

Notes on SPH

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Part I

SPH FOR CONTINUOUM DOMAIN

1. NAVIER-STOKES EQUATIONS IN LAGRANGIAN FORM

The Navier-Stokes equations in the Lagrangian description are[1]

- Continuity equation

$$\frac{D\rho}{Dt} = -\rho \nabla \cdot v$$

This equation is derived from mass conservation inside a infinitesimal control volume δV in Lagrangian description(the control volume is moving in a streamline). This expression assumes that mass is conserved inside the control volume and velocity does not change across the control volume. In the equation, velocity divergence is interpreted as the time rate of volume change per unit volume.

- Momentum en x direction

$$\rho \frac{dv_x}{dt} = -\frac{\partial P}{\partial x} + \frac{\partial \tau_{xx}}{\partial x} + \frac{\partial \tau_{yx}}{\partial y} + \frac{\partial \tau_{zx}}{\partial z} + \rho F_x$$

This equation is derived from Newton second Law for Lagrangian control volume assuming constant mass inside the volume. The forces in consideration here are hydrostatic Pressure P , body forces per volume unit F_x and the forces generated by the stress state of the control volume.

- Energy equation

$$\begin{aligned} \rho \frac{De}{Dt} = & -P \left(\frac{\partial v_x}{\partial x} + \frac{\partial v_y}{\partial y} + \frac{\partial v_z}{\partial z} \right) + \\ & \tau_{xx} \frac{\partial v_x}{\partial x} + \tau_{yx} \frac{\partial v_x}{\partial y} + \tau_{zx} \frac{\partial v_x}{\partial z} \\ & \tau_{xy} \frac{\partial v_y}{\partial x} + \tau_{yy} \frac{\partial v_y}{\partial y} + \tau_{zy} \frac{\partial v_y}{\partial z} \\ & \tau_{xz} \frac{\partial v_z}{\partial x} + \tau_{yz} \frac{\partial v_z}{\partial y} + \tau_{zz} \frac{\partial v_z}{\partial z} \end{aligned}$$

This equation takes in account the effect of work done by isotropic pressure and the energy dissipation of viscous shear forces

Those equations can be expressed in a more compact way. Superscripts denote coordinate directions. The summation of the equations is taken over repeated indices and the total time derivatives are taken in the moving Lagrangian frame.

- Continuity

$$\frac{D\rho}{Dt} = -\rho \frac{\partial v^\beta}{\partial x^\beta}$$

- Momentum

$$\frac{Dv^\alpha}{Dt} = \frac{1}{\rho} \frac{\partial \sigma^{\alpha\beta}}{\partial x^\beta}$$

- Energy

$$\frac{De}{Dt} = \frac{1}{\rho} \sigma^{\alpha\beta} \frac{\partial v^\alpha}{\partial x^\beta}$$

Those equations use the stress tensor $\sigma^{\alpha\beta}$ as a combination of Pressure P and the stress supported by the control volume τ_{ij} .

$$\sigma^{\alpha\beta} = -P\delta^{\alpha\beta} + \tau^{\alpha\beta}$$

In the case of fluid dynamics, the deviatoric stress is given by:

$$\tau^{\alpha\beta} = \mu \epsilon^{\alpha\beta}$$

but, in the case of solid mechanics, the deviatoric stress is proportional to the strain rate tensor[2][3]

There is a little difference with the equation proposed by [4]. The equation on [3] is difficult to see because of the scanning conditions

$$\dot{\tau}^{\alpha\beta} = \mu \epsilon^{\bar{\alpha}\beta} = \mu \left(\dot{\epsilon}^{\alpha\beta} - \frac{1}{3} \delta^{\alpha\beta} \dot{\epsilon}^{\gamma\gamma} \right)$$

where μ is the shear modulus and ϵ is the traceless rate of strain. Rotation terms are needed to allow the transformation of the stress from the reference frame associated with the material to the laboratory reference frame with all other equations are specified[3][4]. The Jaumann rate is the most widely used in codes and we adopt it also. With Jaumann rate our constitutive equation is

$$\dot{\tau}^{\alpha\beta} = \mu \left(\dot{\epsilon}^{\alpha\beta} - \frac{1}{3} \delta^{\alpha\beta} \dot{\epsilon}^{\gamma\gamma} \right) + \tau^{\alpha\gamma} R^{\beta\gamma} + \tau^{\beta\gamma} R^{\alpha\gamma}$$

There are some misunderstandings about the superindexes of R terms) The strain rate and rotation rate that have been used are defined as follows[3][4]

$$\dot{\epsilon}^{\alpha\beta} = \frac{1}{2} \left(\frac{\partial v^\alpha}{\partial x^\beta} + \frac{\partial v^\beta}{\partial x^\alpha} \right)$$

$$R^{\alpha\beta} = \frac{1}{2} \left(\frac{\partial v^\alpha}{\partial x^\beta} - \frac{\partial v^\beta}{\partial x^\alpha} \right)$$

The calculations for each part of σ depend on the equation of state (for pressure) and the material model (for stress). The 'plastic behaviour' can be introduced in the equations using the von Mises yielding criterion. We limit our deviatoric stress tensor by[4]

$$S^{\alpha\beta} \Rightarrow f S^{\alpha\beta}$$

where f is computed from

$$f = \min \left(\frac{Y_o^2}{3J_2}, 1 \right)$$

and J_2 is the second invariant of the deviatoric stress tensor defined as

$$J_2 = \frac{1}{2} S^{\alpha\beta} S^{\alpha\beta}$$

In reference [3] the model is similar but there are some little differences ***I think this could be an important topic to achieve my objectives in the SPH project***

