbf{x} - \chi(\mathbf{X}, t)) d\mathbf{X}$$

- Navier-Stokes solution
- Discretize operators with centered differences on MAC grid.
- Discretize using *forward and backward Euler*

$$\frac{\mathbf{u}^{n+1} - \mathbf{u}^n}{\Delta t} + \mathbf{u}^n \cdot \nabla_h \mathbf{u}^n = -\nabla_h p^{n+1} + \mu \Delta_h \mathbf{u}^{n+1} + \mathbf{f}^n$$

$$\nabla_h \cdot \mathbf{u}^{n+1} = 0$$

- Linear system to solve for $\mathbf{u}^{n+1}$ and p^{n+1} .



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$$\nabla \cdot \mathbf{u} = 0$$

$$\frac{\partial \boldsymbol{\chi}}{\partial t} = \int_{\Omega} \mathbf{u}(\mathbf{x}, t) \delta(\mathbf{x} - \boldsymbol{\chi}(\mathbf{X}, t)) d\mathbf{x}$$

$$\mathbf{f}(\mathbf{x}, t) = \int_{\Gamma} \mathbf{F}(\mathbf{X}, t) \delta(\mathbf{x} - \boldsymbol{\chi}(\mathbf{X}, t)) d\mathbf{X}$$

- Lagrangian discretization:
 - Structure is point cloud (or FE mesh): about 1 Lagrangian point per Eulerian grid cell.
 - Assume $\mathbf{F} = k \frac{\partial^2 \boldsymbol{\chi}}{\partial \mathbf{X}^2}$: tension corresponding to zero rest length springs
 - Forward Euler discretization

$$\boldsymbol{\chi}^{n+1} = \boldsymbol{\chi}^n + \Delta t \mathbf{u}^n(\boldsymbol{\chi}^n)$$

$$\mathbf{F}_{\ell}^n = k \frac{\boldsymbol{\chi}_{\ell+1}^n - 2\boldsymbol{\chi}_{\ell}^n + \boldsymbol{\chi}_{\ell-1}^n}{\Delta X^2}$$

- IB equations:

$$\frac{\partial \mathbf{u}}{\partial t} + \mathbf{u} \cdot \nabla \mathbf{u} = \nabla \cdot \boldsymbol{\sigma} + \mathbf{f}$$

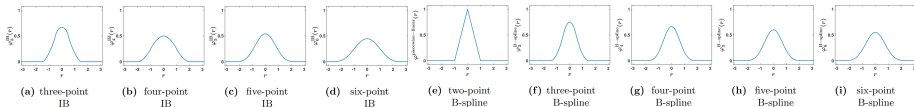
$$\nabla \cdot \mathbf{u} = 0$$

$$\frac{\partial \chi}{\partial t} = \int_{\Omega} \mathbf{u}(\mathbf{x}, t) \delta(\mathbf{x} - \chi(\mathbf{X}, t)) d\mathbf{x}$$

$$\mathbf{f}(\mathbf{x}, t) = \int_{\Gamma} \mathbf{F}(\mathbf{X}, t) \delta(\mathbf{x} - \chi(\mathbf{X}, t)) d\mathbf{X}$$

- Interaction equations:
 - Replace singular delta function with *regularized delta function*

$$\delta(\mathbf{x}) = \delta(x)\delta(y) \approx \delta_h(\mathbf{x}) = \delta_h(x)\delta_h(y)$$



- IB equations:

$$\frac{\partial \mathbf{u}}{\partial t} + \mathbf{u} \cdot \nabla \mathbf{u} = \nabla \cdot \boldsymbol{\sigma} + \mathbf{f}$$

$$\nabla \cdot \mathbf{u} = 0$$

$$\frac{\partial \boldsymbol{\chi}}{\partial t} = \int_{\Omega} \mathbf{u}(\mathbf{x}, t) \delta(\mathbf{x} - \boldsymbol{\chi}(\mathbf{X}, t)) d\mathbf{x}$$

$$\mathbf{f}(\mathbf{x}, t) = \int_{\Gamma} \mathbf{F}(\mathbf{X}, t) \delta(\mathbf{x} - \boldsymbol{\chi}(\mathbf{X}, t)) d\mathbf{X}$$

- Interaction equations:

- Replace singular delta function with *regularized delta function*

$$\delta(\mathbf{x}) = \delta(x)\delta(y) \approx \delta_h(\mathbf{x}) = \delta_h(x)\delta_h(y)$$

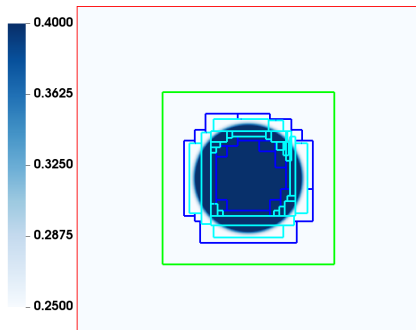
- Interpolation:

$$\mathbf{U}(\mathbf{X}_\ell) = \sum_{i,j} \mathbf{u}(\mathbf{x}_{i,j}) \delta_h(\mathbf{x}_{i,j} - \boldsymbol{\chi}(\mathbf{X}_\ell)) \Delta x^2$$

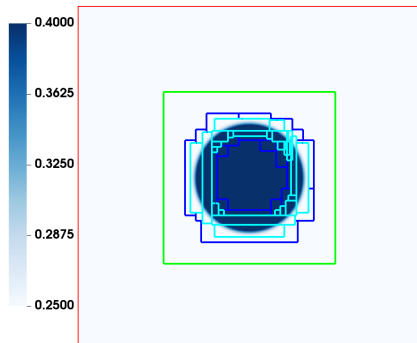
- Spreading:

$$\mathbf{f}(\mathbf{x}_{i,j}) = \sum_{\ell} \mathbf{F}(\mathbf{X}_\ell) \delta_h(\mathbf{x}_{i,j} - \boldsymbol{\chi}(\mathbf{X}_\ell)) \Delta X$$

- Immersed Boundary Adaptive Mesh Refinement (IBAMR) is
 - MPI parallelized: scalable to hundreds of cores
 - Adaptive: can tag cells based on IB location, vorticity, etc.
 - Generalizable: Support for different fluid models, multiple structure types, FSI strategies, Lagrangian discretizations
 - Easy to install: `autoibamr` will build on most systems
 - Easy to use: YMMV, but developers are reachable
- `ibamr.github.io`



