

# Implementation of Crack and Material Velocity Fields with Contact and Nodal Velocity Boundary Conditions

John Nairn

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# Chapter 1

## Velocity Fields

Each node has one to four crack velocity fields denoted [0], [1], [2], and [3]. If the problem has no cracks, then each node has just the single crack velocity field [0]. Each crack velocity field has any number of material velocity fields or just one velocity field if using single material mode.

### 1.1 Crack Velocity Fields

The crack velocity field is determined by drawing a line from the particle to the node. The crack velocity field number  $n$  is denoted by square brackets —  $[n]$ . If the line crosses no cracks, that material point uses field [0]. If the line crosses one crack, that material point uses field [1] (for the first crack found) or field [2] (if a second crack is found for the same node). If the line crosses two cracks, the material point uses field [3]. This scheme can handle one or two cracks at each node. A third crack (*i.e.*, a line that crosses a third crack or one of the cracks in field [3] being different than the cracks for fields [1] and [2]) is handled by picking one of the four fields, but is likely to not geting the physics correct. A warning is issued if a third crack is found.

Figure 1.1 shows schematic view of crack velocity fields for the indicated nodal point. Because field [2] is never allocated before field [1] (*i.e.*, the first single crack crossing is put into field [1] and the second into field [2]), the following crack field situations are all that are possible: [0], [1], [3], [0]&[1], [0]&[3], [1]&[2], [1]&[3], [0]&[1]&[2], [0]&[1]&[3], [1]&[2]&[3], and [0]&[1]&[2]&[3]. The combinations that never occur are the remaining ones or [2], [0]&[2], [2]&[3], and [0]&[2]&[3].

When doing crack contact, the possible field contacts are between [0]&[1] and [2]&[3] on the surface of the first crack (*i.e.*, crack in [1] and [3]) and between [0]&[2] and [1]&[3] on the surface of the second crack (*i.e.*, crack in [2] and [3]). Fields [0]&[3] and [1]&[2] do not contact. These can be visualized by looking at Figure 1.1 where non-contacting fields meet at a point instead of a surface.

### 1.2 Material Velocity Fields

When running in multimaterial mode (which is needed to model material contact), each crack velocity field may have 1 to  $n$  material velocity fields where  $n$  is the number of materials active in the simulations. When not using multimaterial mode,  $n$  is set to 1 and all materials use the same velocity field in each crack velocity field. All these fields are on the node. In these notes, we denote a quantity,  $q$ , in a material velocity field (which is where momentum, forces, *etc.* are stored) on the grid as

$$q_{i,c;m}^{(k)} \tag{1.1}$$

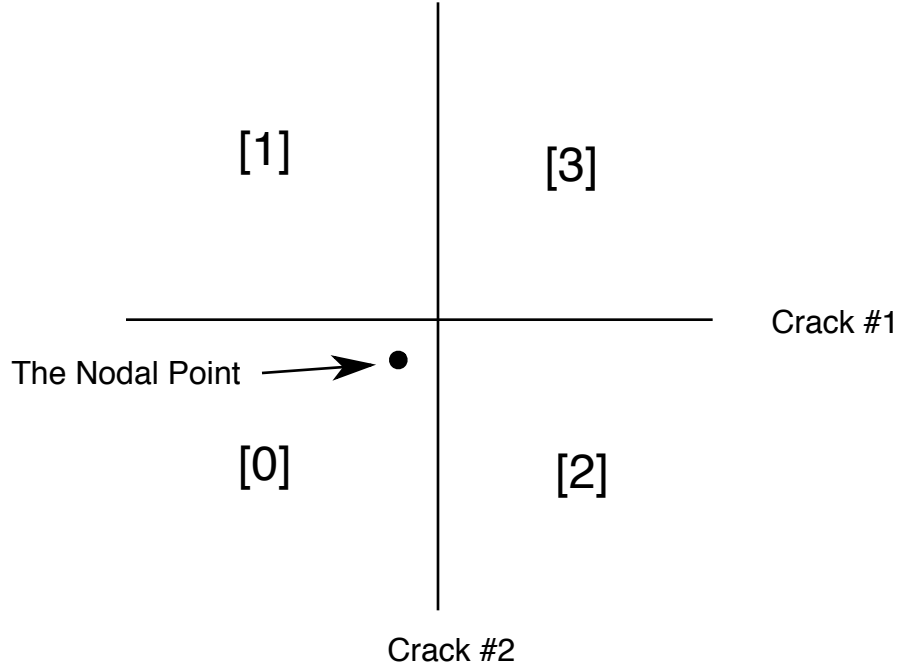


Figure 1.1: Schematic view of crack velocity fields around a single node. The possibility of two interacting cracks leads to at most four velocity fields.

where  $i$  is node number,  $c$  is crack velocity field,  $m$  is material velocity field, and  $k$  indicates time step number. When no cracks are present,  $c$  can be omitted; when in single material model,  $m$  can be omitted, when no cracks and single material mode both  $c$  and  $m$  can be omitted (*i.e.*, each node has only one material velocity field). When analyzing all values within a single time step,  $(k)$  can be omitted for simplicity.

## Chapter 2

# MPM Time Step Tasks

This section lists each task in NairnMPM outlining only those parts relevant to crack and material velocity fields, to contact conditions, to imperfect interfaces, to traction laws on cracks, and to nodal velocity boundary conditions.

### 2.1 Simulation Initialization

Each nodal point is permanently assigned crack field [0]. Other crack fields are dynamically allocated as needed. When in single material velocity field, field [0] is given a single material velocity field (field [0:0] for [crack field number:material field number]). When using multimaterial mode, crack velocity field [0] is assigned pointers to material velocity fields 0 to  $m - 1$  where  $m$  is the number of active materials in the simulations (each index maps to material number defined in the input file). The pointers are initialized to NULL. Any material velocity fields needed in a simulation are dynamically allocated as needed.

### 2.2 Time Step Initialization

All currently allocated crack and material velocity fields are zeroed. Then each material point is examined and compared to all nodes it sees with non-zero shape functions. When cracks are present, a line is traced from the particle to the node and compared to crack surfaces. The end result is determination of the particle's crack velocity field number `vfld` from [0] to [3]. If the node does not have that crack velocity field, it is not allocated on that node. When no cracks are present `vfld` is set to 0. The crack field number, `vfld`, is stored on the particle. This crack field number determination is likely the most time-consuming part of explicit crack calculations.

Note that first crack found on a node by itself goes into field [1] and second goes into field [2], which means never have field [2], without field [1]. No counting a node with zero fields, there are 15 combinations of velocity fields possible on a node. Of these, 5 have no contact, 6 may have contact, and 4 never occur:

- No Contact: [0], [1], [3], [0]&[3], [1]&[2]
- May Have Contact: [0]&[1], [1]&[3], [0]&[1]&[2], [0]&[1]&[3], [1]&[2]&[3], and [0]&[1]&[2]&[3].
- Never Occur: [2], [0]&[2], [2]&[3], [0]&[2]&[3]

A table look up converts material number (from input file) to index `matfld` (or field ID) into the `mvf[]` array of pointers on each crack velocity field. If that pointer `mvf[matfld]` is NULL, a new material velocity field is created.

## 2.3 Momentum and Mass Extrapolation to the Grid

For each material point, determine the appropriate material field (`matfld`) from the material assigned to that material point. Add the momentum and mass to material velocity field `matfld` of crack velocity field `vfld` (where `vfld` is stored on the particle). To support crack and material contact, each node needs to extrapolate particle position or displacement, particle size, and particle gradient (the last two are for normal detection and gradient only when in multimaterial mode). The relevant quantities for material velocity field  $j = \text{matfld}$  in crack velocity field  $c = \text{vfld}$  for particle  $p$  are:

$$m_{i,c;j} = \sum_{p \in j} S_{ip} m_p \quad (2.1)$$

$$\mathbf{p}_{i,c;j} = m_{i,c;j} \mathbf{v}_{i,c;j} = \sum_{p \in j} S_{ip} \mathbf{p}_p \quad (2.2)$$

$$m_{i,c;j} \mathbf{x}_{i,c;j} = \begin{cases} \sum_{p \in j} S_{ip} m_p \mathbf{x}_p & \text{position extrapolation} \\ \sum_{p \in j} S_{ip} m_p \Delta \mathbf{x}_p & \text{displacement extrapolation} \end{cases} \quad (2.3)$$

$$\Omega_{i,c;j} = \begin{cases} \sum_{p \in j} S_{ip} V_p & \text{3D} \\ \sum_{p \in j} S_{ip} A_p & \text{2D and axisymmetric} \end{cases} \quad (2.4)$$

$$\mathbf{g}_{i,c;j} = \begin{cases} \sum_{p \in j} S_{ip} \mathbf{G}_p V_p & \text{3D} \\ \sum_{p \in j} S_{ip} \mathbf{G}_p A_p & \text{2D and axisymmetric} \end{cases} \quad (2.5)$$

The time step ( $k$ ) has been omitted.

After the extrapolations are done, each node is revisited for more calculations and when needed to impose multimaterial and crack contact and imperfect interfaces. The steps are

1. Calculate total mass on number of particles in each velocity field.
2. Copy the initially extrapolated momenta  $\mathbf{p}_{i,c;j}$  for all crack and material velocity fields on all nodes.
3. When in multimaterial mode, check if each node might have contact, which means the nodes has non-zero mass from more than one material. If it does, create a **MaterialContactNode** object. The end result will be a list of all nodes that might have contact in the vector **materialContactNodes** (which is static member of **MaterialContactNode** class).
4. When have cracks, check if each node might have crack contact (see possible fields list above that might have contact). If it does, create a **CrackNode** object. The end result will be a list of all nodes that might have crack contact in the vector **crackContactNodes** (which is static member of **CrackNode** class).

### 2.3.1 Contact Calculations

Once all contacting nodes are found, revisit each node (in the two lists determined above) and impose contact mechanics. The contact mechanics for both multimaterials and cracks are briefly described here. More details are below:



1. First determine and crack or materials and in contact. This calculation requires evaluation of both normal vector at the contacting surface and the separation between the two surface. The accuracy of all contact calculations depends on ability to find normals and separation.
2. If materials or cracks are in contact, calculate the change in momenta needed to reflect the contact physics as  $\Delta p_{i,c:a}$ . Add this momentum to material  $a$  and opposite of this momentum to material  $b$ . Thus momentum is conserved. Some approximate scheme is usually needed to handle multimaterial contact involving more than two materials. For cracks, check contact on surface of first crack (*e.g.*, [0]&[1] and [2]&[3]) and on second cracks (*e.g.*, [0]&[2] and [1]&[3]). Any resulting  $\Delta p_{i,c:a}$  is added to two sides of the target crack. If needed, the added momentum is spread out over all materials (weighted by their mass fraction) if the crack velocity field has more than one material.
3. If materials or cracks are connected by an imperfect interface, the calculations are different. Here calculate the interface force  $\mathbf{f}_{i,c:a}^{(INT)}$  on material  $a$  which will depend on calculated material separation. If normal or tangential force is found, set the contact  $\Delta p_{i,c:a}$  in the respective direction to zero. Some interface laws can become unstable (if too stiff). If they get too stiff, set interface force to zero, but keep corresponding component  $\Delta p_{i,c:a}$  to reflecting “sticking” mechanics. The final  $\Delta p_{i,c:a}$  is added to materials (and crack surfaces), but the interfacial force is not added (it is added later during momentum update instead).

## 2.4 Update Stresses and Strains on Particle

When updating stress and strains before finding forces and updating momenta (needed for **USAVG**± and for **USF** update methods), extrapolate velocity gradient ( $\nabla \mathbf{v}$ ) to the particles and implement the chosen constitutive laws. Because of contact calculations done after mass and momentum extrapolation, the extrapolated velocity gradient will come from velocities that account for contact (but not yet for imperfect interface forces).

## 2.5 Project Forces to the Grid

For each material point, extrapolate  $f_{i,c:m}$  due to particle stresses and  $f_{i,T,c:m}$  due to particle traction loads.. The values from particle  $p$  are added to material velocity field **matfld** of crack velocity field **vfld**, where **vfld** was found during time step initialization and **matfld** is known by the material point’s material type.

After the force extrapolations are done, the following calculations complete the force calculations:

1. Restore momentum original extrapolated to the grid in mass and momentum task.
2. Simple particle force boundary conditions are added for those particles (although these force are actually set in the initialization task).
3. If any cracks have traction laws, forces currently stored on crack particles are interpolated to nodes. Note that this extrapolation uses ordinary grid shape functions because crack particles have no volume and GIMP shape function imply a particle domain. Also not that shape function are renormalized if needed to guarantee 100% of force gets to nodes, which occurs if crack particle sees nodes with zero mass.

4. Add any body forces (such as gravity).

Once all above force steps are done, the force on nodes with grid velocity boundary conditions are changed to be consistent with that condition. The updated momentum on a node with prescribed velocity in the  $\hat{n}$  direction needs to be:

$$\mathbf{p}_{i,c;m}^{(k+1)} = m_{i,c;m} v_i^{(BC)} \hat{n} + \left( \left( \mathbf{p}_{i,c;m}^{(k)} + \mathbf{f}_{i,c;m}^{(tot)} \Delta t \right) \cdot \hat{t} \right) \hat{t} \quad (2.6)$$

$$= \mathbf{p}_{i,c;m}^{(k)} + \left( \mathbf{f}_{i,c;m}^{(tot)} + f_{i,c;m}^{(BC)} \hat{n} \right) \Delta t \quad (2.7)$$

where  $f_{i,c;m}^{(BC)}$  is magnitude of reaction force applied in the  $\hat{n}$  direction to satisfy the boundary condition on that node. The normal component of the updated momentum must give the boundary condition or:

$$\mathbf{p}_{i,c;m}^{(k+1)} \cdot \hat{n} = m_{i,c;m} v_i^{(BC)} = \mathbf{p}_{i,c;m}^{(k)} \cdot \hat{n} + (\mathbf{f}_{i,c;m}^{(tot)} \cdot \hat{n}) \Delta t + f_{i,c;m}^{(BC)} \Delta t \quad (2.8)$$

This equation can be solved to find magnitude of the reaction force:

$$f_{i,c;m}^{(BC)} = \frac{m_{i,c;m} v_i^{(BC)} - \mathbf{p}_{i,c;m}^{(k)} \cdot \hat{n}}{\Delta t} - \mathbf{f}_{i,c;m}^{(tot)} \cdot \hat{n} \quad (2.9)$$

To impose this condition after all other forces are found, revisit all nodes with boundary conditions and add  $f_{i,c;m}^{(BC)} \hat{n}$  to nodal force, which results in

$$\mathbf{f}_{i,c;m}^{(tot)'} = \mathbf{f}_{i,c;m}^{(tot)} - (\mathbf{f}_{i,c;m}^{(tot)} \cdot \hat{n}) \hat{n} + \frac{m_{i,c;m} v_i^{(BC)} - \mathbf{p}_{i,c;m}^{(k)} \cdot \hat{n}}{\Delta t} \hat{n} \quad (2.10)$$

$$= \mathbf{f}_{i,c;m}^{(tot)} - \left( (\mathbf{f}_{i,c;m}^{(tot)} \cdot \hat{n}) + \frac{\mathbf{p}_{i,c;m}^{(k)} \cdot \hat{n}}{\Delta t} \right) \hat{n} + \frac{m_{i,c;m} v_i^{(BC)}}{\Delta t} \hat{n} \quad (2.11)$$

To support superposition of velocity conditions on the same node (but in the same direction  $\hat{n}$ ), the boundary condition value is replaced by

$$m_i^{(n)} v_i^{(BC)} = \sum_k m_i^{(n)} v_i^{(BC,k)} \quad (2.12)$$

where  $v_i^{(BC,k)}$  is velocity for boundary condition  $k$ . A single node can superpose boundary conditions in different directions only if the directions are orthogonal. Note that this approach to velocity boundary conditions (rather than just imposing them in the subsequent momentum update) is required to have the accelerations determined from nodal forces be the correct accelerations needed to move the particle from its initial position in the time step to the proper position that reflects the boundary conditions.

## 2.6 Update Nodal Momenta

This task updates the nodal momenta using

$$\mathbf{p}_{i,c;m}^{(k+1)} = \mathbf{p}_{i,c;m}^{(k)} + \mathbf{f}_{i,c;m} \Delta t \quad (2.13)$$

Now, depending on the problem, this update may cause more contact. Thus the update must be followed by adjusting momenta at all nodes with overlapping materials and at all cracks now in

contact. The logic again is that the final momenta in this step should be identical when doing multiple velocity fields with stick contact to the momenta when using a single velocity field.

The contact calculations are summarized in section 2.3.1. The main difference is that interfacial force is added (to both field momental and field force) when contact done with updated momenta. In additions, the code tracks contact forces (with rigid particles) and friction heating (for frictional contact). These calculations are only done in contact calculations associated with the momentum update.

If contact calculations are done, the velocity boundary conditions are imposed again. If not done, momenta change added in contact calculation might change those conditions. It is easier to reimpose them here then to try and track all times other calculations might change them (and imposing them might eliminate any round off error or of counting on momentum update to end with exact result).

## 2.7 Update the Particles

Update particle position and velocity using nodal velocity (from  $\mathbf{v}_{i,c,m} = \mathbf{p}_{i,c,m}^{(k+1)}/m_{i,c,m}$ ) and acceleration (from  $\mathbf{a}_{i,c,m} = \mathbf{f}_{i,c,m}/m_{i,c,m}$ ) found in material velocity field `matfld` (for type of particle) of crack velocity field `vfld` (determined on current time step for the particle).