2. MATERIAL MODEL - FRACTURE

The fracture model corresponds to the Graddy & Kipp Fragmentation Model.[4]

The model begins by assuming the existence of incipient flaws in brittle solids. For a given failure strain ϵ , the probability distribution

$$n(\epsilon) = k\epsilon^n$$

defines the most number of flaws per unit volume having failure strain lower than ϵ . This Weibull distribution is isotropic, including no information about the location or orientation of any given flaw. The Weibull parameters m and k are material constants. A flaw becomes active when the local tensile strain reaches the flaw's activation threshold. Then, a crack is allowed to grow at constant velocity[4] $c_g = 0.4c_{\text{longitudinal elastic wave}}$. For a crack that is activated at time t' , the half-length at time t is

$$a = c_g(t - t')$$

The growing of a local crack implies that the surrounding material is losing its ability to support stress. This phenomenon is taken into account by introducing a state variable $D\epsilon$ [01] for 'damage'. Damage expresses the reduction in stress under tensile loading[4]

$$\sigma_D = \sigma(1 - D)$$

where σ is the elastic stress in the absence of damage and σ_D is the damaged-relieved stress. A fully damaged material is a cohesionless fluid, feeling no tensile or or shear stress whatsoever. When a particle is fully damaged, it can not support any stress, it is that the material breaks at this particle. To generate a fracture, the system should provide the energy required to deform the particles and then generate the material breaking. I suppose that when the material breaks, the accumulated elastic energy that could not be accumulated in elastic form should convert (partially) into kinetic energy. A crack relieves stress in a volume approximately equal to its circumscribing sphere. Thus

we define D locally as the fractional volume that is relieved of stress by local growing cracks

$$D = \frac{\frac{4}{3}\pi a^3}{V}$$

where $V = \frac{4}{3}\pi R_s^3$ is a subvolume of radius R_s where the crack is growing. I made a modification here. Original text[4] says '...is a volume in which a crack of half-length R_s is growing'. I think this is a writing mistake because in the next paragraph it refers to R_s as a 'a small subvolume of radius R_s '. Using the equation for the half-length of the crack a_g , we can obtain the relation from the growing of the damage D :

$$\frac{dD^{1/3}}{dt} = \frac{C_g}{R_s}$$

This equation expresses the evolution of D in a small subvolume of radius R_s of material subject to tensile strains. The misunderstanding here is: how to define the radius R_s of a subvolume? If I choose the subvolume to be a big portion of the material, D will be small for most of the time. But, If I choose a subvolume to be a very little portion of the material, the damage will be high since the very first steps of the simulation. The overall damage experienced by the material under consideration is given by the distribution of D over the entire object. In an elemental subvolume V , damage is independent of strain once the activation threshold has been exceeded. (It is because of D is defined as a volume relation and the length of the growing crack a_g only depends on the time transpired since the flaw became active) However, over a larger volume it is not, since the number of active flaws are strain dependent. Then, the damage is in fact strain dependent.

Part II

MATHEMATICAL REVIEW

3. CALCULUS

Vector: The term **vector** is used by scientist to indicate a quantity (such as displacement or velocity or force) that has both **magnitude** and **direction**[5]

3.1 Dot Product

Definition If $\vec{a} = \langle a_1, a_2, a_3 \rangle$ and $\vec{b} = \langle b_1, b_2, b_3 \rangle$, then the dot product of \vec{a} and \vec{b} is the number $\vec{a} \cdot \vec{b}$ given by

$$\vec{a} \cdot \vec{b} = a_1b_1 + a_2b_2 + a_3b_3$$

If θ is the angle between the vector \vec{a} and \vec{b} then

$$\vec{a} \cdot \vec{b} = |\vec{a}||\vec{b}|\cos\theta$$

Note that if \vec{a} and \vec{b} are parallel vectors, then $\theta = [0, \pi]$ and $\cos\theta = 1$.

Two vector \vec{a} and \vec{b} are orthogonal if and only if $\vec{a} \cdot \vec{b} = 0$, which means that $\theta = \pi/2$.

3.2 Divergence

If $F = P\hat{i} + Q\hat{j} + R\hat{k}$ is a vector Field in R^3 , then the **divergence of F** is the function of three variables defined by (assuming that the partial derivatives exists):

$$\text{div}(F) = \nabla \cdot F = \frac{\partial P}{\partial x} + \frac{\partial Q}{\partial y} + \frac{\partial R}{\partial z}$$

observe that $\nabla \cdot F$ is a scalar field.

If F is a vector field on R^3 , then $\text{curl}(F)$ is also a vector field on R^3 . As such, we can compute its divergence:

$$\text{div curl}(F) = 0$$

3.3 Divergence Theorem

Let E be a simple solid region and let S be the boundary surface of E , given with positive (outward) orientation. Let F be a vector field whose component functions have continuous partial derivatives on an open region that contains E . Then

$$\iint_S F \cdot dS = \iiint_E \nabla \cdot F dV$$

The divergence theorem states that, under the given conditions, the flux of F across the boundary surface of E is equal to the volume integral of the divergence of F over E i.e the theorem let us change a surface integral to a volume integral.

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