- IB Method requires:
 - Integrating fluid
 - Integrating structural
 - Coupling operators
- Key principles:
 - Eulerian patch hierarchy
 - Integrator classes
 - IB model, forces, and initial position
 - Input file
- First example: `IB/explicit/ex1`: stretched elastic membrane, v0.16.0
- Elastic springs with zero rest length between each Lagrangian point



- IB Method requires:
 - Integrating fluid
 - Integrating structural
 - Coupling operators
- Key principles:
 - Eulerian patch hierarchy
 - Integrator classes
 - IB model, forces, and initial position
 - Input file
- The `PatchHierarchy<NDIM>` object maintains all Eulerian data.
- The `StandardTagAndInitialize<NDIM>` and `GriddingAlgorithm<NDIM>` objects manages AMR and regridding.

example.cpp

```

128 Pointer<CartesianGridGeometry<NDIM> > grid_geometry = new CartesianGridGeometry<NDIM>(
129     "CartesianGeometry", app_initializer->getComponentDatabase("CartesianGeometry"));
130 Pointer<PatchHierarchy<NDIM> > patch_hierarchy = new PatchHierarchy<NDIM>("PatchHierarchy", grid_geometry);
131 Pointer<StandardTagAndInitialize<NDIM> > error_detector =
132     new StandardTagAndInitialize<NDIM>("StandardTagAndInitialize",
133         time_integrator,
134         app_initializer->getComponentDatabase("StandardTagAndInitialize"));
135 Pointer<BergerRigoutsos<NDIM> > box_generator = new BergerRigoutsos<NDIM>();
136 Pointer<LoadBalancer<NDIM> > load_balancer =
137     new LoadBalancer<NDIM>("LoadBalancer", app_initializer->getComponentDatabase("LoadBalancer"));
138 Pointer<GriddingAlgorithm<NDIM> > gridding_algorithm =
139     new GriddingAlgorithm<NDIM>("GriddingAlgorithm",
140         app_initializer->getComponentDatabase("GriddingAlgorithm"),
141         error_detector,
142         box_generator,
143         load_balancer);
    
```


- IB Method requires:
 - Integrating fluid
 - Integrating structural
 - Coupling operators
- Key principles:
 - Eulerian patch hierarchy
 - Integrator classes
 - IB model, forces, and initial position
 - Input file
- HierarchyIntegrator objects can integrate specific equations
- INSStaggeredHierarchyIntegrator and IBExplicitHierarchyIntegrator are the most important for IB

example.cpp

```

102 Pointer<INSHierarchyIntegrator> navier_stokes_integrator;
103 const string solver_type =
104     app_initializer->getComponentDatabase("Main")->getStringWithDefault("solver_type", "STAGGERED");
105 if (solver_type == "STAGGERED")
106 {
107     navier_stokes_integrator = new INSStaggeredHierarchyIntegrator(
108         "INSStaggeredHierarchyIntegrator",
109         app_initializer->getComponentDatabase("INSStaggeredHierarchyIntegrator"));
110 }
111 else if (solver_type == "COLLOCATED")
112 {
113     navier_stokes_integrator = new INSCollocatedHierarchyIntegrator(
114         "INSCollocatedHierarchyIntegrator",
115         app_initializer->getComponentDatabase("INSCollocatedHierarchyIntegrator"));
116 }
117 else
118 {
119     TBOX_ERROR("Unsupported solver type: " << solver_type << "\n"
120               << "Valid options are: COLLOCATED, STAGGERED");
121 }
122 Pointer<IBMethod> ib_method_ops = new IBMethod("IBMethod", app_initializer->getComponentDatabase("IBMethod"));
123 Pointer<IBHierarchyIntegrator> time_integrator =
124     new IBExplicitHierarchyIntegrator("IBHierarchyIntegrator",
125     app_initializer->getComponentDatabase("IBHierarchyIntegrator"),
126     ib_method_ops,
127     navier_stokes_integrator);
    
```

- IB Method requires:
 - Integrating fluid
 - Integrating structural
 - Coupling operators
 - Each IB model has a different class, e.g. IBMethod, IBFEMethod, etc.
 - Initialize and setup force generator
- Key principles:
 - Eulerian patch hierarchy
 - Integrator classes
 - IB model, forces, and initial position
 - Input file

example.cpp

```

122 Pointer<IBMethod> ib_method_ops = new IBMethod("IBMethod", app_initializer->getComponentDatabase("IBMethod"));
123 Pointer<IBHierarchyIntegrator> time_integrator =
124     new IBExplicitHierarchyIntegrator("IBHierarchyIntegrator",
125         app_initializer->getComponentDatabase("IBHierarchyIntegrator"),
126         ib_method_ops,
127         navier_stokes_integrator);

145 // Configure the IB solver.
146 Pointer<IBStandardInitializer> ib_initializer = new IBStandardInitializer(
147     "IBStandardInitializer", app_initializer->getComponentDatabase("IBStandardInitializer"));
148 ib_method_ops->registerLInitStrategy(ib_initializer);
149 Pointer<IBStandardForceGen> ib_force_fcn = new IBStandardForceGen();
150 ib_method_ops->registerIBLagrangianForceFunction(ib_force_fcn);
    
```

- IB Method requires:
 - Integrating fluid
 - Integrating structural
 - Coupling operators
- Key principles:
 - Eulerian patch hierarchy
 - Integrator classes
 - IB model, forces, and initial position
 - Input file
- `IBStandardForceGen` has support for many standard elastic structures

$$\text{Spring (.spring): } \mathbf{F}_i = \kappa \frac{|\mathbf{X}_i - \mathbf{X}_j| - R_0}{|\mathbf{X}_i - \mathbf{X}_j|} (\mathbf{X}_i - \mathbf{X}_j)$$

$$\mathbf{F}_j = -\kappa \frac{|\mathbf{X}_i - \mathbf{X}_j| - R_0}{|\mathbf{X}_i - \mathbf{X}_j|} (\mathbf{X}_i - \mathbf{X}_j)$$

$$\text{Beams (.beam): } \mathbf{F} = K(\mathbf{X}_j + \mathbf{X}_k - 2\mathbf{X}_i) - \gamma$$

$$\mathbf{F}_i = 2\mathbf{F}, \quad \mathbf{F}_j = -\mathbf{F}, \quad \mathbf{F}_k = -\mathbf{F}$$

$$\text{Target (.target): } \mathbf{F}_i = \kappa(\mathbf{X}_0 - \mathbf{X}_i) - EU_i$$

- Note: Scaling for ΔX is not considered here. Discretization for spreading depends on dimension of structure (1/2/3D?)
- Final factor for ΔX should be provided in coefficients.