## 2.8 Optional Update Strain Last

Optionally the calculations can update stain after the momentum update. This calculation is done for methods `USAVG±` and for `USL±`. The - (minus) version updates using the updated velocity and their extrapolated velocity gradient (exactly like the first stress and strain update). The + (plus) method does new extrapolation of momentum to the grid. Because this projection might induce new contact, adjust momenta again at nodes with multiple materials and at crack surfaces (same method as in section 2.3.1. If contact calculations are needed, velocity boundary condition are imposed to insure extrapolated velocity gradients will reflect contact physics. These final set of contact calculations will not include imperfect interface forces of cohesive law forces.

## 2.9 Move Cracks

The crack surfaces and crack plane are updated using the center-of-mass velocity field of each crack velocity field (for those that have multiple materials). In addition, the new crack opening displacements are used to calculation traction loads to be input to the next time step during its force calculations.



## Chapter 3

# Contact Calculations

### 3.1 Finding Normal Vector for Contact and Interfaces

Get the normal correct is vital to accurate contact in MPM. Most problems involving contact can be traced to inaccurate normals and normal are prerequisite to all other contact calculations. Although dynamic contact detection is advantage of MPM, finding the normal is cost of using that advantage. Consider material  $a$  and  $b$  in one crack velocity field on node  $i$ . All except material ID in property subscripts will be omitted for clarity. The goal is to find normal  $\hat{\mathbf{n}}$  directed from  $a$  to  $b$ . Ignoring  $b$ , a unit normal can be estimated from volume gradient:

$$\hat{\mathbf{n}} = \frac{\mathbf{g}_a}{\|\mathbf{g}_a\|} \quad (3.1)$$

Note that code should use volume gradient instead of mass gradient to handle situations involving contact between materials with different densities. This normal may or may not be accurate and may or may not be same as corresponding normal found from material  $b$ . Some MPM codes found normal for each material and looped over materials. But when normals are not equal and opposite, the momentum would not be conserved. A better approach is to use a single normal for the material pair. No one method works best. The options in `OSParticulas` are:

- **AVGGRAD**: Find normal from volume average volume gradients of the two materials:

$$\hat{\mathbf{n}}\|\hat{\mathbf{n}}\| = \mathbf{g}_a + \mathbf{g}_b \quad (3.2)$$

The justification is that averaging is often best and volume weighting counts the one with more information (*i.e.* closest to the contact point) the most. The terms are simply added because  $\mathbf{g}_a$  and  $\mathbf{g}_b$  are already volume-weighted sums (see Eq. (2.5)).

- **MAXGRAD**: Find normal from materials whose volume gradient has the largest magnitude:

$$\hat{\mathbf{n}}\|\hat{\mathbf{n}}\| = \begin{cases} \mathbf{g}_a & \text{if } \|\mathbf{g}_a\| > \|\mathbf{g}_b\| \\ -\mathbf{g}_b & \text{otherwise} \end{cases} \quad (3.3)$$

The justification is that one with more information, represented by higher absolute magnitude, might have more accurate information.

- **MAXVOL**: Find normal from materials whose volume has the largest magnitude:

$$\hat{\mathbf{n}}\|\hat{\mathbf{n}}\| = \begin{cases} \mathbf{g}_a & \text{if } \Omega_a > \Omega_b \\ -\mathbf{g}_b & \text{otherwise} \end{cases} \quad (3.4)$$

The justification is that one with more information, represented by more volume, might have more accurate information.

- **OWN**: Find normal from each material and consider contact effects separately for all materials:

$$\hat{\mathbf{n}} = \frac{\mathbf{g}_a}{\|\mathbf{g}_a\|} \quad (3.5)$$

This method is never recommended and only included in `OSParticulas` to allow comparison to old MPM methods for handling contact.

- **SPECIFY**: Instead of calculating the normals, prescribe a normal that applies to all contact situations. The normal is directed from the lower number material (in the list of defined materials) to the higher one. This approach is only viable when all contact surfaces have the same normal and the normal does not change during the simulation. This situation is uncommon in real problems. Note that if the contact is with a rigid material (see details below), a specified normal mans from non-rigid into the rigid material, regardless of order of the material numbering.
- **LINREG** and **LOGREG**: These regression methods are based on machine learning methods and find normal by looking at geometry of particle that see each node (*i.e.*, they make no use of extrapolated gradients). Details are given in chapter ??.

## 3.2 Detecting Contact

The next task is to decide if materials are in contact or if they are separated. When in contact, the materials are often left to move in the separate velocity fields. When they are in contact, however, the momenta have to change to prevent interpenetration and to implement contact mechanics. If the contact law includes adhesion, it may be necessary to examine separated material as well to decided if they have overcome some input adhesion value. For imperfect interface, forces generally need to be added for both contacted and separated interfaces. All cases need to know which has occurred. Below are the following steps used in `OSParticulas`.

First, a mass-prescreening option is used: The final change in momentum is related to reduced mass on the node, which means the change will be neglible if the mass ratio is high or small. This prescreen skips contact calculations if:

$$\frac{m_a}{m_{tot}} < \epsilon \quad \text{or} \quad \frac{m_a}{m_{tot}} > 1 - \epsilon \quad (3.6)$$

where  $\epsilon$  is a small number. The current version of `OSParticulas` uses  $\epsilon = 1.e^{-5}$  (might change in the future).

A discarded volume prescreening test used to ignore nodes where total volume on a node (sum of material volumes) is compared to total volume associated with that node (*i.e.*, cell volume — for grid with unequal element sizes, it is net volume of cells around node  $i$ ). Let  $V_r$  be ratio of material volumes to node  $i$  cell volume. This prescreen will ignore contact on this node if  $V_r < V_{min}$  where  $V_{min}$  is simulations parameter. The logic is not that contact at interfaces are not present, but rather that this node is likely on the edge of a material and the calculations will be inaccurate. It might better to skip them to to attempt them. This methods was never found to be effective and has now been removed from `OSParticulas`.

If the calculations are nor screened out, the normal is calculated (as explained above) and then contact is determined by one or both of the following checks — an approach velocity check and a

interface separation check. Early MPM work used the first method only, but current thinking is that second is always needed. The option to skip the second check has been removed.

### 3.2.1 Approach Velocity

For two materials with velocities  $\mathbf{v}_a$  and  $\mathbf{v}_b$ , the center of mass velocity is

$$\mathbf{v}_c = \frac{m_a \mathbf{v}_a + m_b \mathbf{v}_b}{m_a + m_b} \quad (3.7)$$

where  $m_a$  and  $m_b$  are the masses for the two materials. Next, define  $\Delta \mathbf{p}_a$  as the momentum change required on material  $a$  for its velocity to change from  $\mathbf{v}_a$  to  $\mathbf{v}_c$  or

$$\mathbf{v}_a + \frac{\Delta \mathbf{p}_a}{m_a} = \mathbf{v}_c \quad \text{which means} \quad \Delta \mathbf{p}_a = -m_a(\mathbf{v}_a - \mathbf{v}_c) \quad (3.8)$$

Inserting the definition of  $\mathbf{v}_c$  leads to

$$\Delta \mathbf{p}_a = -m_a(\mathbf{v}_a - \mathbf{v}_c) = m_b(\mathbf{v}_b - \mathbf{v}_a) = \frac{m_a \mathbf{p}_b - m_b \mathbf{p}_a}{m_a + m_b} \quad (3.9)$$

where  $\mathbf{p}_a$  and  $\mathbf{p}_b$  are the nodal momenta for the two materials. Notice that

$$\Delta \mathbf{v} = \mathbf{v}_b - \mathbf{v}_a = \frac{\Delta \mathbf{p}_a}{m_b} + \frac{\Delta \mathbf{p}_a}{m_a} = \frac{\Delta \mathbf{p}_a}{m_{red}} \quad (3.10)$$

where  $m_{red} = m_a m_b / (m_a + m_b)$  is the reduced nodal mass. Thus the relative velocity of approach of the two surfaces is proportional to, and in the same direction as,  $\Delta \mathbf{p}_a$ . For detecting contact based on approach velocity, the two surfaces must be approaching each other. With a normal directed from  $a$  to  $b$ , this criteria implies  $\Delta \mathbf{v} \cdot \hat{\mathbf{n}} < 0$  or identically that  $\Delta \mathbf{p}_a \cdot \hat{\mathbf{n}} < 0$ .

We thus assume that under all contact situations that  $\Delta \mathbf{p}_a \cdot \hat{\mathbf{n}} < 0$ . In other words, this criterion must always be included. As shown below, this criterion (if used in combination with a separation check) is equivalent to a finding the contact surface have negative traction (*i.e.*, compression stress) in the normal direction. Many contact laws (such as friction) are based on compression state. With this check, those contact laws can assume that  $\Delta \mathbf{p}_a \cdot \hat{\mathbf{n}} < 0$  in their calculations.

### 3.2.2 Determination of Material Separation

The use of approach velocity alone is not enough. It often implies contact even when edges are clearly separated. The detection of contact is required calculation of the separation between the materials in the normal direction. In physical terms, the momentum change in normal direction of  $\Delta \mathbf{p}_a \cdot \hat{\mathbf{n}} < 0$  corresponds to contact compression if and only if the two materials are in contact. If they are separated, that term corresponds to deceleration need to slow the material fields done as if they were a single material. In other words, contact detection is to find nodes with compression stress between the two materials and that determination needs to know both  $\Delta \mathbf{p}_a \cdot \hat{\mathbf{n}}$  and whether to not the surfaces are actually in contact. Three approaches are possible:

1. Extrapolate displacement to the grid. While this works, it requires that the two materials start out in contact such that zero displacement is equal to contact and that total deformation in the problem is small. When the materials come into contact only after motion (such as closing a gap), displacements are not enough.

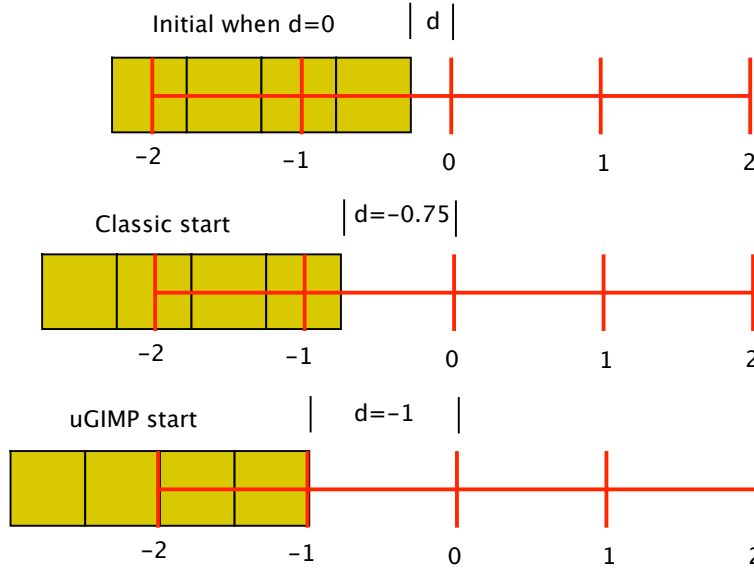


Figure 3.1: One dimensional extrapolation of particle position to grid node. Node  $i$  is at  $x = 0$  ( $x$  in units of cells size).  $d$  is distance from edge of material points to node  $i$ . The “Classic” and “uGIMP” start are the position the extrapolation first becomes nonzero.

2. Extrapolate particle position. While this method is more general, the results have to be correct for grid effects.
3. The regression methods (LINREG and LOGREG) find normal vector from particle around a node. Once normal to interface is found, the separation can be found from perpendicular distance of closest particles to the nodes. This calculation needs to be correct for particle radius. The method is described in chapter ??.

By the extrapolation methods, the interfacial displacement discontinuity on node  $i$  (subscript  $i$  omitted for clarity) is:

$$\delta = \mathbf{x}_b - \mathbf{x}_a \quad (3.11)$$

where  $a$  and  $b$  are two materials (for material contact) or above and below a crack. The separation is then given by

$$\delta_n = \delta \cdot \hat{\mathbf{n}} \quad (3.12)$$

When extrapolating displacements,  $\delta_n \leq 0$  can be used to detect contact, but when extrapolating positions, more calculations are needed.

To correct extrapolated positions for material separation calculations we consider 1D extrapolation near the edge of an object. Figure 3.1 shows group of undeformed particles (when using two particle per element in each direction) approaching node  $i$  at  $x = 0$  with  $d$  as the actual distance from the edge of the material points to that node. The “Classic start” shows that position on node  $i$  will be zero until  $d = -0.75$  or until the center of the first particle enters the element to the left of node  $i$ . Both GIMP and CPDI will start extrapolating to node  $i$  the instant the edge enters that element (see “GIMP start”).

For each  $d_m$ , or actual signed distance from edge of material  $m$  ( $a$  or  $b$ ) to node  $i$ , we can calculate  $d_m^{(ext)} = (\mathbf{x}_m - \mathbf{x}_i) \cdot \hat{\mathbf{n}}$ , or the distance from extrapolated position  $\mathbf{x}_m$  to node  $i$  using



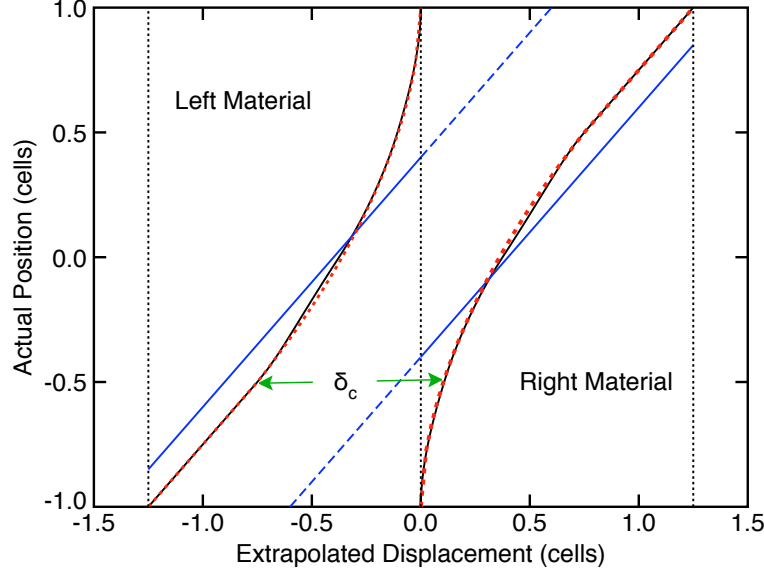


Figure 3.2: Plot of actual distance from edge to node  $i$  at  $x = 0$  ( $d$ ) as a function of extrapolated position ( $d_{ext}$  on node  $i$  when using one particle per cell). The “Left Material” is for edge approaching from the left. The “Right Material” is for an edge departing to the right.

contact normal and a given shape function method. Once these results are found, we invert and the results as shown in Fig. 3.2. The “Left Material” is for edge approaching from the left (if is material  $a$  if  $\hat{n}$  points to the right). When the particle edge first enters the element to left of node  $i$  (assuming GIMP or CPDI), the extrapolated position (from particle center) will be -1.25 from the node. As the edge approaches, the extrapolation will decrease, but remains negative. When particles of material  $m$  completely fill the two elements on opposite sides of the node  $i$ , the extrapolated position will be zero. Thus extrapolation values from -1.25 to 0 are mapping to actual edge distances from  $d = -1$  to 1. The dotted red lines shows this mapping for GIMP (Classic would be piecewise linear approximation starting at (-1,-0.75) and CPDI should be similar to uGIMP, but neither are shown). The dotted red curve for “Right Material” is the same calculation for edge of particles departing to the right (or material  $b$  in  $\hat{n}$  points to the right). It is always positive and not 0 to 1.25 maps to  $d = -1$  to 1. More details are given in supplemental material to our nanoindentation paper [1]

Now imagine two materials (or crack surfaces) near node  $i$ . The normal from material  $a$  to  $b$  is taken in the positive direction, such material  $a$  is the “left” material and  $b$  is the “Right” material. The separation is written as:

$$\delta_n = d_b(d_b^{(ext)}) - d_a(d_a^{(ext)}) \quad \text{where} \quad d_m^{(ext)} = (\mathbf{x}_m - \mathbf{x}_i) \cdot \hat{n} \quad (3.13)$$

is the extrapolated (or apparant) distance from edge to node and  $d_m(d_m^{(ext)})$  converts extrapolated distance to actual distance. The final  $\delta_n$  is actual separation and  $\delta_n \leq 0$  is required for contact.