- IB Method requires:
 - Integrating fluid
 - Integrating structural
 - Coupling operators
 - Input file is parsed and provides runtime setup
 - Set physical parameters (viscosity, density)
- Key principles:
 - Eulerian patch hierarchy
 - Integrator classes
 - IB model, forces, and initial position
 - Input file

input2d

```

1 // constants
2 PI = 3.14159265358979
3
4 // physical parameters
5 L = 1.0
6 MU = 1.0e-2
7 RHO = 1.0
8 K = 1.0
9
10 // grid spacing parameters
11 MAX_LEVELS = 1 // maximum number of levels in locally refined grid
12 REF_RATIO = 4 // refinement ratio between levels
13 N = 64 // actual number of grid cells on coarsest grid level
14 NFINEST = (REF_RATIO^(MAX_LEVELS - 1))*N // effective number of grid cells on finest grid level
15 DX_FINEST = L/NFINEST
16
17 // solver parameters
18 DELTA_FUNCTION = "IB_4"
19 SOLVER_TYPE = "STAGGERED" // the fluid solver to use (STAGGERED or COLLOCATED)
20 START_TIME = 0.0e0 // initial simulation time
21 END_TIME = (3.0*log10(K))*0.55/K // final simulation time
22 GROW_DT = 2.0e0 // growth factor for timesteps
23 NUM_CYCLES = 1 // number of cycles of fixed-point iteration
24 CONVECTIVE_TS_TYPE = "ADAMS_BASHFORTH" // convective time stepping type
25 CONVECTIVE_OP_TYPE = "PPM" // convective differencing discretization type
26 CONVECTIVE_FORM = "ADVECTIVE" // how to compute the convective terms
27 NORMALIZE_PRESSURE = TRUE // whether to explicitly force the pressure to have mean zero
28 CFL_MAX = 0.3 // maximum CFL number
29 DT = (1.0/K)*1.6e-2*DX_FINEST // maximum timestep size
30 ERROR_ON_DT_CHANGE = TRUE // whether to emit an error message if the time step size changes
    
```

- IB Method requires:
 - Integrating fluid
 - Integrating structural
 - Coupling operators
- Each IB model has a different class, e.g. `IBMethod`, `IBFEMethod`, etc.
- Input file specifies structure name and level placement.

input2d

```

96 IBStandardInitializer {
97     max_levels      = MAX_LEVELS
98     structure_names = "curve2d_64"
99
100     beta  = 0.35
101     alpha = 0.25^2/beta
102
103     A = PI*alpha*beta // area of ellipse
104     R = sqrt(A/PI)    // radius of disc with equivalent area as the ellipse
105     perim = 2*PI*R    // perimeter of the equivalent disc
106
107     dx = L/NFINEST
108     dx_64 = L/64
109     num_node_circum = (dx_64/dx)*ceil(perim/(dx_64/3)/4)*4
110     ds = 2.0*PI*R/num_node_circum
111
112     curve2d_64 {
113         level_number = MAX_LEVELS - 1
114         uniform_spring_stiffness = K/ds
115     }
116 }
    
```

- IB Method requires:
 - Integrating fluid
 - Integrating structural
 - Coupling operators
- Each IB model has a different class, e.g. IBMethod, IBFEMethod, etc.
- IB structures can be setup as text files. Details on format can be found ibamr.github.io/docs: search for IBStandardInitializer

- Key principles:

- Eulerian patch hierarchy
- Integrator classes
- IB model, forces, and initial position
- Input file

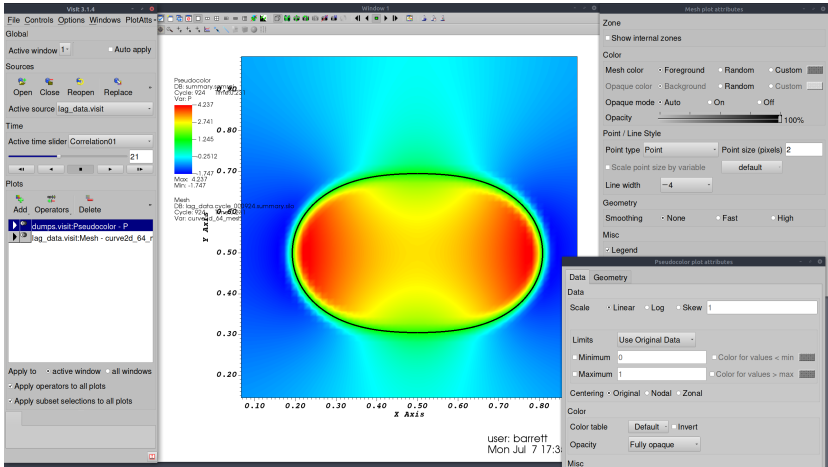
curve2d_64.vertex

```
304
6.7857142857142860e-01 5.0000000000000000e-01
6.7853328871213048e-01 5.0723341542964395e-01
6.7841888542630380e-01 5.1446374098708458e-01
6.7822826758319077e-01 5.2168788812000810e-01
6.7796151660833393e-01 5.2890277091531634e-01
6.7761874644879327e-01 5.3610530741732487e-01
6.7720010352447235e-01 5.4329242094427166e-01
6.7670576666557192e-01 5.5046104140257135e-01
6.7613594703620039e-01 5.5760810659825688e-01
6.7549088804417134e-01 5.6473056354504547e-01
6.7477086523702756e-01 5.7182536976847198e-01
6.7397618618433708e-01 5.7888949460553218e-01
6.7310719034630906e-01 5.8591992049927966e-01
6.7216424892878812e-01 5.9291364428782534e-01
6.7114776472468796e-01 5.9986767848718681e-01
6.7005817194193118e-01 6.0677905256744202e-01
6.6889593601797048e-01 6.1364481422163919e-01
6.6766155342096956e-01 6.2046203062692351e-01
6.6635555143772840e-01 6.2722778969734039e-01
```