OSParticulas has two methods to find  $\delta_n$ . The first method (as published in my papers) is to assume the  $d_m()$  functions are linear and given by:

$$d_m(d_m^{(ext)}) = \begin{cases} d_a^{(ext)} + 0.4\Delta x & \text{left or } m = a \\ d_b^{(ext)} - 0.4\Delta x & \text{right or } m = b \end{cases} \quad (3.14)$$

where  $\Delta x$  is an effective cell size (and equal to cell size with square (2D) or cubic (3D) cells, but changed to  $h_\perp$  for rectangular cells and changed for arbitrary neighbor sizes for a Tartan grid). With this equation, the material separation becomes

$$\delta_n = (\mathbf{x}_b - \mathbf{x}_i) \cdot \hat{\mathbf{n}} - 0.4\Delta x - ((\mathbf{x}_a - \mathbf{x}_i) \cdot \hat{\mathbf{n}} + 0.4\Delta x) = \boldsymbol{\delta} \cdot \hat{\mathbf{n}} - 0.8\Delta x \quad (3.15)$$

My paper describes this approach as subtracting a contact constant, but the net effect of this approach is same as assuming the above linear conversion form  $d_m^{(ext)}$  to  $d_m$ . This net subtraction of  $0.8\Delta x$  was found by a weighted average of  $d_b^{(ext)} - d_a^{(ext)}$  when edges are in contact and counted the results near  $d = 0$  more. The weighting logic was that momentum change is greatest when  $d$  is close to zero or contact is near a node. Although the subtraction value can be varied to any value, it appears results far from  $0.8\Delta x$  cells give poor results or that  $0.8\Delta x$  cells appears to be universal value when using this linear method and when have two particles per cell.

An alternate approach is to fit the inverted curves derived from shape function calculations. The solid black lines show the fits:

$$\frac{d_m(d_m^{(ext)})}{\Delta x} = \begin{cases} 1 - 2 \left( \frac{-d_a^{(ext)}}{1.25\Delta x} \right)^{0.58} & \text{left or } m = a \\ 2 \left( \frac{d_b^{(ext)}}{1.25\Delta x} \right)^{0.58} - 1 & \text{right or } m = b \end{cases} \quad (3.16)$$

Thus in contact calculations, one finds  $d_m(d_m^{(ext)})$  for left ( $a$ ) and right ( $b$ ) materials using this equation, and then subtracts for the actual  $\delta_n$  or surface separation. Contact occurs when  $\delta_n \leq 0$ . This option is selected in `OSParticulas` by setting the `ContactPosition` to a negative number and the absolute value of that number becomes the exponent in the above fits. Like 0.8 method, the exponent can be varied, but it appears that -0.58 (in the command, which uses exponent of 0.58) is universal value that should not be adjusted.

A slightly better fit can be found by recognizing that the shape function calculation is exactly linear over a small region. Using piecewise function with part linear and the rest a power law gives:

$$\frac{d_m(d_m^{(ext)})}{\Delta x} = \begin{cases} \begin{cases} \frac{d_a^{(ext)}}{\Delta x} + 0.25 & d_a^{(ext)} < -0.75\Delta x \\ 1 - \frac{3}{2} \left( \frac{-4d_a^{(ext)}}{3\Delta x} \right)^{0.5546} & d_a^{(ext)} > -0.75\Delta x \end{cases} & \text{left or } m = a \\ \begin{cases} \frac{3}{2} \left( \frac{4d_b^{(ext)}}{3\Delta x} \right)^{0.5546} - 1 & d_b^{(ext)} < 0.75\Delta x \\ \frac{d_b^{(ext)}}{\Delta x} - 0.25 & d_b^{(ext)} > 0.75\Delta x \end{cases} & \text{right or } m = b \end{cases} \quad (3.17)$$

The difference with first power law in Eq. (3.16) is small. This revised form is not yet implemented in code.

The above calculations assume two particles per cell. A displacement criterion to detect contact, however, depends on the number of particles. For one particle per cell, the linear method changes to:

$$d_m(d_m^{(ext)}) = \begin{cases} d_a^{(ext)} + 0.535\Delta x & \text{left or } m = a \\ d_b^{(ext)} - 0.535\Delta x & \text{right or } m = b \end{cases} \quad (3.18)$$

or

$$\delta_n = \boldsymbol{\delta} \cdot \hat{\mathbf{n}} - 1.07\Delta x \quad (3.19)$$

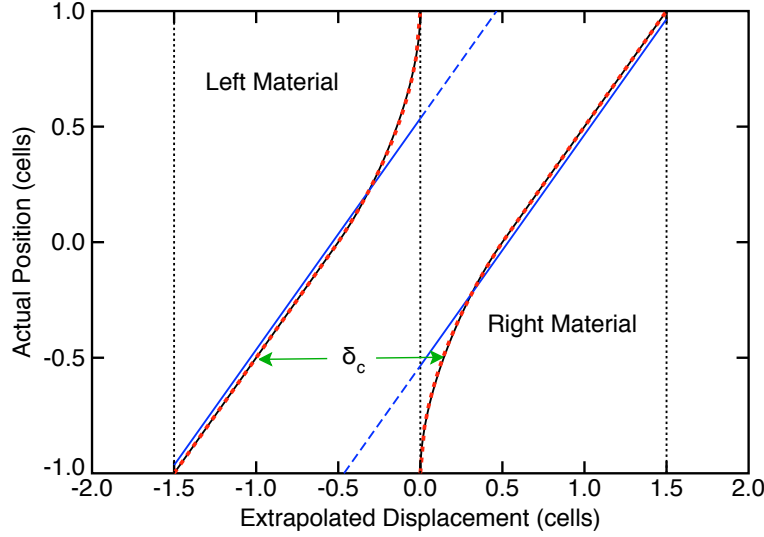


Figure 3.3: Plot of actual distance from edge to node  $i$  at  $x = 0$  ( $d$ ) as a function of extrapolated position ( $d_{ext}$  on node  $i$  when using one particle per cell. The “Left Material” is for edge approaching from the left. The “Right Material” is for an edge departing to the right.

A power law fit can also be derived. For one particle, a significant portion of the curve is exactly linear. A piecewise fit with a linear portion and power law portion is given by:

$$\frac{d_m(d_m^{(ext)})}{\Delta x} = \begin{cases} \begin{cases} \frac{d_a^{(ext)}}{\Delta x} + 0.5 & d_a^{(ext)} < -0.5\Delta x \\ 1 - \left(\frac{-2d_a^{(ext)}}{\Delta x}\right)^{0.5387} & d_a^{(ext)} > -0.5\Delta x \end{cases} & \text{left or } m = a \\ \begin{cases} \left(\frac{2d_b^{(ext)}}{\Delta x}\right)^{0.5387} - 1 & d_b^{(ext)} < 0.5\Delta x \\ \frac{d_b^{(ext)}}{\Delta x} - 0.5 & d_b^{(ext)} > 0.5\Delta x \end{cases} & \text{right or } m = b \end{cases} \quad (3.20)$$

A plot of the linear and power-law functions for one particle per cell is in Fig. 3.3

A final addition is needed when  $\delta_n$  is found after the momentum update. Because  $\delta_n$  is found from initially extrapolated positions, it should be updated for interfacial movement during the time step. The correction is

$$\delta'_n = \delta_n + (\mathbf{v}_b - \mathbf{v}_a) \cdot \hat{\mathbf{n}} \Delta t = \delta_n + \frac{\Delta \mathbf{p}_a \cdot \hat{\mathbf{n}} \Delta t}{m_{red}} \quad (3.21)$$

In summary, the three extrapolation options are in `OSParticulas` are:

1. No `ContactPosition` command — if contact position is never set, contact is determined by extrapolated displacements and no correction is applied to the extrapolated values. (Note: this is **not** the same as setting contact position to 0, which is to extrapolate position instead of displacement but not correct the values; these results would be very bad).
2. Positive `ContactPosition` — contact is determined by extrapolated positions which are adjusted by constant value provided (should probably always be 0.8 for two particles per cell or 1.07 for two.)

3. Negative **ContactPosition** — contact is determined by extrapolated positions which are adjusted by Eq. (3.16) where exponent is absolute value of value provided (should probably always be -0.58). This option is for two particle only. The piecewise curves for two or one particle per cell are not yet implemented.

The fourth option is to find separation for regression results (see chapter ??). The **ContactPosition** setting has no affect on this lat option.

### 3.3 Changes in Momentum for Contact

The relatively easy part s(compared to finding normal and detecting contact) are changing momenta and finding proper force needed to enforce a contact law. The basic analysis is for material  $a$  and contact with material  $b$ , but to handle general problems where a node may see more that two materials (although it may not be handled rigorously), the following algorithm is used. At a node with any number of materials, the center of mass results are:

$$\mathbf{p}_c = \sum_i \mathbf{p}_i \quad m_c = \sum_i m_i \quad \mathbf{v}_c = \frac{\mathbf{p}_c}{m_c} \quad (3.22)$$

We define a *virtual* material  $b$  derived from all mass and momenta not associated with material  $a$ :

$$\mathbf{p}_b = \mathbf{p}_c - \mathbf{p}_a \quad m_b = m_c - m_a \quad \mathbf{v}_b = \frac{\mathbf{p}_b}{m_b} \quad (3.23)$$

The equations for  $\Delta \mathbf{p}_a$  and  $\Delta \mathbf{v}$  from above are unchanged, but in terms of just  $a$  and  $c$ , they become

$$\Delta \mathbf{p}_a = \frac{m_a}{m_c} \mathbf{p}_c - \mathbf{p}_a, \quad \Delta \mathbf{v} = \frac{\Delta \mathbf{p}_a}{m_{red}}, \quad \text{and} \quad m_{red} = \frac{m_a(m_c - m_a)}{m_c} \quad (3.24)$$

All these results reduces exactly to two-material case for nodes with two materials.

To implement contact conditions, with inclusion of nodes with three or more materials, each material is considered relative to the center of mass velocity and  $\Delta \mathbf{p}_a$  and  $\Delta \mathbf{v}$  are found. Stick conditions are trival. The momentum change of  $\Delta \mathbf{p}_a$  is applied to material  $a$ . These momenta changes correspond to normal and tangential forces of

$$\mathbf{f}_n = \frac{(\Delta \mathbf{p}_a \cdot \hat{\mathbf{n}})}{\Delta t} \hat{\mathbf{n}} \quad \text{and} \quad \mathbf{f}_t = \frac{(\Delta \mathbf{p}_a \cdot \hat{\mathbf{t}})}{\Delta t} \hat{\mathbf{t}} \quad (3.25)$$

where  $\Delta t$  is time step and  $\hat{\mathbf{t}}$  is unit vector in the tangential direction of motion. We choose  $\hat{\mathbf{t}}$  to be in the same direction as the relative motion in the tangential direction such that  $(\mathbf{v}_b - \mathbf{v}_a) \cdot \hat{\mathbf{t}} > 0$  and identically such that  $\Delta \mathbf{p}_a \cdot \hat{\mathbf{t}} > 0$  or that the angle between  $\hat{\mathbf{t}}$  and  $\mathbf{v}_b - \mathbf{v}_a$  is between -90 and 90 degrees. Give  $\hat{\mathbf{n}}$ ,  $\hat{\mathbf{t}}$  can be found from

$$\hat{\mathbf{t}} = \frac{\Delta \mathbf{p}_a - (\Delta \mathbf{p}_a \cdot \hat{\mathbf{n}}) \hat{\mathbf{n}}}{\|\Delta \mathbf{p}_a - (\Delta \mathbf{p}_a \cdot \hat{\mathbf{n}}) \hat{\mathbf{n}}\|} \quad (3.26)$$

The positive normal contact force is thus  $f_n = -(\Delta \mathbf{p}_a \cdot \hat{\mathbf{n}})/\Delta t$  and the positive sliding force for interfaces to stick (move in same velocity field) is  $f_t = (\Delta \mathbf{p}_a \cdot \hat{\mathbf{t}})/\Delta t$ . We write the normal pressure ( $N$ ) and sticking traction ( $S_{stick}$ ) as

$$N = -\frac{\Delta \mathbf{p}_a \cdot \hat{\mathbf{n}}}{A_c \Delta t} \quad \text{and} \quad S_{stick} = \frac{\Delta \mathbf{p}_a \cdot \hat{\mathbf{t}}}{A_c \Delta t} \quad (3.27)$$

where  $A_c$  is contact area (calculated by methods in [2]). Implementation of general contact laws for sliding force  $S_{slide}$  as a function of normal pressure  $N$  is done as follows:

1. If  $S_{stick} < S_{slide}$ , then the driving forces for frictional sliding are too small to overcome sticking meaning the two materials stick rather than slide. Set final momentum change to  $\Delta \mathbf{p}'_a = \Delta \mathbf{p}_a$ .
2. But, if  $S_{stick} > S_{slide}$ , the surface would rather slide at the lower  $S_{slide}$  then stick. Set final momentum change to

$$\Delta \mathbf{p}'_a = d_n \hat{\mathbf{n}} + (S_{slide} A_c \Delta t) \hat{\mathbf{t}} \quad (3.28)$$

where  $d_n = \Delta \mathbf{p}_a \cdot \hat{\mathbf{n}} = -f_n \Delta t$

3. Change the momenta for materials  $a$  and  $b$  to

$$\mathbf{p}'_a = \mathbf{p}_a + \Delta \mathbf{p}'_a \text{ and } \mathbf{p}'_b = \mathbf{p}_b - \Delta \mathbf{p}'_a \quad (3.29)$$

A subtle point, which is only found by requiring two identical materials in contact at a perfect interface to match single material mode (see below), is that when consistent forces are found for grid boundary conditions, they must be based on  $\mathbf{p}'_a$  and not on the originally extrapolated  $\mathbf{p}_a$ . The forces are found from stresses in materials that are based on contact conditions being satisfied and therefore boundary conditions force must do the same.

4. When momentum change is found *after* finding forces and updating momenta, the updated forces need to be kept with changed momenta. The nodal forces should be changed to:

$$\mathbf{f}'_a = \mathbf{f}_a + \frac{\Delta \mathbf{p}_a^{+'}}{\Delta t} \quad (3.30)$$

where  $\Delta \mathbf{p}_a^{+'}$  is momentum change in contact calculations using updated momenta.

5. If three or more materials are present, some modifications are needed (see below).

See our paper on contact laws for more details and handling various types of contact laws (such as those with velocity dependent laws or with adhesion). When frictional sliding occurs, the energy dissipated can be found from frictional work and input to heat conduction calculations. This option is also described in generalized contact paper [3].

### 3.3.1 Material Velocity Field Updates

When the calculations update momenta,  $\mathbf{p}_a$  will have changed to  $\mathbf{p}'_a$  (similar for material  $b$ ) and  $\mathbf{f}_a$  will have been extrapolated to the grid; the update will be:

$$\mathbf{p}_a^+ = \mathbf{p}'_a + \mathbf{f}_a \Delta t \quad (3.31)$$

Here  $\mathbf{f}_a$  are the forces extrapolated from particle stresses that are found when interfaces satisfy boundary conditions. After updating momentum, new contact calculations will find  $\Delta \mathbf{p}_a^{+'}$  meaning that the final particle momentum and force are:

$$\mathbf{p}_a^{+'} = \mathbf{p}_a^+ + \Delta \mathbf{p}_a^{+'} \quad \text{and} \quad \mathbf{f}'_a = \mathbf{f}_a + \frac{\Delta \mathbf{p}_a^{+'}}{\Delta t} \quad (3.32)$$

The total particle update becomes

$$\mathbf{p}_a^{+'} = \mathbf{p}'_a + \mathbf{f}'_a \Delta t = \mathbf{p}'_a + \left( \mathbf{f}_a + \frac{\Delta \mathbf{p}_a^{+'}}{\Delta t} \right) \Delta t \quad (3.33)$$

The final velocity and acceleration for material  $a$  are:

$$\mathbf{v}_a^+ = \frac{\mathbf{p}_a^{+'}}{m_a} \quad \text{and} \quad \mathbf{a}_a^+ = \frac{\mathbf{f}'_a}{m_a} \quad (3.34)$$

To validate this process, show that for two identical materials with a perfect interface translates into material velocity and acceleration equal to center of mass velocity and acceleration. When two identical materials are in contact and the interface is perfect, we know that

$$\mathbf{p}'_a = \frac{m_a}{m_i} \mathbf{p}_i \quad \text{and} \quad \mathbf{p}_a^{+'} = \frac{m_a}{m_i} (\mathbf{p}_i + \mathbf{f}_i \Delta t), \quad (3.35)$$

and

$$\Delta \mathbf{p}_a^{+'} = \mathbf{p}_a^{+'} - \mathbf{p}_a^+ = \frac{m_a}{m_i} (\mathbf{p}_i + \mathbf{f}_i \Delta t) - (\mathbf{p}'_a + \mathbf{f}_a \Delta t) = \frac{m_a}{m_i} \mathbf{f}_i \Delta t - \mathbf{f}_a \Delta t \quad (3.36)$$

Solving for forces gives:

$$\frac{\mathbf{f}_i}{m_i} = \frac{1}{m_a} \left( \mathbf{f}_a + \frac{\Delta \mathbf{p}_a^{+'}}{\Delta t} \right) \quad (3.37)$$

The end result is that

$$\mathbf{v}_a^+ = \frac{\mathbf{p}_a^{+'}}{m_a} = \frac{1}{m_i} (\mathbf{p}_i + \mathbf{f}_i \Delta t) \quad \text{and} \quad \mathbf{a}_a^+ = \frac{\mathbf{f}'_a}{m_a} = \frac{\mathbf{f}_i}{m_i} \quad (3.38)$$

In other words the material velocity and acceleration both match the single-material mode velocity and acceleration. Any other handling of momenta and force would cause a mismatch between velocity and/or acceleration.