curve2d_64.spring

```
304
0 1 1.9353241079974475e+02 0.0000000000000000e+00
1 2 1.9353241079974475e+02 0.0000000000000000e+00
2 3 1.9353241079974475e+02 0.0000000000000000e+00
3 4 1.9353241079974475e+02 0.0000000000000000e+00
4 5 1.9353241079974475e+02 0.0000000000000000e+00
5 6 1.9353241079974475e+02 0.0000000000000000e+00
6 7 1.9353241079974475e+02 0.0000000000000000e+00
7 8 1.9353241079974475e+02 0.0000000000000000e+00
8 9 1.9353241079974475e+02 0.0000000000000000e+00
9 10 1.9353241079974475e+02 0.0000000000000000e+00
10 11 1.9353241079974475e+02 0.0000000000000000e+00
11 12 1.9353241079974475e+02 0.0000000000000000e+00
12 13 1.9353241079974475e+02 0.0000000000000000e+00
13 14 1.9353241079974475e+02 0.0000000000000000e+00
14 15 1.9353241079974475e+02 0.0000000000000000e+00
15 16 1.9353241079974475e+02 0.0000000000000000e+00
16 17 1.9353241079974475e+02 0.0000000000000000e+00
17 18 1.9353241079974475e+02 0.0000000000000000e+00
18 19 1.9353241079974475e+02 0.0000000000000000e+00
```

- Use VisIt or Paraview to visualize data
- `dumps.visit` contains Eulerian data, `lag_data.visit` contains Lagrangian data



- To use AMR, increase MAX_LEVELS.
- GRADIENT_DETECTOR will refine based on IB location and vorticity values. Add REFINE_BOXES to specify portion of domain, see `navier_stokes/ex0`. Change MAX_LEVELS=2.
- Note: Make sure structure is refined as well: set structure to `curve2d_256`

```

10 // grid spacing parameters
11 MAX_LEVELS = 1 // maximum number of levels in locally refined grid
12 REF_RATIO = 4 // refinement ratio between levels
13 N = 64 // actual number of grid cells on coarsest grid level
14 NFINEST = (REF_RATIO^(MAX_LEVELS - 1))*N // effective number of grid cells on finest grid level
15 DX_FINEST = L/NFINEST

```

```

198 GriddingAlgorithm {
199     max_levels = MAX_LEVELS
200     ratio_to_coarser {
201         level_1 = REF_RATIO, REF_RATIO
202         level_2 = REF_RATIO, REF_RATIO
203         level_3 = REF_RATIO, REF_RATIO
204         level_4 = REF_RATIO, REF_RATIO
205         level_5 = REF_RATIO, REF_RATIO
206     }
207     largest_patch_size {
208         level_0 = 512, 512 // all finer levels will use same values as level_0
209     }
210     smallest_patch_size {
211         level_0 = 8, 8 // all finer levels will use same values as level_0
212     }
213     efficiency_tolerance = 0.85e0 // min % of tag cells in new patch level
214     combine_efficiency = 0.85e0 // chop box if sum of volumes of smaller boxes < efficiency * vol of large box
215 }
216
217 StandardTagAndInitialize {
218     tagging_method = "GRADIENT_DETECTOR"
219 }

```


- Example:

IB/explicit/ex0:

$$\mathbf{u}(\mathbf{x}, 0) = 0,$$

$$\chi(\theta, 0) = (X_0, Y_0) + R(\cos \theta, \sin \theta),$$

$$\mathbf{F}(\theta, 0) = \kappa \frac{\partial^2 \chi}{\partial \theta^2}$$

$$p(\mathbf{x}, 0) = \kappa H(\|\mathbf{x} - (X_0, Y_0)\|) + C$$

- Same as `ex1`, but exact initial conditions.
- Note: Despite exact initial conditions, spurious flows around structure

- Example:

IB/explicit/ex0:

$$\mathbf{u}(\mathbf{x}, 0) = 0,$$

$$\chi(\theta, 0) = (X_0, Y_0) + R(\cos \theta, \sin \theta),$$

$$\mathbf{F}(\theta, 0) = \kappa \frac{\partial^2 \chi}{\partial \theta^2}$$

$$p(\mathbf{x}, 0) = \kappa H(\|\mathbf{x} - (X_0, Y_0)\|) + C$$

- Same as `ex1`, but exact initial conditions.
- Note: Despite exact initial conditions, spurious flows around structure

- One problem: Discrete interpolated velocities are not divergence free.

$$\nabla_h \cdot \mathbf{u} = 0 \not\Rightarrow$$

$$\nabla_h \cdot \sum_{i,j} \mathbf{u}_{i,j} \delta_h(\mathbf{x}_{i,j} - \chi) h^2 = 0$$

- Example:

IB/explicit/ex0:

$$\mathbf{u}(\mathbf{x}, 0) = 0,$$

$$\chi(\theta, 0) = (X_0, Y_0) + R(\cos \theta, \sin \theta),$$

$$\mathbf{F}(\theta, 0) = \kappa \frac{\partial^2 \chi}{\partial \theta^2}$$

$$p(\mathbf{x}, 0) = \kappa H(\|\mathbf{x} - (X_0, Y_0)\|) + C$$

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$$\nabla_h \cdot \mathbf{u} = 0 \not\Rightarrow$$

$$\nabla_h \cdot \sum_{i,j} \mathbf{u}_{i,j} \delta_h(\mathbf{x}_{i,j} - \chi) h^2 = 0$$

- Another problem: $\mathbf{f}$ corresponds to a gradient (pressure)

$$\mathbf{f} = \int \mathbf{F}(\theta) \delta(\mathbf{x} - \chi) d\theta = \nabla \phi$$

but

$$\int \mathbf{F}(\theta) \delta_h(\mathbf{x} - \chi) d\theta \neq \nabla_h \phi$$

- One problem: Discrete interpolated velocities are not divergence free.

$$\nabla_h \cdot \mathbf{u} = 0 \not\Rightarrow$$

$$\nabla_h \cdot \sum_{i,j} \mathbf{u}_{i,j} \delta_h(\mathbf{x}_{i,j} - \chi) h^2 = 0$$

- Example:
IB/explicit/ex0:
 $\mathbf{u}(\mathbf{x}, 0) = 0,$
 $\chi(\theta, 0) = (X_0, Y_0) + R(\cos \theta, \sin \theta),$
 $\mathbf{F}(\theta, 0) = \kappa \frac{\partial^2 \chi}{\partial \theta^2}$
 $p(\mathbf{x}, 0) = \kappa H(\|\mathbf{x} - (X_0, Y_0)\|) + C$

- Another problem: $\mathbf{f}$ corresponds to a gradient (pressure)

$$\mathbf{f} = \int \mathbf{F}(\theta) \delta(\mathbf{x} - \chi) d\theta = \nabla \phi$$

but

$$\int \mathbf{F}(\theta) \delta_h(\mathbf{x} - \chi) d\theta \neq \nabla_h \phi$$

- Same as ex1, but exact initial conditions.
- Note: Despite exact initial conditions, spurious flows around structure

- Choice of delta function can dramatically affect results

- Example:

IB/explicit/ex0:

$$\mathbf{u}(\mathbf{x}, 0) = 0,$$

$$\chi(\theta, 0) = (X_0, Y_0) + R(\cos \theta, \sin \theta),$$

$$\mathbf{F}(\theta, 0) = \kappa \frac{\partial^2 \chi}{\partial \theta^2}$$

$$p(\mathbf{x}, 0) = \kappa H(\|\mathbf{x} - (X_0, Y_0)\|) + C$$

- Same as `ex1`, but exact initial conditions.
- Note: Despite exact initial conditions, spurious flows around structure

- One problem: Discrete interpolated velocities are not divergence free.

$$\nabla_h \cdot \mathbf{u} = 0 \not\Rightarrow$$

$$\nabla_h \cdot \sum_{i,j} \mathbf{u}_{i,j} \delta_h(\mathbf{x}_{i,j} - \chi) h^2 = 0$$

- Another problem: $\mathbf{f}$ corresponds to a gradient (pressure)

$$\mathbf{f} = \int \mathbf{F}(\theta) \delta(\mathbf{x} - \chi) d\theta = \nabla \phi$$

but

$$\int \mathbf{F}(\theta) \delta_h(\mathbf{x} - \chi) d\theta \neq \nabla_h \phi$$

- Change `IB_4` to
`COMPOSITE_BSPLINE_54`

- Track the centerline of the rod, denoted by χ .
- Introduce *director vectors* that account for twist of body.
- Use rod theories to describe both force and torque on thin bodies.

$$\rho \frac{D\mathbf{u}}{Dt} = \nabla \cdot \boldsymbol{\sigma} + \mathbf{f}(\mathbf{x}, t) + \frac{1}{2} \nabla \times \mathbf{n}(\mathbf{x}, t)$$

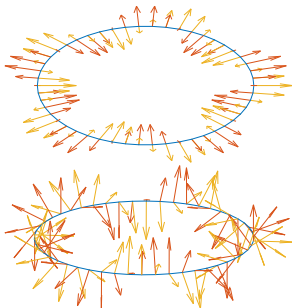
$$\nabla \cdot \mathbf{u}(\mathbf{x}, t) = 0$$

$$\mathbf{f}(\mathbf{x}, t) = \int_{\Omega} \mathbf{F}(s, t) \phi(\mathbf{x} - \chi(s, t)) ds$$

$$\mathbf{n}(\mathbf{x}, t) = \int_{\Omega} \mathbf{N}(s, t) \phi(\mathbf{x} - \chi(s, t)) ds$$

$$\frac{\partial \chi}{\partial t}(s, t) = \int_U \mathbf{u}(\mathbf{x}, t) \phi(\mathbf{x} - \chi(s, t)) d\mathbf{x}$$

$$\frac{\partial \mathbf{D}^i}{\partial t}(s, t) = \left(\frac{1}{2} \int_U \nabla \times \mathbf{u}(\mathbf{x}, t) \phi(\mathbf{x} - \chi(s, t)) d\mathbf{x} \right) \times \mathbf{D}^i(s, t)$$



- Track the centerline of the rod, denoted by χ .
- Introduce *director vectors* that account for twist of body.
- Use rod theories to describe both force and torque on thin bodies.

$\mathbf{F}(s, t)$ and $\mathbf{N}(s, t)$ come from a energy formulation:

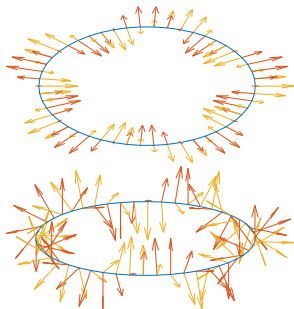
$$E_{\text{twist}} = \frac{1}{2} \int_0^L a_3 (\Omega_3 - \tau_1)^2 ds$$

$$E_{\text{bend}} = \frac{1}{2} \int_0^L \left(a_1 (\Omega_1 - \kappa_1)^2 + a_2 (\Omega_2 - \kappa_2)^2 \right) ds$$

$$E_{\text{shear}} = \frac{1}{2} \int_0^L \left(b_1 \left(\frac{\partial \chi}{\partial s} \cdot \mathbf{D}^1 \right)^2 + b_2 \left(\frac{\partial \chi}{\partial s} \cdot \mathbf{D}^2 \right)^2 \right) ds$$

$$E_{\text{stretch}} = \frac{1}{2} \int_0^L b_3 \left(\frac{\partial \chi}{\partial s} \cdot \mathbf{D}^3 - 1 \right)^2 ds$$

$$\text{with } \Omega_1 = \frac{\partial \mathbf{D}^2}{\partial s} \cdot \mathbf{D}^3, \Omega_2 = \frac{\partial \mathbf{D}^3}{\partial s} \cdot \mathbf{D}^1, \Omega_3 = \frac{\partial \mathbf{D}^1}{\partial s} \cdot \mathbf{D}^2.$$