### 3.3.2 Force-Based Contact Calculations

Some early MPM contact claims improved results by looking at contact force rather than a momentum change, but the approach can be shown identical to above momentum change as long as one is careful to synchronize changes in momentum with changes in force. The approach is to consider velocity updates of the two velocity fields as

$$\mathbf{v}_a = \frac{\mathbf{p}_a + (\mathbf{f}_a + \mathbf{f}_c) \Delta t}{m_a} \quad \text{and} \quad \mathbf{v}_b = \frac{\mathbf{p}_b + (\mathbf{f}_b - \mathbf{f}_c) \Delta t}{m_b} \quad (3.39)$$

where  $\mathbf{f}_c$  is contact force needed to impose contact mechanics. Equating the final velocities in the normal direction (same equations apply if equating velocities in tangential direction or in any direction defined by a unit vector) is easily solved to be:

$$\mathbf{f}_c = \left( \frac{m_a (\mathbf{p}_b + \mathbf{f}_b \Delta t) - m_b (\mathbf{p}_a + \mathbf{f}_a \Delta t)}{(m_a + m_b) \Delta t} \cdot \hat{\mathbf{n}} \right) \hat{\mathbf{n}} \quad (3.40)$$

$$\mathbf{f}_c = \left( \frac{\Delta \mathbf{p}_a}{\Delta t} \cdot \hat{\mathbf{n}} + \Delta \mathbf{f}_a \cdot \hat{\mathbf{n}} \right) \hat{\mathbf{n}} \quad (3.41)$$

But realizing that updated momentum before this calculation (*e.g.*,  $\mathbf{p}'_a = \mathbf{p}_a + \mathbf{f}_a \Delta t$ ), this result can be written as:

$$\mathbf{f}_c \Delta t = \left( \frac{m_a \mathbf{p}'_b - m_b \mathbf{p}'_a}{(m_a + m_b)} \cdot \hat{\mathbf{n}} \right) \hat{\mathbf{n}} = (\Delta \mathbf{p}_a \cdot \hat{\mathbf{n}}) \hat{\mathbf{n}} \quad (3.42)$$

where last form is using the updated momenta. This result is identical to the momentum method in previous section for stick in normal direction if you correctly add to nodal force as well as nodal momentum.

### 3.3.3 Low Mass Situations

The above algorithm never divides by a possibly low mass, but that does not necessarily mean it is not affected by one of the masses being very low. In GIMP calculations it will be common for a node one element away from the contact region to have nearly all mass in one material and a very low mass in the other material. When that occurs:

$$\text{if } m_a \ll m_b, \text{ then } \Delta \mathbf{p}_a \approx m_a \mathbf{v}_b - \mathbf{p}_a \quad \text{or} \quad \text{if } m_b \ll m_a, \text{ then } \Delta \mathbf{p}_a \approx \mathbf{p}_b - m_b \mathbf{v}_a \quad (3.43)$$

In other words,  $\Delta \mathbf{p}_a$  will be based on the unreliable nodal momentum (since it is likely to be from one particle on the edge of a material). Since  $\Delta \mathbf{p}_a$  is used to determine contact and to adjust the momenta, the entire algorithm will be unreliable. Now, since  $\Delta \mathbf{p}_a$  will also be very small, it might not make a difference, but in one calculation sliding down an incline plane, the calculation stopped when accepting all contact situations and was fixed by screening out those with  $m_a/(m_a + m_n)$  or  $m_b/(m_a + m_n)$  less than  $10^{-5}$  (might change in the future). Thus in step one above, if the mass ratio meets these criteria, ignore contact for that node.

### 3.3.4 Rigid Material Contact

Specify normals changes and add rigid bias option.

If material  $b$  is rigid, contact can be handled by a special case of the two-material contact, A rigid material corresponds to  $m_b \rightarrow \infty$ . The basic contact equations change to

$$\mathbf{v}_c = \mathbf{v}_b \quad (3.44)$$

$$\Delta \mathbf{p}_a = -m_a(\mathbf{v}_a - \mathbf{v}_b) \quad (3.45)$$

$$\Delta \mathbf{v} = \mathbf{v}_b - \mathbf{v}_a = \frac{\Delta \mathbf{p}_a}{m_a} \quad (3.46)$$

$$m_{red} = m_a \quad (3.47)$$

After these changes, the contact methods above can be applied the same way except the momentum of the rigid particle is not changed. In `NairnMPM`, the density of rigid materials is set to  $1000 \text{ g/cm}^3$  such that material point mass found in `NairnMPM::PreliminaryCalcs()` is equal to the volume for the rigid material point in  $\text{mm}^3$ . Thus the mass gradient needed in the contact algorithm is equal to the volume gradient. The nodal mass, however, for material velocity field of rigid particles is set to zero as a way to determine a rigid particle velocity field. The nodal  $\mathbf{v}_k$  is always the correct velocity and is calculated when the first rigid particle is extrapolated to that node.

### 3.3.5 Three of More Materials

Several options. Lumping seems best. Need more detail.





## Chapter 4

# Imperfect Interface Calculations

### 4.1 Interface Forces

The displacement at node  $i$  for materials  $a$  and  $b$  in crack field  $c$  after current time step can be written as:

$$\mathbf{u}_a(\Delta t) = \mathbf{u}_a(0) + \frac{\mathbf{p}_a \Delta t}{m_a} + \frac{(\mathbf{f}_a + \mathbf{f}_a^{(INT)})(\Delta t)^2}{2m_a} \quad (4.1)$$

$$\mathbf{u}_b(\Delta t) = \mathbf{u}_b(0) + \frac{\mathbf{p}_b \Delta t}{m_b} + \frac{(\mathbf{f}_b - \mathbf{f}_a^{(INT)})(\Delta t)^2}{2m_b} \quad (4.2)$$

where node  $i$  and crack field  $c$  have been omitted (for clarity) and  $\mathbf{f}_a^{(INT)}$  is force that must be added to  $a$  and subtracted from  $b$  to impose interface conditions. Here  $\mathbf{f}_a$  is force found in the current time step. When interface calculations are done after the momentum update, the initial momentum needed in above expression is found from updated momentum ( $\mathbf{p}'_a$ ) using:

$$\mathbf{p}_a = \mathbf{p}'_a - \mathbf{f}_a \Delta t \quad \text{and} \quad \mathbf{p}_b = \mathbf{p}'_b - \mathbf{f}_b \Delta t \quad (4.3)$$

The same holds for material  $b$ . Alternatively, displacement could be written in terms of fully updated momenta as:

$$\mathbf{u}_a(\Delta t) = \mathbf{u}_a(0) + \frac{\mathbf{p}''_a \Delta t}{m_a} - \frac{(\mathbf{f}_a + \mathbf{f}_a^{(INT)})(\Delta t)^2}{2m_a} \quad (4.4)$$

$$\mathbf{u}_b(\Delta t) = \mathbf{u}_b(0) + \frac{\mathbf{p}''_b \Delta t}{m_b} - \frac{(\mathbf{f}_b - \mathbf{f}_a^{(INT)})(\Delta t)^2}{2m_b} \quad (4.5)$$

where

$$\mathbf{p}''_a = \mathbf{p}'_a + \mathbf{f}_a^{(INT)} \Delta t = \mathbf{p}_a + (\mathbf{f}_a + \mathbf{f}_a^{(INT)}) \Delta t \quad (4.6)$$

$$\mathbf{p}''_b = \mathbf{p}'_b - \mathbf{f}_a^{(INT)} \Delta t = \mathbf{p}_b + (\mathbf{f}_b - \mathbf{f}_a^{(INT)}) \Delta t \quad (4.7)$$

or partially updated momenta as

$$\mathbf{u}_a(\Delta t) = \mathbf{u}_a(0) + \frac{\mathbf{p}'_a \Delta t}{m_a} - \frac{(\mathbf{f}_a - \mathbf{f}_a^{(INT)})(\Delta t)^2}{2m_a} \quad (4.8)$$

$$\mathbf{u}_b(\Delta t) = \mathbf{u}_b(0) + \frac{\mathbf{p}'_b \Delta t}{m_b} - \frac{(\mathbf{f}_b + \mathbf{f}_a^{(INT)})(\Delta t)^2}{2m_b} \quad (4.9)$$

Any above approach is correct; these notes use the first.

For interface modeling, we need to follow the displacement discontinuity or  $\delta(\Delta t) = \mathbf{u}_b(\Delta t) - \mathbf{u}_a(\Delta t)$ :

$$\delta(\Delta t) = \delta(0) + \frac{\Delta \mathbf{p}_a \Delta t}{m_{red}} + \frac{\Delta \mathbf{f}_a(\Delta t)^2}{2m_{red}} - \frac{\mathbf{f}_a^{(INT)}(\Delta t)^2}{2m_{red}} \quad (4.10)$$

where

$$\Delta \mathbf{p}_a = \frac{m_a \mathbf{p}_b - m_b \mathbf{p}_a}{m_a + m_b}, \quad \Delta \mathbf{f}_a = \frac{m_a \mathbf{f}_b - m_b \mathbf{f}_a}{m_a + m_b}, \quad \text{and} \quad m_{red} = \frac{m_a m_b}{m_a + m_b} \quad (4.11)$$

But the displacement discontinuity condition must also follow an equation of motion determined by interfacial traction law that superposes the forces found in MPM with interfacial forces, or:

$$\delta(t) = \delta(0) + \frac{\Delta \mathbf{p}_a t}{m_{red}} + \frac{\Delta \mathbf{f}_a t^2}{2m_{red}} - \frac{1}{m_{red}} \int_0^t dt_2 \int_0^{t_2} dt_1 \mathbf{F}_{int}(\delta(t_1)) \quad (4.12)$$

where the interfacial force is

$$\mathbf{F}_{int}(\delta(t)) = F_n(\delta(t)) \hat{\mathbf{n}} + F_t(\delta(t)) \hat{\mathbf{t}} \quad (4.13)$$

where  $F_n(\delta(t))$  and  $F_t(\delta(t))$  are normal and tangential imperfect interface laws. Equating the two interfacial discontinuity equations at  $t = \Delta t$ , the interfacial force we must add (and subtract) becomes

$$\mathbf{f}_a^{(INT)} = \frac{2}{(\Delta t)^2} \int_0^{\Delta t} dt_2 \int_0^{t_2} dt_1 \mathbf{F}_{int}(\delta(t_1)) \quad (4.14)$$

The integral term is found by solving the interfacial equations of motion. The equations for normal and tangential motion are:

$$\delta_n(t) = \delta_n(0) + \frac{d_n t}{m_{red}} + \frac{f_n t^2}{2m_{red}} - \frac{1}{m_{red}} \int_0^t dt_2 \int_0^{t_2} dt_1 F_n(\delta(t_1)) \quad (4.15)$$

$$\delta_t(t) = \delta_t(0) + \frac{d_t t}{m_{red}} + \frac{f_t t^2}{2m_{red}} - \frac{1}{m_{red}} \int_0^t dt_2 \int_0^{t_2} dt_1 F_t(\delta(t_1)) \quad (4.16)$$

where  $d_n = \Delta \mathbf{p}_a \cdot \hat{\mathbf{n}}$ ,  $d_t = \Delta \mathbf{p}_a \cdot \hat{\mathbf{t}}$ ,  $f_n = \Delta \mathbf{f}_a \cdot \hat{\mathbf{n}}$ , and  $f_t = \Delta \mathbf{f}_a \cdot \hat{\mathbf{t}}$ . These equations provide integral equations to be solved (for given interface law) to find  $\delta_n(t)$  and  $\delta_t(t)$ . The analysis can be converted to differential equations:

$$\frac{d^2 \delta_n(t)}{dt^2} + \frac{F_n(\delta(t))}{m_{red}} = \frac{f_n}{m_{red}} \quad \text{and} \quad \frac{d^2 \delta_t(t)}{dt^2} + \frac{F_t(\delta(t))}{m_{red}} = \frac{f_t}{m_{red}} \quad (4.17)$$

Once these are solved, the solutions are used to find an *effective* interface force that must be added such that:

$$\delta_n(\Delta t) = \delta_n(0) + \frac{d_n \Delta t}{m_{red}} + \frac{f_n(\Delta t)^2}{2m_{red}} - \frac{f_{n,a}^{(INT)}(\Delta t)^2}{2m_{red}} \quad (4.18)$$

where the effective force is found from

$$f_{n,a}^{(INT)} = f_n + \frac{2}{\Delta t} \left( \frac{m_{red}(\delta_n(0) - \delta_n(\Delta t))}{\Delta t} + d_n \right) \quad (4.19)$$

A similar equation is used for tangential force as well. The total force added to material  $a$  is then

$$\mathbf{f}_{i,c;a}^{(INT)} = f_{n,a}^{(INT)} \hat{\mathbf{n}} + f_{t,a}^{(INT)} \hat{\mathbf{t}} \quad (4.20)$$

And equal an opposite force is applied to material  $b$ . The equations above are coupled, but in same laws they would decouple if laws can be expressed as  $F_n(\delta_n(t))$  and  $F_t(\delta_t(t))$  where  $\delta_n = \boldsymbol{\delta} \cdot \hat{\mathbf{n}}$  and  $\delta_t = \boldsymbol{\delta} \cdot \hat{\mathbf{t}}$  are normal and tangential components of the interface separation. In other words, the decouple when normal force depends only on normal displacement and tangential force depends only on tangential displacement. Most of the following notes assume this decoupling and therefore might need revision if they do not decouple.

## 4.2 Interfacial Energy

For calculations of composite properties, it is useful to track interfacial energy. Two approaches are to calculate total interfacial energy on each time step. This approach only works for elastic interfaces (need not be linear). The second approach is to find incremental interfacial energy.

The total interfacial energy is:

$$E_{int} = \int_0^{\delta_n} F_n(\delta) d\delta + \int_0^{\delta_t} F_t(\delta) d\delta \quad (4.21)$$

The incremental interfacial energy for one component

$$\Delta E_{int} = \int_0^{\Delta t} F_n(\delta_n(t)) \frac{d\delta_n(t)}{dt} dt + \int_0^{\Delta t} F_t(\delta_t(t)) \frac{d\delta_t(t)}{dt} dt \quad (4.22)$$

The derivative is found after solving for  $\delta_n(t)$

## 4.3 Linear Imperfect Interface

For a linear imperfect interface, such as tangential component of linear interface material is `OSParticulas`, the differential equations can be solved in closed form. The tangential interface law is  $F_t(\delta_t(t)) = D_t A \delta_t(t)$ , where  $D_t$  is imperfect interface parameter and  $A$  is contact area. The differential equation and its solution are:

$$\frac{d^2 \delta_t(t)}{dt^2} + \frac{D_t A \delta_t(t)}{m_{red}} = \frac{f_t}{m_{red}} \quad (4.23)$$

$$\delta_t(t) = \frac{f_t}{D_t A} + C_1 \sin \kappa_t t + C_2 \cos \kappa_t t \quad (4.24)$$

$$\text{where } \kappa_t = \sqrt{\frac{D_t A}{m_{red}}} \quad (4.25)$$

is frequency of the resulting harmonic oscillator. The initial conditions are

$$\delta_t(0) = C_2 + \frac{f_t}{D_t A} \quad \text{and} \quad \frac{d\delta_t(0)}{dt} = C_1 \kappa_t = \frac{d_t}{m_{red}} \quad (4.26)$$

The final solution becomes

$$\delta_t(t) = \frac{d_t}{m_{red} \kappa_t} \sin \kappa_t t + \left( \delta_t(0) - \frac{f_t}{D_t A} \right) \cos \kappa_t t + \frac{f_t}{D_t A} \quad (4.27)$$

$$= \frac{d_t}{m_{red} \kappa_t} \sin \kappa_t t + \left( \delta_t(0) - \frac{f_t}{m_{red} \kappa_t^2} \right) \cos \kappa_t t + \frac{f_t}{m_{red} \kappa_t^2} \quad (4.28)$$

The effective force is found using:

$$\delta_t(\Delta t) = \frac{d_t}{m_{red}\kappa_t} \sin \kappa_t \Delta t + \delta_t(0) \cos \kappa_t \Delta t + \frac{f_t}{m_{red}\kappa_t^2} (1 - \cos \kappa_t \Delta t) \quad (4.29)$$

which leads to

$$f_{t,a}^{(INT)} = f_t + \frac{2}{\Delta t} \left( \frac{m_{red}}{\Delta t} (\delta_t(0) - \delta_t(\Delta t)) + d_t \right) \quad (4.30)$$

$$= f_t + \frac{2}{\Delta t} \left( d_t \left( 1 - \frac{\sin \kappa_t \Delta t}{\kappa_t \Delta t} \right) + \left( \frac{m_{red} \delta_t(0)}{\Delta t} - \frac{f_t}{\kappa_t^2 \Delta t} \right) (1 - \cos \kappa_t \Delta t) \right) \quad (4.31)$$

$$= \frac{2d_t}{\Delta t} \left( 1 - \frac{\sin \kappa_t \Delta t}{\kappa_t \Delta t} \right) + \frac{2m_{red} \delta_t(0) (1 - \cos \kappa_t \Delta t)}{(\Delta t)^2} + f_t \left( 1 - \frac{2(1 - \cos \kappa_t \Delta t)}{\kappa_t^2 (\Delta t)^2} \right) \quad (4.32)$$

As  $D_t \rightarrow 0$  or  $\kappa_t \rightarrow 0$  and using L'Hopital's rule:

$$\lim_{\kappa_t \rightarrow 0} \frac{\sin \kappa_t \Delta t}{\kappa_t \Delta t} = \lim_{\kappa_t \rightarrow 0} \frac{\Delta t \cos \kappa_t \Delta t}{\Delta t} = 1 \quad (4.33)$$

$$\lim_{\kappa_t \rightarrow 0} \frac{2(1 - \cos \kappa_t \Delta t)}{\kappa_t^2 (\Delta t)^2} = \lim_{\kappa_t \rightarrow 0} \frac{2\Delta t \sin \kappa_t \Delta t}{2\kappa_t (\Delta t)^2} = 1 \quad (4.34)$$

which leads to  $f_{t,a}^{(INT)} \rightarrow 0$  as expected for a debonded interface. But as  $D_t \rightarrow \infty$ ,  $f_{t,a}^{(INT)}$  oscillates at increasing frequency. The issue is that MPM analysis cannot handle frequencies beyond its current resolution. Some calculations suggested stability requires  $\kappa_t \Delta t < 1$  (or something on that order and certainly must have  $\kappa_t \Delta t < \pi/2$ ).