- Example: IB/explicit/ex4: twisted elastic ring
- Register a `GeneralizedIBMethod` object with the integrator.
- Physical parameters specified in input files.

```
161     Pointer<GeneralizedIBMethod> ib_method_ops = new GeneralizedIBMethod(  
162         "GeneralizedIBMethod", app_initializer->getComponentDatabase("GeneralizedIBMethod"));  
163     Pointer<IBHierarchyIntegrator> time_integrator =  
164         new IBExplicitHierarchyIntegrator("IBHierarchyIntegrator",  
165             app_initializer->getComponentDatabase("IBHierarchyIntegrator"),  
166             ib_method_ops,  
167             navier_stokes_integrator);  
185     // Configure the IB solver.  
186     Pointer<IBStandardInitializer> ib_initializer = new IBStandardInitializer(  
187         "IBStandardInitializer", app_initializer->getComponentDatabase("IBStandardInitializer"));  
188     ib_method_ops->registerInitStrategy(ib_initializer);  
189     Pointer<IBKirchhoffRodForceGen> ib_force_and_torque_fcn = new IBKirchhoffRodForceGen();  
190     ib_method_ops->registerIBKirchhoffRodForceGen(ib_force_and_torque_fcn);
```


- Example: IB/explicit/ex4: twisted elastic ring
- Register a GeneralizedIBMethod object with the integrator.
- Physical parameters specified in input files.

```
64 for r = 0:nr-1
65     theta = 2*pi*r/nr;
66
67     R = [cos(theta) , sin(theta) , 0];
68     Theta = [-sin(theta), cos(theta) , 0];
69     Z = [0 , 0 , 1];
70
71     X = r1*R;
72     vertices(r+1,:) = X;
73
74     fprintf(vertex_fid, '%1.16e %1.16e %1.16e\n', X(1), X(2), X(3));
75
76     r_curr = r;
77     r_next = mod(r+1,nr);
78
79     fprintf(rod_fid, ['%6d %6d %1.16e %1.16e %1.16e %1.16e %1.16e %1.16e ' ...
80                     '%1.16e %1.16e %1.16e %1.16e\n'], r_curr, r_next, ds, ...
81             a1, a2, a3, b1, b2, b3, kappa1, kappa2, tau);
82
83     D3 = cos(beta)*Theta + sin(beta)*Z;
84     E = -sin(beta)*Theta + cos(beta)*Z;
85     D1 = cos(p*theta + epsilon*sin(theta))*E + sin(p*theta + epsilon*sin(theta))*R;
86     D2 = -sin(p*theta + epsilon*sin(theta))*E + cos(p*theta + epsilon*sin(theta))*R;
87
88     D1_vals(r+1,:) = D1;
89     D2_vals(r+1,:) = D2;
90     D3_vals(r+1,:) = D3;
91
92     fprintf(director_fid, '%1.16e %1.16e %1.16e\n', D1(1), D1(2), D1(3));
93     fprintf(director_fid, '%1.16e %1.16e %1.16e\n', D2(1), D2(2), D2(3));
94     fprintf(director_fid, '%1.16e %1.16e %1.16e\n', D3(1), D3(2), D3(3));
95 end %for
```

- Viscoelastic fluids exhibit both *elastic* and *viscous* responses to forces.

Viscoelastic Fluid Models

- Viscoelastic fluids exhibit both *elastic* and *viscous* responses to forces.

Springs:

$$\sigma = G\gamma$$

Dashpot:

$$\sigma = \mu\dot{\gamma}$$

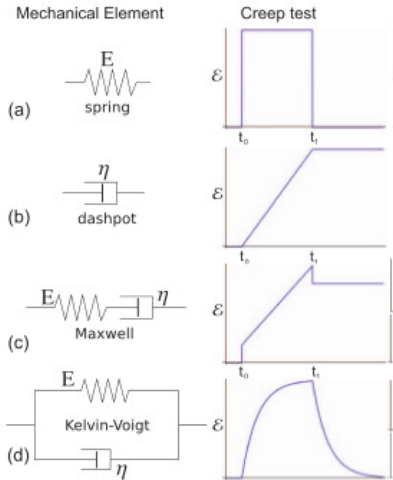
Maxwell Element:

$$\dot{\sigma} = \frac{\mu}{\lambda}\dot{\gamma} - \frac{1}{\lambda}\sigma$$

$$\lambda = \frac{\mu}{G}$$

Kelvin-Voigt:

$$\sigma = G\gamma + \mu\dot{\gamma}$$



- Maxwell model:

$$\lambda \overset{\nabla}{\tau}_p = \eta \dot{\gamma} - \tau_p$$

- Oldroyd type models: let $\sigma = -p\mathbb{I} + \mu_s \dot{\gamma} + \tau_p$
 - Oldroyd-B model:

$$\tau_p + \lambda \overset{\nabla}{\tau}_p = \mu \dot{\gamma}$$

- Upper-convected derivative:

$$\overset{\nabla}{\tau} = \lambda \left(\frac{\partial \tau}{\partial t} + \mathbf{u} \cdot \nabla \mathbf{u} - \nabla \mathbf{u} \cdot \tau - \tau \cdot \nabla \mathbf{u}^T \right)$$

- Maxwell model:

$$\lambda \overset{\nabla}{\tau}_p = \eta \dot{\mathbb{C}} - \tau_p$$

- Oldroyd type models: let $\sigma = -p\mathbb{I} + \mu_s \dot{\mathbb{C}} + \tau_p$
 - Oldroyd-B model:

$$\tau_p + \lambda \overset{\nabla}{\tau}_p = \mu \dot{\mathbb{C}}$$

- Upper-convected derivative:

$$\overset{\nabla}{\tau} = \lambda \left(\frac{\partial \tau}{\partial t} + \mathbf{u} \cdot \nabla \mathbf{u} - \nabla \mathbf{u} \cdot \tau - \tau \cdot \nabla \mathbf{u}^T \right)$$

- Giesekus model:

$$\tau_p + \lambda \overset{\nabla}{\tau}_p + \alpha \frac{\lambda}{\mu_p} \tau_p \cdot \tau_p = \mu \dot{\mathbb{C}}$$

- Here, $\tau_p \cdot \tau_p$ is an *anisotropic drag term*

- Complete Model:

$$\rho \left(\frac{\partial \mathbf{u}}{\partial t} + \mathbf{u} \cdot \nabla \mathbf{u} \right) = \nabla \cdot \boldsymbol{\sigma} + \mathbf{f}$$

$$\nabla \cdot \mathbf{u} = 0$$

$$\boldsymbol{\sigma} = -p\mathbb{I} + \mu_s \dot{\boldsymbol{\gamma}} + \boldsymbol{\tau}$$

$$\lambda \dot{\boldsymbol{\tau}} = \mu_p \dot{\boldsymbol{\gamma}} - \boldsymbol{\tau}$$

- Relevant dimensionless numbers:

$$\text{Re} = \frac{\rho UL}{\mu}, \quad \text{Wi} = \dot{\gamma} \lambda, \quad \text{De} = \frac{\lambda}{t_p}$$

- Example: `examples/complex_fluids/ex2`, flow past a confined cylinder.
- Note: `set OUTPUT_STRAIN = FALSE`.