This observed stability condition is

$$\Delta t < \frac{1}{\kappa_t} = \sqrt{\frac{m_{red}}{D_t A}} \quad (4.35)$$

These values vary in a simulation, but maximum  $m_{red} \sim m_i/2 = \rho V_i/2 = \rho \Delta x^3/2$  and  $A \sim \Delta x^2$ . Thus stability requires

$$\Delta t < \sqrt{\frac{\rho \Delta x^3}{2D_t \Delta x^2}} = \Delta x \sqrt{\frac{\rho}{2D_t \Delta x}} = C_{int} \frac{\Delta x}{v_{int}} \quad (4.36)$$

where  $C_{int}$  is a CFL factor for interfaces and  $v_{int}$  is an effective wave speed for an interface and given by:

$$v_{int} = \sqrt{\frac{D_t \Delta x}{\rho}} \quad (4.37)$$

Here  $\rho$  would be minimum density of the materials in contact. This velocity depends on grid spacing. Imagine simulation with time step  $\Delta t = C \Delta x/v$  (where  $C$  is CFL factor and  $v$  is maximum material wave speed), the maximum  $D_t$  that can be simulated would be

$$v_{int} < \frac{C_{int} v}{C}, \quad \sqrt{\frac{D_t \Delta x}{\rho}} < \frac{C_{int} v}{C}, \quad \text{or} \quad D_t < \left( \frac{C_{int}}{C} \right)^2 \frac{\rho v^2}{\Delta x} \quad (4.38)$$

Thus, to simulate high  $D_t$  for given materials ( $\rho$  and  $v$ ) and an acceptable interface CFL ( $C_{int}$ ), the simulation needs to have sufficiently low  $\Delta x$  and/or  $C$ . To allow any  $D_t$  in a simulation, each calculation can check  $\kappa_t \Delta t$ . If it is larger than 1 (or maybe up to  $\pi/2$ ), the calculation can revert to stick conditions. In other words, the calculation is converted to stick contact (which is like infinite

$D_t$ ). If the conversion to stick occurs when  $D_t$  has not reached the stick limit, a better simulation would need to increase resolution or decrease time step. A simulation could pre-screen interface laws before starting a calculation, but if that screening was based on  $m_{red} \sim m_i/2$ , the simulation should still become unstable at node with smaller  $m_{red}$ . It is better to check stability on each interface calculation.

#### 4.3.1 Linear Interfacial Energy

The total interfacial energy is

$$E_{int} = \int_0^{\delta_t(\Delta t)} D_t A x dx = \frac{1}{2} D_t A (\delta_t(\Delta t))^2 \quad (4.39)$$

For incremental energy, the results are:

$$\Delta E_{int} = D_t A \int_0^{\Delta t} \delta_t(t) \frac{d\delta_t(t)}{dt} dt = \frac{1}{2} D_t A \int_0^{\Delta t} \frac{d(\delta_t(t))^2}{dt} dt \quad (4.40)$$

$$= \frac{1}{2} D_t A (\delta_t(t))^2 \Big|_0^{\Delta t} = \frac{1}{2} D_t A ((\delta_t(\Delta t))^2 - (\delta_t(0))^2) \quad (4.41)$$

#### 4.3.2 Linear interface Summary and Implementation

Summarizing linear interfaces, set  $\phi = \kappa_t \Delta t$  (dimensionless) and  $m = m_{red}/\Delta t$  (mass/time), then:

$$f_{t,a}^{(INT)} = \frac{2m\delta_t(0)(1 - \cos \phi)}{\Delta t} + \frac{2d_t}{\Delta t} \left(1 - \frac{\sin \phi}{\phi}\right) + f_t \left(1 - \frac{2(1 - \cos \phi)}{\phi^2}\right) \quad (4.42)$$

$$\delta_t(\Delta t) = \delta_t(0) \cos \phi + \frac{d_t}{m} \frac{\sin \phi}{\phi} + \frac{f_t \Delta t (1 - \cos \phi)}{m \phi^2} \quad (4.43)$$

The final discontinuity can be used in either total or incremental energies above. When calculations are done after updating the momenta:

$$d_t = d'_t - f_t \Delta t \quad \text{where} \quad d'_t = \Delta \mathbf{p}'_a \cdot \hat{\mathbf{t}} \quad (4.44)$$

and  $\Delta \mathbf{p}'_a$  is the updated momentum term. The force and displacement become:

$$f_{t,a}^{(INT)} = \frac{2m\delta_t(0)(1 - \cos \phi)}{\Delta t} + \frac{2d'_t}{\Delta t} \left(1 - \frac{\sin \phi}{\phi}\right) + f_t \left(\frac{2 \sin \phi}{\phi} - \frac{2(1 - \cos \phi)}{\phi^2} - 1\right) \quad (4.45)$$

$$\delta_t(\Delta t) = \delta_t(0) \cos \phi + \frac{d'_t}{m} \frac{\sin \phi}{\phi} + \frac{f_t \Delta t}{m} \left(\frac{1 - \cos \phi}{\phi^2} - \frac{\sin \phi}{\phi}\right) \quad (4.46)$$

The options for implementation are:

1. The interface calculations can be done after the mass and momentum calculation, which is before forces are known. This calculation can use the above analysis with  $f_t = 0$ . This approach is correct to first order in time step. It has been used in `OSParticulas` with good results.
2. Alternatively, the calculations can be done after the momentum update and then add the second order terms involving  $f_t$ . New calculations suggest the post-update method is better by a small amount, which corresponds to it being a second order calculation of interfacial force. It is the current `OSParticulas` method.

3. Attempts to combine the first two (*i.e.*, update momenta by interfacial force in mass and momentum update for input to constitutive laws, and then again after momentum and add energy calculations) gives very poor results. It is not clear why (because first phase momenta are restored prior to update), but it is a very large effect (and poor results). In other words, use method 1 or 2, but not some combination of the two methods.
4. Note that contact calculations, which include crack and multimaterial nodes modeling imperfect interfaces, are done two or three times per time step. For interfaces, each of these calculations takes momentum change to stick as input. If the interface can find stable normal or tangential force, the component of momentum change in those direction are set to zero. If stability requires conversion to stick or debond, the momentum change is set to model that condition and interface forces are set to zero. Interfacial forces, when determined, are applied to the node only when contact is done after updating momenta. In other words, interface calculations are superposition of interfaces with forces and interfaces acting like stick or debonded contact. They are always one or the other and never both.
5. The toggling between stick (or debonded) and interface forces is essential for stiff interfaces or for when  $\phi = \kappa_t \Delta t$  is too large. This condition arises often is interfacial stiffness is too high and at nodes with very low  $m_{red}$  (even when interfacial stiffness is not too high). Interface calculations must therefore screen out all high frequency interfaces by switching to stick conditions whenever  $\phi > \phi_{max}$ . The results do not appear very sensitive to cutoff for values of  $\phi_{max} < \pi/2$ , although  $\phi_{max} = \pi$  shows instabilities. Current code uses  $\phi < \pi/2$ . When  $\phi > \phi_{max}$ , the interface uses stick conditions (by leaving stick momentum in the input  $\Delta \mathbf{p}_a$  and by setting interface force and energy to zero). For non-linear interfaces, unstable conditions could be converted to a debonded interfaces if it is softening, but that is not relevant to linear interfaces.
6. Incremental energy calculations are very bad. Again, the reasons are not certain but perhaps are caused by uncertainty in finding initial separation. In other words, subtracting  $\delta_t(0)^2$  does not compensate enough for actual initial energy and accumulation of errors grow with time step. The calculations with the two layer example are probably a worst-case problem because interfacial energy is small overall.

### 4.3.3 Compare to Old Method

The interface force in my paper was calculated using only the initial discontinuity or:

$$f_{t,a}^{(INT,0)} = D_t A \delta_t(0) = m_{red} \kappa_t^2 \delta_t(0) = \frac{m \delta_t(0) \phi^2}{\Delta t} \quad (4.47)$$

This is identical to the  $\delta(0)$  term in harmonic oscillator analysis by expanding  $(1 - \cos \phi)$  to two terms. The old method limited forces to value equal to mean of high-frequency forces, but it may not be best way to limit forces. The limit based on  $\phi$  appears better.

If above is considered zeroth order, maybe improved result, and one that would work for non-linear laws, could be to use the zeroth order force to estimate change in discontinuity during the time step. Substituting the force:

$$\delta_t(t) = \delta_t(0) \left( 1 - \frac{\phi^2 t^2}{(\Delta t)^2} \right) + \frac{d_t t}{m_{red}} + \frac{f_t t^2}{2m_{red}} \quad (4.48)$$

The mean displacement during the time step is:

$$\frac{1}{\Delta t} \int_0^{\Delta t} \delta_t(t) dt = \delta_t(0) \left(1 - \frac{\phi^2}{6}\right) + \frac{d_t}{2m} + \frac{f_t \Delta t}{6m} \quad (4.49)$$

A force based on mean displacement is:

$$f_{t,a}^{(INT,1)} = \frac{2m\delta_t(0)}{\Delta t} \left(\frac{\phi^2}{2} - \frac{\phi^4}{12}\right) + \frac{2d_t}{\Delta t} \left(\frac{\phi^2}{4}\right) + f_t \left(\frac{\phi^2}{6}\right) \quad (4.50)$$

A result that plots better is

$$f_{t,a}^{(INT,1')} = \frac{f_{t,a}^{(INT,0)} + f_{t,a}^{(INT,1)}}{2} = \frac{2m\delta_t(0)}{\Delta t} \left(\frac{\phi^2}{2} - \frac{\phi^4}{24}\right) + \frac{2d_t}{\Delta t} \left(\frac{\phi^2}{8}\right) + f_t \left(\frac{\phi^2}{12}\right) \quad (4.51)$$

Using updated force, the displacement is

$$\delta_t(t) = \delta_t(0) \left(1 - \frac{t^2}{(\Delta t)^2} \left(\frac{\phi^2}{2} - \frac{\phi^4}{24}\right)\right) + \frac{d_t t}{m_{red}} \left(1 - \frac{\phi^2 t}{8\Delta t}\right) + \frac{f_t t^2}{2m_{red}} \left(1 - \frac{\phi^2}{12}\right) \quad (4.52)$$

I tried another round, but it did not improve the results. The above first order looks like a good option to full solution.

Figure 4.1 compares first order force,  $f_{t,a}^{(INT,1')}$ , to full method found by solving differential equation over the full range of  $\phi$  allowed for stability.. The curves give interface force for several relative values of  $\delta_t(0)$ ,  $d_t$ , and  $f_t$ . For a wide range of values, the approximate method is always close to the full harmonic oscillator solution. The two solutions are virtually identical for  $d_t$ ,  $f_t \ll \delta_t(0)$  (*i.e.*, the (0,0) curves). The solutions are also always close when  $\phi \sim 0$  as well (*i.e.*, compliant interface with low interface forces). Note that old zeroth method is the dotted line. It falls within all data, is good for compliant interface and for  $d_t$ ,  $f_t \ll \delta_t(0)$  provided  $\phi < 1$ .

The final results for implementation (based on updated momenta and using position at end from final force) are:

$$f_{t,a}^{(INT,1')} = \frac{m\delta_n(0)\phi^2}{\Delta t} \left(1 - \frac{\phi^2}{12}\right) + \frac{d'_t}{\Delta t} \left(\frac{\phi^2}{4}\right) - f_t \left(\frac{\phi^2}{6}\right) \quad (4.53)$$

$$\delta_t(\Delta t) = \delta_t(0) \left(1 - \frac{\phi^2}{2} \left(1 - \frac{\phi^2}{12}\right)\right) + \frac{d'_t}{m} \left(1 - \frac{\phi^2}{8}\right) - \frac{f_t \Delta t}{m} \left(\frac{1}{2} - \frac{\phi^2}{12}\right) \quad (4.54)$$

## 4.4 Extension to Non-Linear Laws

The revised old method in previous section is OK, but could just use full harmonic oscillator analysis instead. For non-linear laws, however, the differential equation probably cannot be solved and the method of the previous section could guide an approximate method. Repeating the steps for arbitrary  $F_t(\delta(t))$  lead to:

1. Define incremental frequency as:

$$\kappa_t = \sqrt{\frac{|F'_t(\delta_t(0))|}{m_{red}}} \quad \text{and} \quad \phi = \kappa_t \Delta t \quad (4.55)$$

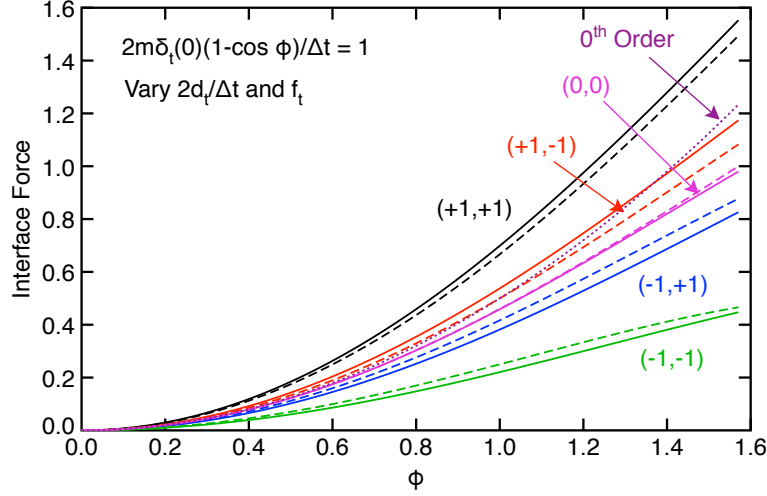


Figure 4.1: Plot of sample force curves for  $2m\delta_t(0)\phi^2/\Delta t = 1$  and various values of  $(2d_t/\Delta t, f_t)$  as indicated. The zeroth order result does not depend on  $d_t$  or  $f_t$  and is therefore a single curve.

If  $\phi > \pi/2$  (or maybe a lower cutoff) proceed with stick conditions. But, if  $F'_t(\delta_t(0)) < 0$  (*i.e.*, law with softening region), the linearized differential equation would involve exponential terms that would blow up for large  $\phi$ . In this case, the physical interpretation is that high frequency corresponds to failure and the interface should be treated as debonded (both force and momentum change set to zero). For low  $\phi$ , should be possible to stably model softening.

2. Find the zeroth order force,  $f_{t,a}^{(INT,0)} = F_t(\delta_t(0))$ , and calculate displacement discontinuity and its mean value:

$$\delta_t(t) = \delta_t(0) - \frac{F_t(\delta_t(0))t^2}{2m_{red}} + \frac{d_t t}{m_{red}} + \frac{f_t t^2}{2m_{red}} \quad (4.56)$$

$$\langle \delta_t^{(1)} \rangle = \delta_t(0) - \frac{F_t(\delta_t(0))\Delta t}{6m} + \frac{d_t}{2m} + \frac{f_t \Delta t}{6m} \quad (4.57)$$

3. Find force based on updated mean displacement

$$f_{t,a}^{(INT,1)} = F_t \left( \langle \delta_t^{(1)} \rangle \right) \quad (4.58)$$

Also check stability by finding  $F'_t \left( \langle \delta_t^{(1)} \rangle \right)$  and switch to stick (or debonding if softening) if needed.

4. Find adjusted force (this combination was based on best approach for linear problem and one that gives exact first two Taylor series terms for the  $\cos \phi$  term):

$$f_{t,a}^{(INT,1')} = \frac{F_t(\delta(0)) + F_t \left( \langle \delta_t^{(1)} \rangle \right)}{2} \quad (4.59)$$

5. Find displacement at end of time step based on adjusted force:

$$\delta_t(\Delta t) = \delta_t(0) - \frac{f_{t,a}^{(INT,1')} \Delta t}{2m} + \frac{d_t}{m} + \frac{f_t \Delta t}{2m} \quad (4.60)$$



Also check stability by finding  $F'_t(\delta_t(\Delta t))$  and switch to stick (or debonding if softening) if needed.

6. Calculate total energy based on final discontinuity (note this approach only allows elastic laws and here would find area under curve up to final displacement discontinuity). Energy is only needed for other calculations. It works for composite analysis, but not anything else.
7. Summarizing and using post-update terms:

$$\langle \delta_t^{(1)} \rangle = \delta_t(0) - \frac{F_t(\delta_t(0))\Delta t}{6m} + \frac{d'_t}{2m} - \frac{f_t\Delta t}{3m} \quad (4.61)$$

$$f_{t,a}^{(INT,1')} = \frac{F_t(\delta(0)) + F_t(\langle \delta_t^{(1)} \rangle)}{2} \quad (4.62)$$

$$\delta_t(\Delta t) = \delta_t(0) - \frac{f_{t,a}^{(INT,1')}\Delta t}{2m} + \frac{d'_t}{m} - \frac{f_t\Delta t}{2m} \quad (4.63)$$

with stability checks done for  $\delta(0)$ ,  $\langle \delta_t^{(1)} \rangle$ , and  $\delta_t(\Delta t)$ , although maybe mean one can be skipped. Implementation could use these equations, or just stope with zeroth order force and displacements (or an option to pick). Class to implement any non-linear law only needs to provide  $F_t(\delta)$ ,  $F'_t(\delta)$ , and methods to check stability (for stick, stable, or debonding) and to find energy.

## 4.5 Bilinear Imperfect Interface

The linear interface contact law in `OSParticulas` is bilinear in the normal direction (and therefore nonlinear) with different interface parameters in tension ( $D_{nt}$ ) and compression ( $D_{nc}$ ). We start with the interfaces initially separated or  $\delta_n(0) > 0$ . The situation evolves as a linear interface and remains that way as long as  $\delta_n(t)$  remains positive. Assuming all linear, final separation is

$$\delta_t(\Delta t) = \frac{d_n}{m_{red}\kappa_{nt}} \sin \kappa_{nt}\Delta t + \left( \delta_n(0) - \frac{f_n}{m_{red}\kappa_{nt}^2} \right) \cos \kappa_{nt}\Delta t + \frac{f_n}{m_{red}\kappa_{nt}^2} \quad (4.64)$$

$$\kappa_{nt} = \sqrt{\frac{D_{nt}A}{m_{red}}} \quad (4.65)$$

If  $\delta_t(\Delta t) > 0$ , use fully linear analysis in previous section (using normal direction parameters). Similarly, if start with  $\delta_n(0) < 0$ , the interface evolves as a linear interface as long as  $\delta_n(t)$  remains negative. The equations are same as above, but use compression stiffness.

$$\kappa_{nc} = \sqrt{\frac{D_{nc}A}{m_{red}}} \quad (4.66)$$

and check if  $\delta_t(\Delta t) < 0$  before using linear analysis.

The above calculations ignores possible vibrations of the interface opening and closing during the time step but ending up the same way it started.

### 4.5.1 Intrastep Contact

If  $\delta_t(\Delta t)$  goes from positive to negative during the time step, the separation becomes zero during the time step at  $0 < t_c < \Delta t$  found by solving

$$0 = \frac{d_n}{m_{red}\kappa_{nt}} \sin \kappa_{nt}t_c + \left( \delta_n(0) - \frac{f_n}{m_{red}\kappa_{nt}^2} \right) \cos \kappa_{nt}t_c + \frac{f_n}{m_{red}\kappa_{nt}^2} \quad (4.67)$$

If  $\delta_t(\Delta t)$  goes from negative to positive during the time step, we solve the same equation, but use  $\kappa_{nc}$ . Two forms useful below are:

$$\delta_n(0) - \frac{f_n}{m_{red}\kappa_{nt}^2} = -\frac{d_n}{m_{red}\kappa_{nt}} \tan \kappa_{nt}t_c - \frac{f_n}{m_{red}\kappa_{nt}^2} \sec \kappa_{nt}t_c \quad (4.68)$$

$$\frac{d_n}{m_{red}\kappa_{nt}} = -\left(\delta_n(0) - \frac{f_n}{m_{red}\kappa_{nt}^2}\right) \cot \kappa_{nt}t_c - \frac{f_n}{m_{red}\kappa_{nt}^2} \csc \kappa_{nt}t_c \quad (4.69)$$

Based on stability arguments, we know that  $0 < \kappa_{nt}t_c < \pi/2$  and therefore both  $\sin \kappa_{nt}t_c$  and  $\cos \kappa_{nt}t_c$  are positive. We can solve (with  $d_n = d'_n - f_n\Delta t$ ) as follows:

$$-\left(\delta_n(0) - \frac{f_n}{m_{red}\kappa_{nt}^2}\right) \cos \kappa_{nt}t_c = \frac{d_n}{m_{red}\kappa_{nt}} \sin \kappa_{nt}t_c + \frac{f_n}{m_{red}\kappa_{nt}^2} \quad (4.70)$$

$$-(m_{red}\kappa_{nt}^2\delta_n(0) - f_n) \cos \kappa_{nt}t_c = \kappa_{nt}d_n \sin \kappa_{nt}t_c + f_n \quad (4.71)$$

$$(m_{red}\kappa_{nt}^2\delta_n(0) - f_n)^2 (1 - \sin^2 \kappa_{nt}t_c) = \kappa_{nt}^2 d_n^2 \sin^2 \kappa_{nt}t_c + 2\kappa_{nt}f_n d_n \sin \kappa_{nt}t_c + f_n^2 \quad (4.72)$$

or a quadratic equation

$$0 = a \sin^2 \kappa_{nt}t_c + b \sin \kappa_{nt}t_c + c \quad (4.73)$$

$$a = (m_{red}\kappa_{nt}^2\delta_n(0) - f_n)^2 + \kappa_{nt}^2 d_n^2 \quad (4.74)$$

$$b = 2\kappa_{nt}f_n d_n \quad (4.75)$$

$$c = f_n^2 - (m_{red}\kappa_{nt}^2\delta_n(0) - f_n)^2 \quad (4.76)$$

The solution is positive term between 0 and  $\Delta t$  from:

$$\sin \kappa_{nt}t_c = \frac{-\kappa_{nt}f_n d_n \pm (m_{red}\kappa_{nt}^2\delta_n(0) - f_n) \sqrt{(m_{red}\kappa_{nt}^2\delta_n(0) - f_n)^2 + \kappa_{nt}^2 d_n^2 - f_n^2}}{(m_{red}\kappa_{nt}^2\delta_n(0) - f_n)^2 + \kappa_{nt}^2 d_n^2} \quad (4.77)$$

$$\cos \kappa_{nt}t_c = \frac{-f_n (m_{red}\kappa_{nt}^2\delta_n(0) - f_n) \mp \kappa_{nt}d_n \sqrt{(m_{red}\kappa_{nt}^2\delta_n(0) - f_n)^2 + \kappa_{nt}^2 d_n^2 - f_n^2}}{(m_{red}\kappa_{nt}^2\delta_n(0) - f_n)^2 + \kappa_{nt}^2 d_n^2} \quad (4.78)$$

In pre-update calculations (when  $f_n = 0$  and therefore  $d_n < 0$  for all contact cases), a simpler solution is

$$\tan \kappa_{nt}t_c = \frac{-m_{red}\kappa_{nt}\delta_n(0)}{d_n} \quad \text{or} \quad t_c = \frac{1}{\kappa_{nt}} \tan^{-1} \left( \frac{-m_{red}\kappa_{nt}\delta_n(0)}{d_n} \right) \quad (4.79)$$

Or from general solution with  $f_n = 0$

$$\sin \kappa_{nt}t_c = \frac{m_{red}\kappa_{nt}\delta_n(0)}{\sqrt{(m_{red}\kappa_{nt}\delta_n(0))^2 + d_n^2}} \quad \text{and} \quad \cos \kappa_{nt}t_c = \frac{-d_n}{\sqrt{(m_{red}\kappa_{nt}\delta_n(0))^2 + d_n^2}} \quad (4.80)$$

#### 4.5.2 Intrastep Contact of Nearly Debonded Interface

The fully debonded interface is considered in section 4.6.1. Here consider contact of nearly debonded interface or Eq. (4.73) when  $\kappa_{nt}$  (or  $\kappa_{nc}$ ) is small. If  $f_n \neq 0$  and  $\sin \kappa_{nt}t_c \approx \kappa_{nt}t_c$  and resulting

constants divided by  $2\kappa_{nt}m_{red}f_n$ , the results reduce to:

$$0 = at_c^2 + bt_c + c \quad (4.81)$$

$$a = \frac{f_n}{2m_{red}} + \left( \frac{d_n^2}{2m_{red}f_n} - \delta_n(0) \left( 1 - \frac{m_{red}\delta_n(0)}{2f_n} \kappa_{nt}^2 \right) \right) \kappa_{nt}^2 \quad (4.82)$$

$$b = \frac{d_n}{m_{red}} \quad (4.83)$$

$$c = \delta_n(0) \left( 1 - \frac{m_{red}\delta_n(0)}{2f_n} \kappa_{nt}^2 \right) \quad (4.84)$$

Neglecting terms  $\kappa_{nt}^2$  and higher and multiplying back by  $2m_{red}$  gives:

$$0 = f_n t_c^2 + 2d_n t_c + 2m_{red}\delta_n(0) \quad (4.85)$$

$$t_c = \frac{-d_n \pm \sqrt{d_n^2 - 2f_n m_{red}\delta_n(0)}}{f_n} \quad (4.86)$$

If almost debonded ( $\kappa_{nt}t_c$  small) and  $f_n = 0$  and constants divided by  $m_{red}^2\kappa_{nt}^4$ , the results reduce to:

$$0 = at_c^2 + c, \quad a = \frac{d_n^2}{m_{red}^2} + \kappa_{nt}^2 \delta_n(0)^2, \quad c = \delta_n(0)^2 \quad (4.87)$$

$$t_c = \frac{|\delta_n(0)|}{\sqrt{\frac{d_n^2}{m_{red}^2} + \kappa_{nt}^2 \delta_n(0)^2}} \quad (4.88)$$

where absolute value are to use this starting opened or closed. Expanding in series and neglecting terms  $\kappa_{nt}^2$  and higher gives:

$$t_c = \frac{m_{red}|\delta_n(0)|}{d_n} \quad (4.89)$$

which is solution to quadratic equation above when  $f_n = 0$ . A stable quadratic equation solver gets both results. These forms are used in code when  $\phi$  is very small. Otherwise,  $t_c$  is found from the full Eq. (4.73).

### 4.5.3 Post Contact with new Interface Parameter

The displacement discontinuity allow for contact change at  $t_c$  within the time step is and relation for zero contact at  $t_c$  are:

$$\begin{aligned} \delta_n(t) = & \delta(0) + \frac{d_n t}{m_{red}} + \frac{f_n t^2}{2m_{red}} - \frac{1}{m_{red}} \int_0^{t_c} dt_2 \int_0^{t_1} dt_1 D_{nt} A \delta_n(t_1) \\ & - \frac{1}{m_{red}} \int_{t_c}^t dt_2 \left( \int_0^{t_c} dt_1 D_{nt} A \delta_n(t_1) + \int_{t_c}^{t_2} dt_1 D_{nc} A \delta_n(t_1) \right) \end{aligned} \quad (4.90)$$

$$\delta_n(t_c) = 0 = \delta(0) + \frac{d_n t_c}{m_{red}} + \frac{f_n t_c^2}{2m_{red}} - \frac{1}{m_{red}} \int_0^{t_c} dt_2 \int_0^{t_1} dt_1 D_{nt} A \delta_n(t_1) \quad (4.91)$$

Note that this post-contact analysis assumes that the interface remains in contact and does not oscillate between opened and closed. That later might be possible due the harmonic oscillations

inherent in final solution. Subtracting the second equation (which is zero) from the first gives:

$$\begin{aligned}
\delta_n(t) &= \frac{d_n(t-t_c)}{m_{red}} + \frac{f_n(t^2-t_c^2)}{2m_{red}} - \frac{1}{m_{red}} \int_{t_c}^t dt_2 \left( \int_0^{t_c} dt_1 D_{nt} A \delta_n(t_1) + \int_{t_c}^{t_2} dt_1 D_{nc} A \delta_n(t_1) \right) \\
&= \frac{d_n(t-t_c)}{m_{red}} + \frac{f_n(t^2-t_c^2)}{2m_{red}} - \frac{t-t_c}{m_{red}} \int_0^{t_c} dt_1 D_{nt} A \delta_n(t_1) - \frac{1}{m_{red}} \int_{t_c}^t dt_2 \int_{t_c}^{t_2} dt_1 D_{nc} A \delta_n(t_1) \\
&= \frac{d_n^{(c)}(t-t_c)}{m_{red}} + \frac{f_n(t-t_c)^2}{2m_{red}} - \frac{1}{m_{red}} \int_{t_c}^t dt_2 \int_{t_c}^{t_2} dt_1 D_{nc} A \delta_n(t_1)
\end{aligned} \tag{4.92}$$

where  $d_n^{(c)}$  is found by equating the last two equations above

$$\frac{d_n^{(c)}}{m_{red}} = \frac{d_n}{m_{red}} + \frac{f_n t_c}{m_{red}} - \frac{1}{m_{red}} \int_0^{t_c} dt_1 D_{nt} A \delta_n(t_1) \tag{4.93}$$

$$\begin{aligned}
&= \frac{d_n}{m_{red}} + \frac{f_n t_c}{m_{red}} - \frac{D_{nt} A}{m_{red} \kappa_{nt}^2} \left( -\frac{d_n}{m_{red}} \cos \kappa_{nt} t_1 \right. \\
&\quad \left. + \left( \delta_n(0) - \frac{f_n}{m_{red} \kappa_{nt}^2} \right) \kappa_{nt} \sin \kappa_{nt} t_1 + \frac{f_n t}{m_{red}} \right) \Big|_0^{t_c}
\end{aligned} \tag{4.94}$$

$$= \frac{d_n}{m_{red}} \cos \kappa_{nt} t_c - \left( \delta_n(0) - \frac{f_n}{m_{red} \kappa_{nt}^2} \right) \kappa_{nt} \sin \kappa_{nt} t_c \tag{4.95}$$

Note that  $d_n^{(c)}$  is also the momentum associated with separation velocity and the time of contact found by differentiating Eq. (4.64) using  $\Delta t = t_c$ :

$$\frac{d_n^{(c)}}{m_{red}} = \frac{d\delta_n(t_c)}{dt} = \frac{d_n}{m_{red}} \cos \kappa_{nt} t_c - \left( \delta_n(0) - \frac{f_n}{m_{red} \kappa_{nt}^2} \right) \kappa_{nt} \sin \kappa_{nt} t_c \tag{4.96}$$

which leads to an identical result. If using post-update terms then

$$\frac{d_n^{(c)}}{m_{red}} = \frac{d'_n - f_n \Delta t}{m_{red}} \cos \kappa_{nt} t_c - \delta_n(0) \kappa_{nt} \sin \kappa_{nt} t_c + \frac{f_n}{m_{red} \kappa_{nt}} \sin \kappa_{nt} t_c \tag{4.97}$$

$$= \frac{d'_n}{m_{red}} \cos \kappa_{nt} t_c - \delta_n(0) \kappa_{nt} \sin \kappa_{nt} t_c + \frac{f_n t_c}{m_{red}} \left( \frac{\sin \kappa_{nt} t_c}{\kappa_{nt} t_c} - \frac{\Delta t}{t_c} \cos \kappa_{nt} t_c \right) \tag{4.98}$$

We next shift the time axis to  $t' = t - t_c$ , to get

$$\delta_n(t') = \frac{d_n^{(c)} t'}{m_{red}} + \frac{f_n t'^2}{2m_{red}} - \frac{1}{m_{red}} \int_0^{t'} dt'_2 \int_0^{t_2} dt'_1 D_{nc} A \delta_n(t'_1) \tag{4.99}$$

The general solution to the differential equation is

$$\delta_n(t') = \frac{f_n}{m_{red} \kappa_{nc}^2} + C_1 \sin \kappa_{nc} t' + C_2 \cos \kappa_{nc} t' \quad \text{where} \quad \kappa_{nc} = \sqrt{\frac{D_{nc} A}{m_{red}}} \tag{4.100}$$

The initial conditions are

$$\delta_n(0) = 0 = \frac{f_n}{m_{red} \kappa_{nc}^2} + C_2 \quad \text{and} \quad \frac{d\delta_n(0)}{dt} = C_1 \kappa_{nc} = \frac{d_n^{(c)}}{m_{red}} \tag{4.101}$$

The final solution becomes

$$\delta_n(t') = \frac{d_n^{(c)}}{m_{red}\kappa_{nc}} \sin \kappa_{nc}t' - \frac{f_n}{m_{red}\kappa_{nc}^2} \cos \kappa_{nc}t' + \frac{f_n}{m_{red}\kappa_{nc}^2} \quad (4.102)$$

$$\delta_n(t) = \frac{d_n^{(c)}}{m_{red}\kappa_{nc}} \sin \kappa_{nc}(t - t_c) - \frac{f_n}{m_{red}\kappa_{nc}^2} \cos \kappa_{nc}(t - t_c) + \frac{f_n}{m_{red}\kappa_{nc}^2} \quad (4.103)$$

Then using Eq. (4.19), the effective interface force for entire step along with the final discontinuity become:

$$f_{n,a}^{(INT)} = \frac{2}{\Delta t} \left( m\delta_n(0) + d_n - d_n^{(c)} \frac{\sin \kappa_{nc}(\Delta t - t_c)}{\kappa_{nc}\Delta t} \right) + f_n \left( 1 - \frac{2(1 - \cos \kappa_{nc}(\Delta t - t_c))}{\kappa_{nc}^2(\Delta t)^2} \right) \quad (4.104)$$

$$\delta_t(\Delta t) = \frac{d_n^{(c)}}{m_{red}} \frac{\sin \kappa_{nc}(\Delta t - t_c)}{\kappa_{nc}} + \frac{f_n}{m_{red}} \left( \frac{1 - \cos \kappa_{nc}(\Delta t - t_c)}{\kappa_{nc}^2} \right) \quad (4.105)$$

These are the equations implements in the linear interface contact low. The next section expands the trigonometric terms for perhaps a more basic result, but the result equations are more complex and less convenient for programming.