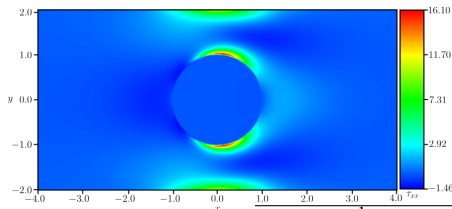
- Example: `examples/complex_fluids/ex2`
- Advect a stress tensor, apply an extra force to momentum

```
243     Pointer<INSHierarchyIntegrator> navier_stokes_integrator;  
244     const string solver_type = app_initializer->getComponentDatabase("Main")->getString("solver_type");  
245     if (solver_type == "STAGGERED")  
246     {  
247         navier_stokes_integrator = new INStaggeredHierarchyIntegrator(  
248             "INStaggeredHierarchyIntegrator",  
249             app_initializer->getComponentDatabase("INStaggeredHierarchyIntegrator"));  
250     }  
251     else if (solver_type == "COLLOCATED")  
252     {  
253         navier_stokes_integrator = new INSCollocatedHierarchyIntegrator(  
254             "INSCollocatedHierarchyIntegrator",  
255             app_initializer->getComponentDatabase("INSCollocatedHierarchyIntegrator"));  
256     }  
257     else  
258     {  
259         TBOX_ERROR("Unsupported solver type: " << solver_type << "\n"  
260                     << "Valid options are: COLLOCATED, STAGGERED");  
261     }  
262  
263     // Create the advection diffusion integrator for the extra stress.  
264     Pointer<AdvDiffSemiImplicitHierarchyIntegrator> adv_diff_integrator;  
265     adv_diff_integrator = new AdvDiffSemiImplicitHierarchyIntegrator(  
266         "AdvDiffSemiImplicitHierarchyIntegrator",  
267         app_initializer->getComponentDatabase("AdvDiffSemiImplicitHierarchyIntegrator"));  
268     navier_stokes_integrator->registerAdvDiffHierarchyIntegrator(adv_diff_integrator);  
269  
270     // Generate the extra stress component and set up necessary components. The constructor registers the extra  
271     // stress with the advection diffusion integrator.  
272     Pointer<CFINSForcing> polymericStressForcing;  
273     double viscosity = 0.0, relaxation_time = 0.0;  
274     if (input_db->keyExists("ComplexFluid"))  
275     {  
276         polymericStressForcing = new CFINSForcing("PolymericStressForcing",  
277             app_initializer->getComponentDatabase("ComplexFluid"),  
278             navier_stokes_integrator,  
279             grid_geometry,  
280             adv_diff_integrator,  
281             visit_data_writer);  
282         time_integrator->registerBodyForceFunction(polymericStressForcing);  
283         viscosity = input_db->getDatabase("ComplexFluid")->getDouble("viscosity");  
284         relaxation_time = input_db->getDatabase("ComplexFluid")->getDouble("relaxation_time");  
285     }  
286 }
```

- Example: `examples/complex_fluids/ex2`
- *Impose* no motion by applying force $\mathbf{F} = \kappa(\chi^0 - \chi) - \eta \frac{\partial \chi}{\partial t}$
- As $\kappa \rightarrow \infty$, we must have that $\chi \rightarrow \chi^0$.

```
64 void
65 tether_force_function(VectorValue<double>& F,
66                      const libMesh::VectorValue<double>& /*n*/,
67                      const libMesh::VectorValue<double>& /*N*/,
68                      const TensorValue<double>& /*FF*/,
69                      const libMesh::Point& x,
70                      const libMesh::Point& X,
71                      Elem* const /*elem*/,
72                      unsigned short int /*side*/,
73                      const vector<const vector<double>>>& var_data,
74                      const vector<const vector<VectorValue<double>>>& /*grad_var_data*/,
75                      double /*time*/,
76                      void* /*ctx*/)
77 {
78     const std::vector<double>& U = *var_data[0];
79     for (unsigned int d = 0; d < NDIM; ++d)
80     {
81         F(d) = kappa_s * (X(d) - x(d)) - eta_s * U[d];
82     }
83     std::vector<double> d(NDIM);
84     d[0] = std::abs(x(0) - X(0));
85     d[1] = std::abs(X(1) - x(1));
86     // Quit the simulation if the structure has moved more than a quarter of the grid spacing.
87     if (ERROR_ON_MOVE && ((d[0] > 0.25 * dx) || (d[1] > 0.25 * dx)))
88     {
89         TBOX_ERROR("Structure has moved too much.\n");
90     }
91
92     // Configure the IBFE solver.
93     ib_method_ops->initializeFEEquationSystems();
94     std::vector<int> vars(NDIM);
95     for (unsigned int d = 0; d < NDIM; ++d) vars[d] = d;
96     vector<SystemData> sys_data(1, SystemData(IBFESurfaceMethod::VELOCITY_SYSTEM_NAME, vars));
97     IBFESurfaceMethod::LagSurfaceForceFcnData body_fcn_data(tether_force_function, sys_data);
98     ib_method_ops->registerLagSurfaceForceFunction(body_fcn_data);
```


Immersed Boundary Method



	L_1 rate	L_2 rate	L_∞ rate
$\mathbf{u}$	1.693	1.470	1.020
τ_{xx}	1.173	0.537	-0.093
τ_{xy}	1.341	0.632	0.130
τ_{yy}	1.254	0.872	0.017
p	1.275	1.204	-0.541

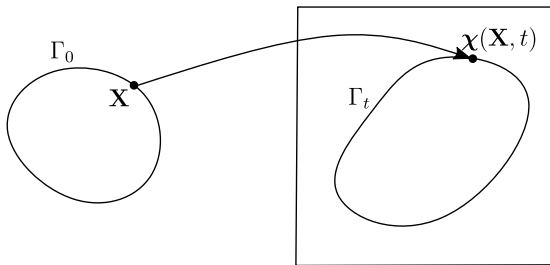
- Drag coefficient

$$C_d = \int_{\Gamma} \sigma \cdot \mathbf{n} \cdot \mathbf{e}_x ds$$

- IB stresses fail to converge pointwise!

$M_{\text{fac}} = \frac{1}{2}$	BS3	IB3	PL	IB4	IB6
Coarse	6.15%	9.55%	12.92%	10.50%	11.16%
Medium	2.78%	4.11%	5.39%	4.73%	5.02%
Fine	1.39%	2.04%	2.71%	2.32%	2.38%
$M_{\text{fac}} = 1$					
Coarse	5.45%	6.20%	4.77%	9.51%	11.16%
Medium	2.61%	2.92%	2.54%	4.37%	5.02%
Fine	1.24%	1.42%	1.19%	2.07%	2.38%
$M_{\text{fac}} = 2$					
Coarse	4.57%	5.86%	1.57%	9.06%	11.16%
Medium	2.12%	2.66%	1.03%	4.16%	5.02%
Fine	1.01%	1.28%	0.46%	1.99%	2.38%

Immersed Boundary Method

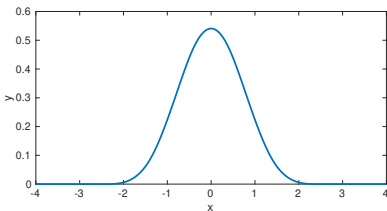


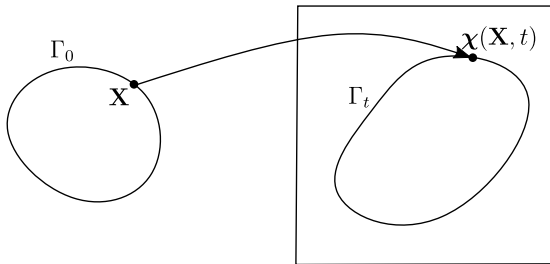
$$\rho \frac{D\mathbf{u}(\mathbf{x}, t)}{Dt} = \nabla \cdot \boldsymbol{\sigma}(\mathbf{x}, t) + \mathbf{f}(\mathbf{x}, t)$$

$$\nabla \cdot \mathbf{u}(\mathbf{x}, t) = 0$$

$$\mathbf{f}(\mathbf{x}, t) = \int_{\Gamma_0} \mathbf{F}(\mathbf{X}, t) \delta(\mathbf{x} - \chi(\mathbf{X}, t)) d\mathbf{X}$$

$$\frac{\partial \chi(\mathbf{X}, t)}{\partial t} = \int_{\Omega} \mathbf{u}(\mathbf{x}, t) \delta(\mathbf{x} - \chi(\mathbf{X}, t)) d\mathbf{x}$$





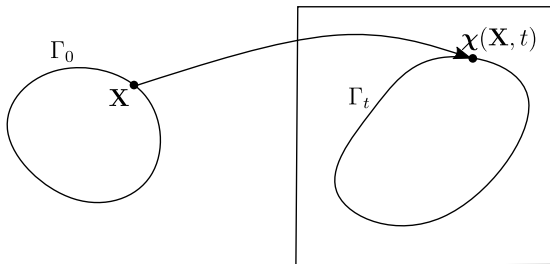
$$\rho \frac{D\mathbf{u}(\mathbf{x}, t)}{Dt} = \nabla \cdot \boldsymbol{\sigma}(\mathbf{x}, t) + \mathbf{f}(\mathbf{x}, t)$$

$$\nabla \cdot \mathbf{u}(\mathbf{x}, t) = 0$$

$$[[\boldsymbol{\sigma}]] \cdot \mathbf{n} = -\mathbf{F}\mathbf{J}^{-1}$$

$$\frac{\partial \boldsymbol{\chi}(\mathbf{X}, t)}{\partial t} = \int_{\Omega} \mathbf{u}(\mathbf{x}, t) \delta(\mathbf{x} - \boldsymbol{\chi}(\mathbf{X}, t)) d\mathbf{x}$$

- Singular delta function induces jumps in the stress.
- Regular finite differences have $\mathcal{O}(1)$ errors for discontinuous variables
- Immersed Interface: replace δ_h with corrected differences.



- If $\phi(x)$ has a discontinuity at $\hat{x} \in [x_{i+\frac{1}{2}}, x_{i+1}]$, then

$$\frac{\partial \phi}{\partial x} \left(x_{i+\frac{1}{2}} \right) = \frac{\phi_{i+1} - \phi_i}{h} - \frac{1}{h} \sum_{m=0}^2 \frac{(x_{i+1} - \hat{x})^m}{m!} \left[\left[\frac{\partial^m \phi}{\partial x^m} \right] (\hat{x}) + \mathcal{O}(h^2) \right]$$

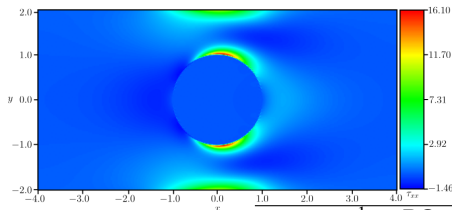
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$$\nabla \cdot \mathbf{u}(\mathbf{x}, t) = 0$$

$$[[\boldsymbol{\sigma}]] \cdot \mathbf{n} = -\mathbf{F}\mathbf{J}^{-1}$$

$$\frac{\partial \boldsymbol{\chi}(\mathbf{X}, t)}{\partial t} = \int_{\Omega} \mathbf{u}(\mathbf{x}, t) \delta(\mathbf{x} - \boldsymbol{\chi}(\mathbf{X}, t)) d\mathbf{x}$$

Immersed Boundary Method



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- Drag coefficient

$$C_d = \int_{\Gamma} \boldsymbol{\sigma} \cdot \mathbf{n} \cdot \mathbf{e}_x ds$$

- IB stresses fail to converge pointwise!
- Averaged quantities are accurate.

$M_{\text{fac}} = \frac{1}{2}$	BS3	IB3	PL	IB4	IB6
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Fine	1.01%	1.28%	0.46%	1.99%	2.38%

- Derived conservation of mass and momentum equations
- Developed the IB equations:
 - Lagrangian structure moves at local fluid velocity
 - Deformations induce elastic forces which are applied to momentum equation
 - Fluid viscosity exists throughout entire domain.
- IBAMR provides a framework to implement different IB strategies strategies