In programming, a special case for  $\kappa_{nc}$  is small (needed most when it is zero) neglecting terms  $\kappa_{nc}^2$  and higher is:

$$f_{n,a}^{(INT)} = \frac{2}{\Delta t} \left( m\delta_n(0) + d_n - d_n^{(c)} \left( 1 - \frac{t_c}{\Delta t} \right) \right) + f_n \left( 1 - \left( 1 - \frac{t_c}{\Delta t} \right)^2 \right) \quad (4.106)$$

$$\delta_t(\Delta t) = \frac{d_n^{(c)}}{m_{red}} (\Delta t - t_c) + \frac{f_n}{2m_{red}} (\Delta t - t_c)^2 \quad (4.107)$$

#### 4.5.4 Expand to Initial Time Step Values

Expanding the trigonometric terms:

$$\begin{aligned} \delta_n(t') &= \frac{d_n^{(c)}}{m_{red}\kappa_{nc}} (\sin \kappa_{nc}t \cos \kappa_{nc}t_c - \sin \kappa_{nc}t_c \cos \kappa_{nc}t) \\ &\quad - \frac{f_n}{m_{red}\kappa_{nc}^2} (\cos \kappa_{nc}t \cos \kappa_{nc}t_c + \sin \kappa_{nc}t \sin \kappa_{nc}t_c) + \frac{f_n}{m_{red}\kappa_{nc}^2} \end{aligned} \quad (4.108)$$

which simplifies to:

$$\delta_n(t) = \frac{d_n^{(c)} \cos \kappa_{nc}t_c - \frac{f_n \sin \kappa_{nc}t_c}{\kappa_{nc}}}{m_{red}\kappa_{nc}} \sin \kappa_{nc}t - \frac{d_n^{(c)} \sin \kappa_{nc}t_c + \frac{f_n \cos \kappa_{nc}t_c}{\kappa_{nc}}}{m_{red}\kappa_{nc}} \cos \kappa_{nc}t + \frac{f_n}{m_{red}\kappa_{nc}^2} \quad (4.109)$$

The coefficients simplify with the contact conditions as follows:

$$\begin{aligned} \frac{d_n^{(c)} \cos \kappa_{nc}t_c - \frac{f_n \sin \kappa_{nc}t_c}{\kappa_{nc}}}{m_{red}\kappa_{nc}} &= \left[ \frac{d_n}{m_{red}} \cos \kappa_{nt}t_c - \left( \delta_n(0) - \frac{f_n}{m_{red}\kappa_{nt}^2} \right) \kappa_{nt} \sin \kappa_{nt}t_c \right] \frac{\cos \kappa_{nc}t_c}{\kappa_{nc}} \\ &\quad - \frac{f_n \sin \kappa_{nc}t_c}{m_{red}\kappa_{nc}^2} \end{aligned} \quad (4.110)$$

$$\begin{aligned} &= \left[ \frac{d_n}{m_{red}} \cos \kappa_{nt}t_c + \left( \frac{d_n}{m_{red}} \tan \kappa_{nt}t_c + \frac{f_n}{m_{red}\kappa_{nt}} \sec \kappa_{nt}t_c \right) \sin \kappa_{nt}t_c \right] \frac{\cos \kappa_{nc}t_c}{\kappa_{nc}} - \frac{f_n \sin \kappa_{nc}t_c}{m_{red}\kappa_{nc}^2} \\ &= \frac{1}{m_{red}\kappa_{nc}} \left[ \left( d_n + \frac{f_n \sin \kappa_{nt}t_c}{\kappa_{nt}} \right) \frac{\cos \kappa_{nc}t_c}{\cos \kappa_{nt}t_c} - \frac{f_n \sin \kappa_{nc}t_c}{\kappa_{nc}} \right] \end{aligned} \quad (4.111)$$

and

$$\frac{d_n^{(c)} \sin \kappa_{nc} t_c + \frac{f_n \cos \kappa_{nc} t_c}{\kappa_{nc}}}{m_{red} \kappa_{nc}} = \left[ \frac{d_n}{m_{red}} \cos \kappa_{nt} t_c - \left( \delta_n(0) - \frac{f_n}{m_{red} \kappa_{nt}^2} \right) \kappa_{nt} \sin \kappa_{nt} t_c \right] \frac{\sin \kappa_{nc} t_c}{\kappa_{nc}} + \frac{f_n \cos \kappa_{nc} t_c}{m_{red} \kappa_{nc}^2} \quad (4.112)$$

$$= - \left( \delta_n(0) - \frac{f_n}{m_{red} \kappa_{nt}^2} + \frac{f_n \cos \kappa_{nt} t_c}{m_{red} \kappa_{nt}^2} \right) \frac{\kappa_{nt} \sin \kappa_{nc} t_c}{\kappa_{nc} \sin \kappa_{nt} t_c} + \frac{f_n \cos \kappa_{nc} t_c}{m_{red} \kappa_{nc}^2} \quad (4.113)$$

If the problem was linear ( $\kappa_{nt} = \kappa_{nc} = \kappa_n$ ) then

$$\frac{d_n^{(c)} \cos \kappa_{nc} t_c - \frac{f_n \sin \kappa_{nc} t_c}{\kappa_{nc}}}{m_{red} \kappa_{nc}} = \frac{d_n}{\kappa_n m_{red}} \quad (4.114)$$

$$\frac{d_n^{(c)} \sin \kappa_{nc} t_c + \frac{f_n \cos \kappa_{nc} t_c}{\kappa_{nc}}}{m_{red} \kappa_{nc}} = - \left( \delta_n(0) - \frac{f_n}{m_{red} \kappa_n^2} \right) \quad (4.115)$$

These results recover the linear result as it should.

The net effective force is found from

$$\delta_n(\Delta) = \frac{d_n^{(c)} \cos \kappa_{nc} t_c - \frac{f_n \sin \kappa_{nc} t_c}{\kappa_{nc}}}{m_{red} \kappa_{nc}} \sin \kappa_{nc} \Delta t - \frac{d_n^{(c)} \sin \kappa_{nc} t_c + \frac{f_n \cos \kappa_{nc} t_c}{\kappa_{nc}}}{m_{red} \kappa_{nc}} \cos \kappa_{nc} \Delta t + \frac{f_n}{m_{red} \kappa_{nc}^2} = \delta_n(0) + \frac{d_n \Delta t}{m_{red}} + \frac{f_n (\Delta t)^2}{2 m_{red}} - \frac{f_{n,a}^{(INT)} (\Delta t)^2}{2 m_{red}} \quad (4.116)$$

$$f_{n,a}^{(INT)} = f_n + \frac{2}{\Delta t} \left[ d_n \left( 1 - \frac{d_n^{(c)} \cos \kappa_{nc} t_c - \frac{f_n \sin \kappa_{nc} t_c}{\kappa_{nc}}}{d_n} \frac{\sin \kappa_{nc} \Delta t}{\kappa_{nc} \Delta t} \right) - \frac{f_n}{\kappa_{nc}^2 \Delta t} + \frac{\delta_n(0) m_{red}}{\Delta t} \left( 1 + \frac{d_n^{(c)} \sin \kappa_{nc} t_c + \frac{f_n \cos \kappa_{nc} t_c}{\kappa_{nc}}}{\delta_n(0) m_{red} \kappa_{nc}} \cos \kappa_{nc} \Delta t \right) \right] \quad (4.117)$$

This result reduces to linear result when  $\kappa_{nt} = \kappa_{nc} = \kappa_n$ . Inserting revised coefficients for bilinear gives:

$$f_{n,a}^{(INT)} = f_n + \frac{2}{\Delta t} \left[ \left( d_n - \left[ \left( d_n + \frac{f_n \sin \kappa_{nt} t_c}{\kappa_{nt}} \right) \frac{\cos \kappa_{nc} t_c}{\cos \kappa_{nt} t_c} - \frac{f_n \sin \kappa_{nc} t_c}{\kappa_{nc}} \right] \frac{\sin \kappa_{nc} \Delta t}{\kappa_{nc} \Delta t} \right) - \frac{f_n}{\kappa_{nc}^2 \Delta t} + \left( \frac{m_{red} \delta_n(0)}{\Delta t} - \left[ \left( \frac{m_{red} \delta_n(0)}{\Delta t} - \frac{f_n}{\kappa_{nt}^2 \Delta t} (1 - \cos \kappa_{nt} t_c) \right) \frac{\kappa_{nt} \sin \kappa_{nc} t_c}{\kappa_{nc} \sin \kappa_{nt} t_c} - \frac{f_n \cos \kappa_{nc} t_c}{\kappa_{nc}^2 \Delta t} \right] \cos \kappa_{nc} \Delta t \right) \right] \quad (4.118)$$

In pre-update when  $f_n = 0$ , this results reduces to:

$$f_{n,a}^{(INT)} = \frac{2}{\Delta t} \left[ d_n \left( 1 - \frac{\cos \kappa_{nc} t_c}{\cos \kappa_{nt} t_c} \frac{\sin \kappa_{nc} \Delta t}{\kappa_{nc} \Delta t} \right) + \frac{\delta_n(0) m_{red}}{\Delta t} \left( 1 - \frac{\kappa_{nt} \sin \kappa_{nc} t_c}{\kappa_{nc} \sin \kappa_{nt} t_c} \cos \kappa_{nc} \Delta t \right) \right] \quad (4.119)$$

Get  $f_n$  term by setting  $d_n = \delta_n(0) = 0$  or

$$f_{n,a}^{(INT)} = f_n \left[ 1 - \frac{2}{\Delta t} \left( \frac{\sin \kappa_{nt} t_c}{\kappa_{nt}} \frac{\cos \kappa_{nc} t_c}{\cos \kappa_{nt} t_c} - \frac{\sin \kappa_{nc} t_c}{\kappa_{nc}} \right) \frac{\sin \kappa_{nc} \Delta t}{\kappa_{nc} \Delta t} - \frac{2}{\kappa_{nc}^2 (\Delta t)^2} \left( 1 - \left( (1 - \cos \kappa_{nt} t_c) \frac{\kappa_{nt} \sin \kappa_{nc} t_c}{\kappa_{nc} \sin \kappa_{nt} t_c} + \cos \kappa_{nc} t_c \right) \cos \kappa_{nc} \Delta t \right) \right] \quad (4.120)$$

For energy calculations, we can find

$$\begin{aligned}\delta_n(\Delta) = & \frac{1}{m_{red}\kappa_{nc}} \left[ \left( d_n + \frac{f_n \sin \kappa_{nt} t_c}{\kappa_{nt}} \right) \frac{\cos \kappa_{nc} t_c}{\cos \kappa_{nt} t_c} - \frac{f_n \sin \kappa_{nc} t_c}{\kappa_{nc}} \right] \sin \kappa_{nc} \Delta t + \frac{f_n}{m_{red}\kappa_{nc}^2} \\ & + \left[ \left( \delta_n(0) - \frac{f_n}{m_{red}\kappa_{nt}^2} + \frac{f_n \cos \kappa_{nt} t_c}{m_{red}\kappa_{nt}^2} \right) \frac{\kappa_{nt} \sin \kappa_{nc} t_c}{\kappa_{nc} \sin \kappa_{nt} t_c} - \frac{f_n \cos \kappa_{nc} t_c}{m_{red}\kappa_{nc}^2} \right] \cos \kappa_{nc} \Delta t\end{aligned}\quad (4.121)$$

In pre-update when  $f_n = 0$ , this results reduces to:

$$\delta_n(\Delta t) = \frac{d_n}{m_{red}\kappa_{nc}} \frac{\cos \kappa_{nc} t_c}{\cos \kappa_{nt} t_c} \sin \kappa_{nc} \Delta t + \delta_n(0) \frac{\kappa_{nt} \sin \kappa_{nc} t_c}{\kappa_{nc} \sin \kappa_{nt} t_c} \cos \kappa_{nc} \Delta t \quad (4.122)$$

Get  $f_n$  term by setting  $d_n = \delta_n(0) = 0$  or

$$\begin{aligned}m_{red}\delta_t(\Delta) = & f_n \left[ \left( \frac{\sin \kappa_{nt} t_c}{\kappa_{nt}} \frac{\cos \kappa_{nc} t_c}{\cos \kappa_{nt} t_c} - \frac{\sin \kappa_{nc} t_c}{\kappa_{nc}} \right) \frac{\sin \kappa_{nc} \Delta t}{\kappa_{nc}} \right. \\ & \left. + \frac{1}{\kappa_{nc}^2} \left( 1 - \left( (1 - \cos \kappa_{nt} t_c) \frac{\kappa_{nt} \sin \kappa_{nc} t_c}{\kappa_{nc} \sin \kappa_{nt} t_c} + \cos \kappa_{nc} t_c \right) \cos \kappa_{nc} \Delta t \right) \right]\end{aligned}\quad (4.123)$$

Total energy is found at the end of the time step. An incremental approach needs to look at results before and after the contact time.

$$\Delta E_{int} = \frac{1}{2} D_{nt} A \int_0^{t_c} \frac{d(\delta_t(t))^2}{dt} dt + \frac{1}{2} D_{nc} A \int_{t_c}^{\Delta t} \frac{d(\delta_t(t))^2}{dt} dt \quad (4.124)$$

$$= \frac{1}{2} D_{nt} A (\delta_t(t))^2 \Big|_0^{t_c} + \frac{1}{2} D_{nc} A (\delta_t(t))^2 \Big|_{t_c}^{\Delta t} = \frac{1}{2} A (D_{nc}(\delta_t(\Delta t))^2 - D_{nt}(\delta_t(0))^2) \quad (4.125)$$

Summarizing, the interfacial force with  $\phi = \kappa_{nc} \Delta t$  and  $m = m_{red}/\Delta t$  is

$$\begin{aligned}f_{n,a}^{(INT)} = & \frac{2}{\Delta t} \left[ d_n \left( 1 - \frac{\cos \kappa_{nc} t_c}{\cos \kappa_{nt} t_c} \frac{\sin \phi}{\phi} \right) + m \delta_n(0) \left( 1 - \frac{\kappa_{nt} \sin \kappa_{nc} t_c}{\kappa_{nc} \sin \kappa_{nt} t_c} \cos \phi \right) \right] \\ & + f_n \left[ 1 - \frac{2}{\Delta t} \left( \frac{\sin \kappa_{nt} t_c}{\kappa_{nt}} \frac{\cos \kappa_{nc} t_c}{\cos \kappa_{nt} t_c} - \frac{\sin \kappa_{nc} t_c}{\kappa_{nc}} \right) \frac{\sin \phi}{\phi} \right. \\ & \left. - \frac{2 \left( 1 - \left( (1 - \cos \kappa_{nt} t_c) \frac{\kappa_{nt} \sin \kappa_{nc} t_c}{\kappa_{nc} \sin \kappa_{nt} t_c} + \cos \kappa_{nc} t_c \right) \cos \phi \right)}{\phi^2} \right]\end{aligned}\quad (4.126)$$

and interface separation is

$$\begin{aligned}\delta_n(\Delta t) = & \delta_n(0) \frac{\kappa_{nt} \sin \kappa_{nc} t_c}{\kappa_{nc} \sin \kappa_{nt} t_c} \cos \phi + \frac{d_n}{m} \frac{\cos \kappa_{nc} t_c}{\cos \kappa_{nt} t_c} \frac{\sin \phi}{\phi} \\ & + \frac{f_n}{m} \left[ \left( \frac{\sin \kappa_{nt} t_c}{\kappa_{nt}} \frac{\cos \kappa_{nc} t_c}{\cos \kappa_{nt} t_c} - \frac{\sin \kappa_{nc} t_c}{\kappa_{nc}} \right) \frac{\sin \phi}{\phi} \right. \\ & \left. + \frac{\Delta t \left( 1 - \left( (1 - \cos \kappa_{nt} t_c) \frac{\kappa_{nt} \sin \kappa_{nc} t_c}{\kappa_{nc} \sin \kappa_{nt} t_c} + \cos \kappa_{nc} t_c \right) \cos \phi \right)}{\phi^2} \right]\end{aligned}\quad (4.127)$$

An interface that starts in contact and separates would use same equation but interchange compression and tension terms.

## 4.6 Debonding Interface and Perfect in Compression

### 4.6.1 Debonded Interface

Imagine a debonded interface with no force when separated. If not debonded in compression, it is bilinear. If perfect in compression, it is should be a new approach to contact. We start with finite stiffness in compression and then try to deduce results for perfect interface.

Starting with  $\delta_n(0) > 0$ , the only concern is if contact occurs in the time step. The discontinuity without contact forces is:

$$\delta_n(t) = \delta_n(0) + \frac{d_n t}{m_{red}} + \frac{f_n t^2}{2m_{red}} \quad (4.128)$$

If  $\delta_n(\Delta t) > 0$ , interface remains debonded with zero force and energy for time step. If negative, then

$$\delta_n(0) < -\frac{(d_n + \frac{1}{2}f_n\Delta t)\Delta t}{m_{red}} \quad (4.129)$$

When this applies, contact occurs at:

$$0 = 2m_{red}\delta_n(0) + 2d_n t_c + f_n t_c^2 \quad (4.130)$$

$$t_c = \frac{-d_n \pm \sqrt{d_n^2 - 2m_{red}\delta_n(0)f_n}}{f_n} \quad (4.131)$$

$$= -\frac{d_n}{f_n} \pm \text{sign}(f_n) \sqrt{\left(\frac{d_n}{f_n}\right)^2 - \frac{m_{red}\delta_n(0)}{f_n}} \quad (4.132)$$

The velocity at time of contact is

$$\frac{d\delta_n(t_c)}{dt} = \frac{d_n}{m_{red}} + \frac{f_n t_c}{m_{red}} < 0 \quad \text{or} \quad d_n^{(c)} = d_n + f_n t_c \quad (4.133)$$

These results imply that one or the other of  $d_n$  and  $f_n$  must be negative (or both). The three possibilities are:

$$\begin{aligned} d_n < 0 \text{ and } f_n < 0 &\implies -\frac{d_n}{f_n} < 0 (\text{sign}(f_n) = -1) \implies \text{use } - \text{ root} \\ d_n > 0 \text{ and } f_n < 0 &\implies -\frac{d_n}{f_n} < t_c \implies \text{use } - \text{ root} \\ d_n < 0 \text{ and } f_n > 0 &\implies -\frac{d_n}{f_n} > t_c \implies \text{use } - \text{ root} \end{aligned} \quad (4.134)$$

The contact velocity becomes:

$$m_{red} \frac{d\delta_n(t_c)}{dt} = -\sqrt{d_n^2 - m_{red}\delta_n(0)f_n} \quad (4.135)$$

A special case when  $f_n = 0$ , which implies  $d_n < 0$ , and leads to

$$t_c = -\frac{m_{red}\delta_n(0)}{d_n} \quad \text{and} \quad \frac{d\delta_n(t_c)}{dt} = \frac{d_n}{m_{red}} \quad \text{or} \quad d_n^{(c)} = d_n \quad (4.136)$$

In summary, for all possibilities:

$$t_c = \begin{cases} -\frac{m_{red}\delta_n(0)}{d_n} & f_n = 0 \\ \frac{-d_n \pm \sqrt{d_n^2 - 2m_{red}\delta_n(0)f_n}}{f_n} & \text{otherwise} \end{cases} \quad (4.137)$$

$$d_n^{(c)} = d_n + f_n t_c = d'_n - f_n \Delta t' = -\sqrt{d_n^2 - m_{red}\delta_n(0)f_n} \quad (4.138)$$

where  $\Delta t' = \Delta t - t_c$ . To finish the step, proceed to post-contact methods.



### 4.6.2 Post Contact when Debonded Interface

For finite compression stiffness, the interfacial force and displacement from Eq. (4.104) and previous section become

$$\begin{aligned} f_{n,a}^{(INT)} &= \frac{2}{\Delta t} \left[ m\delta_n(0) + d'_n \left( 1 - \frac{\Delta t' \sin \phi'}{\Delta t \phi'} \right) \right] + f_n \left[ 2 \left( \frac{\Delta t'}{\Delta t} \right)^2 \left( \frac{\sin \phi'}{\phi'} - \frac{(1 - \cos \phi')}{(\phi')^2} \right) - 1 \right] \\ \delta_t(\Delta t) &= \frac{\Delta t'}{\Delta t} \left[ \frac{d'_n \sin \phi'}{m \phi'} + \frac{f_n \Delta t'}{m} \left( \frac{1 - \cos \phi'}{(\phi')^2} - \frac{\sin \phi'}{\phi'} \right) \right] \end{aligned} \quad (4.139)$$

where  $\phi' = \kappa_{nc}\Delta t'$ . This result reduces to linear interface (see Eq. (4.45)) when  $t_c = 0$ ,  $\Delta t' = \Delta t$ ,  $\phi' = \phi$ , and  $\delta_n(0) = 0$  (this last because contact occurred at time zero).

### 4.6.3 Post Contact When Perfect in Compression

If compression is a perfect interface, the contact force would be zero up to the point of contact. We then assume that there is no penetration and that it does not bounce off within this time step. In other words  $\delta_n(t) = 0$  for  $t > t_c$ . For it to remain zero, Eq. (4.92) implies that:

$$\int_0^{\Delta t'} dt_2 \int_0^{t_2} dt_1 F_n(0) = d_n^{(c)} \Delta t' + \frac{f_n (\Delta t')^2}{2} \quad (4.140)$$

where  $\Delta t' = \Delta t - t_c$  or the time interval after contact and hypothetical force is evaluated only at  $\delta_n(t) = 0$ . The final separation during the time step become:

$$\delta_n(t) = \begin{cases} \delta(0) + \frac{d_n t}{m_{red}} + \frac{f_n t^2}{2m_{red}} & t < t_c \\ 0 & t > t_c \end{cases} \quad (4.141)$$

The interfacial force to mimic this evolution in the MPM time step would be

$$f_{n,a}^{(INT)} = f_n + \frac{2}{\Delta t} \left( \frac{m_{red}\delta_n(0)}{\Delta t} + d_n \right) = \frac{2}{\Delta t} \left( \frac{m_{red}\delta_n(0)}{\Delta t} + d'_n \right) - f_n \quad (4.142)$$

This result is also found by taking limit of Eq. (4.139) as  $\phi' \rightarrow \infty$ . This result does not depend on the contact time,  $t_c$ . Note that this could be alternative force for contact if done after the momentum update. The current method finds  $f_{n,a}^{(INT)} \Delta t = d'_n$ . The new momentum change (added to particle momentum and force) would be

$$\Delta \mathbf{p}'_a \cdot \hat{\mathbf{n}} = f_{n,a}^{(INT)} \Delta t = \frac{2m_{red}\delta_n(0)}{\Delta t} + 2d_n + f_n \Delta t = d'_n + d_n + \frac{2m_{red}\delta_n(0)}{\Delta t} \quad (4.143)$$

### 4.6.4 Other?

For interface starting in compression and moving into tension, repeat the above analysis but interchange tension and compression interface properties.

For interface starting perfect ( $D < 0$  in `OSParticulas`, it stays perfect as stick contact). For imperfect interface moving into contact, can apply force to reach contact and then adjust momentum to stick after contact using  $d_n^{(c)}$ . In other words, change input momentum change to the new value at the point of contact and also apply the force.

## 4.7 Use Imperfect Interface to Soften Contact Calculations

Contact calculations treat debonded as having zero stiffness with contact has infinite stiffness. This abrupt change, could be interpreted as a bilinear interface leading to following revised contact algorithm:

1. Find displacement at end of current time step as

$$\delta_n(\Delta t) = \delta(0) + \frac{1}{m} \left( d'_n - \frac{f_n \Delta t}{2} \right) \quad (4.144)$$

where  $m = m_{red}/\Delta t$ . But, if not in contact calculations just after momentum updates, simply set  $\delta_n(\Delta t) = \delta_n(0)$ .

2. Because interface is considered as perfect interface when in contact, we set the interfacial normal force to be

$$f_{n,a}^{(INT)} \Delta t = d'_n \quad (4.145)$$

which is simply  $d_n$  when not just after momentum updates. It is possible force should use perfect interface results instead or

$$f_{n,a}^{(INT)} = 2 \left( \frac{m_{red} \delta_n(0)}{\Delta t} + d'_n \right) - f_n \Delta t \quad (4.146)$$

although some trials with this force did not look as good.

3. Find tangential interface force to stick or force if was perfect interface in shear

$$f_{n,t}^{(INT)} \Delta t = d'_t \quad (4.147)$$

4. If interface has adhesion, then if  $\delta_n(\Delta t) > 0$ , check failure based on traction stresses implied by  $f_{n,a}^{(INT)}$  and  $f_{n,t}^{(INT)}$ . If they say adhesion not failure, the contact treated as stick conditions, but if they say failed, treat as not in contact with no momentum changes.
5. If  $\delta_n(\Delta t) > 0$  the interface is not in contact and no momentum changes are done. If  $\delta_n(\Delta t) < 0$ , the two option are to a) assume in contact, apply  $f_{n,a}^{(INT)}$  to node, and use contact law to find final tangential force. b) check sign of  $f_{n,a}^{(INT)}$ . If positive, assume not in contact, but if negative, assume compression and proceed like option a. Option “a” is the new “displacementOnly” option in code. Note that this option might end up applying tensile force to the node even if no tangential force. Option “b” never applies tensile normal force to the node.

A revised used of imperfect interfaces for revised contact is the find interfacial force and displacement using a finite contact stiffness as an imperfect interface parameter  $D_c$  for linear contact stiffness. The following algorithm only applies in contact calculations just after momentum update when interfacial force is known. First, find  $m = m_{red}/\Delta t$ ,  $d = D_c A_c \Delta t$ , and  $\phi = \sqrt{d/m}$

1. If  $\phi > \pi/2$ , contact is treated as perfect and finds terms using above contact algorithm or:

$$\delta_n(\Delta t) = \delta(0) + \frac{1}{m} \left( d'_n - \frac{f_n \Delta t}{2} \right) \quad \text{and} \quad f_{n,a}^{(INT)} \Delta t = d'_n \quad (4.148)$$

but, if  $\phi < \pi/2$ , find these terms instead

$$\delta_n(\Delta t) = \delta_n(0) \cos \phi + \frac{d'_n}{m} \frac{\sin \phi}{\phi} + \frac{f_n \Delta t}{m} \left( \frac{1 - \cos \phi}{\phi^2} - \frac{\sin \phi}{\phi} \right) \quad (4.149)$$

$$f_{n,a}^{(INT)} \Delta t = 2 [m(\delta_n(0) - \delta_n(\Delta t)) + d'_n] - f_n \Delta t \quad (4.150)$$

2. Proceed to step 3 of previous algorithm. Only  $\delta_n(\Delta t)$  and  $f_{n,a}^{(INT)} \Delta t$  may have changed.
3. Contact calculations done other places can use the previous algorithm with no interface calculations.



## Chapter 5

# MPM Time Step Tasks For Perfect Interface

As a check/evaluation of contact methods, this section contrasts the MPM time step tasks for a single-material mode calculation to a simulation with two identical with contact law that assumes a perfect interface (*e.g.*, the `IgnoreContact` law) and with no cracks.

### 5.1 Simulation and Velocity Field Initialization

Each nodal point is permanently assigned crack field [0]. When in single material velocity field, field [0] is given a single material velocity field [0]. When using multimaterial mode, crack velocity field [0] is assigned pointers to material velocity fields for materials *a* and *b*. The pointers are initialized to `NULL`. Any material velocity fields needed in a simulation are dynamically allocated as needed.

### 5.2 Momentum and Mass Extrapolation to the Grid

In single material mode, there is only one material velocity field with nodal values:

$$m_i = \sum_p S_{ip} m_p \quad \text{and} \quad \mathbf{p}_i = m_i \mathbf{v}_i = \sum_p S_{ip} \mathbf{p}_p \quad (5.1)$$

In multimaterial mode

$$m_{i,j} = \sum_{p \in j} S_{ip} m_p \quad \text{and} \quad \mathbf{p}_{i,j} = m_{i,j} \mathbf{v}_{i,j} = \sum_{p \in j} S_{ip} \mathbf{p}_p \quad (5.2)$$

Because  $j = a$  or  $b$  only here

$$m_i = m_{i,a} + m_{i,b} \quad \text{and} \quad \mathbf{p}_i = \mathbf{p}_{i,a} + \mathbf{p}_{i,b} \quad \text{but} \quad \mathbf{v}_i \neq \mathbf{v}_{i,a} \neq \mathbf{v}_{i,b} \quad (5.3)$$

Multimaterial mode does more extrapolations for position, volume, and volume gradient, but when using `IgnoreContact`, those quantities are not used.

### 5.3 Revised XPIC Task

If using revised XPIC( $m$ ) for  $m > 1$ , the mass and momentum tasks are followed by XPIC tasks that replaces grid velocities with  $\mathbf{v}_i(m) = m(\mathbf{v}_i - \mathbf{v}_i^*)$ . In vector form:

$$\mathbf{v}(m) = \sum_{k=1}^m (-1)^{k+1} \mathbf{v}_k \quad (5.4)$$

where

$$\mathbf{v}_k = \frac{m-k+1}{k} \mathbf{S}^+ \mathbf{S} \mathbf{v}_{k-1} \quad (5.5)$$

starting with

$$\mathbf{v}_1 = m \mathbf{v} \quad (5.6)$$

In single material mode, the explicit sums are:

$$(\mathbf{v}_k)_i = \frac{m-k+1}{k} \sum_p \sum_{j \in \Omega_p} \frac{M_p S_{pi} S_{pj}}{m_i} (\mathbf{v}_{k-1})_j \quad \forall i \in \Omega_p \quad (5.7)$$

In multimaterial mode,  $\mathbf{v}(m)$  will be found in each material velocity field. Explicit sums for one material are:

$$(\mathbf{v}_k)_{i,a} = \frac{m-k+1}{k} \sum_{p \in a} \sum_{j \in \Omega_p} \frac{M_p S_{pi} S_{pj}}{m_{i,a}} (\mathbf{v}_{k-1})_{j,a} \quad \forall i \in \Omega_p \quad (5.8)$$

Single material results when modeling two materials can be written:

$$m_i (\mathbf{v}_k)_i = \frac{m-k+1}{k} \sum_p \sum_{j \in \Omega_p} M_p S_{pi} S_{pj} \frac{(\mathbf{p}_{k-1})_{j,a} + (\mathbf{p}_{k-1})_{j,b}}{m_{j,a} + m_{j,b}} \quad (5.9)$$

Note that prior to XPIC task that

$$m_i \mathbf{v}_i = \mathbf{p}_i = \mathbf{p}_{i,a} + \mathbf{p}_{i,b} = m_{i,a} \mathbf{v}_{i,a} + m_{i,b} \mathbf{v}_{i,b} \quad (5.10)$$

but after replacing grid velocities with  $\mathbf{v}(m)$ , it appears that

$$m_i \mathbf{v}(m)_i \neq m_{i,a} \mathbf{v}(m)_{i,a} + m_{i,b} \mathbf{v}(m)_{i,b} \quad (5.11)$$

If true, multimaterial mode would not be able to reduce to single material mode for the case of a perfect interface.

### 5.4 Post Momentum and Mass Extrapolation Tasks

After the extrapolations are done, each node is revisited for more calculations to impose multimaterial effects (in that mode). The steps are

1. Calculate total mass and number of particles in each velocity field. In the process copy  $\mathbf{p}_i$  and  $\mathbf{p}_{i,j}$  for all material velocity fields on all nodes. Also make a list of nodes that have both materials as list of `MaterialContactNode` objects.

2. In multimaterial mode, change momenta for fields  $a$  and  $b$  by

$$\pm \Delta \mathbf{p}_a = \pm \left( \frac{m_a}{m_i} \mathbf{p}_i - \mathbf{p}_a \right) \quad (5.12)$$

Adding to material  $a$  (and subtracting from  $b$ ) leads to

$$\mathbf{p}'_{i,a} = \frac{m_{i,a}}{m_i} \mathbf{p}_i = m_{i,a} \mathbf{v}_i \quad \text{and} \quad \mathbf{p}'_{i,a} = \frac{m_{i,b}}{m_i} \mathbf{p}_i = m_{i,b} \mathbf{v}_i \quad (5.13)$$

Note that materials  $a$  and  $b$  now have the same velocity ( $= \mathbf{v}_i$ ) but have different momenta. These changes are not done by setting new momenta based of  $\mathbf{v}_i$ , but by using

$$\Delta \mathbf{p}_a = \frac{m_{i,a} \mathbf{p}_i - m_i \mathbf{p}_{i,a}}{m_i} = \frac{m_{i,a} \mathbf{p}_{i,b} - m_{i,b} \mathbf{p}_{i,a}}{m_i} \quad (5.14)$$

The velocity for material  $a$  (for example) would be

$$\mathbf{v}_{i,a} = \frac{\mathbf{p}_{i,a}}{m_{i,a}} + \frac{m_{i,a} \mathbf{p}_{i,b} - m_{i,b} \mathbf{p}_{i,a}}{m_{i,a} m_i} \quad (5.15)$$

Perhaps this equation would lose accuracy when  $m_{i,a} \ll m_{i,b}$ ?

3. Note that  $\mathbf{p}'_{i,a}$  replaces the store momenta (to be restored later) and the force is seeded with

$$\mathbf{f}_{i,a} = - \frac{\Delta \mathbf{p}_a}{\Delta t} \quad (5.16)$$

4. Impose velocity boundary conditions by setting  $\mathbf{v}_i = \mathbf{v}^{(BC)}$ , which which would changed  $a$  and  $b$  in multimaterial mode.

## 5.5 Update Stresses and Strains on Particle

When updating stresses and strains before finding forces and updating momenta (needed for **USAVG** $\pm$  and for **USF** update methods), extrapolate velocity gradient ( $\nabla \mathbf{v}$ ) to the particles and implement the chosen constitutive laws. The extrapolations are:

$$\nabla \mathbf{v}_p = \sum_i \mathbf{v}_i \otimes \mathbf{G}_{ip} \quad \text{and} \quad \nabla \mathbf{v}_{p \in j} = \sum_i \mathbf{v}_{i,j} \otimes \mathbf{G}_{ip} \quad (5.17)$$

Because  $\mathbf{v}_{i,a} = \mathbf{v}_{i,b} = \mathbf{v}_i$  for a perfect interface, both modes should extrapolate the same velocity to the particles resulting in same updates of stress and strain on particles.

## 5.6 Project Forces to the Grid

For each material point, extrapolate  $\mathbf{f}_{i,j}$  due to particle stresses particle forces (later not used often):

$$\mathbf{f}_{i,j} = \sum_{p \in j} \left( -m_p \frac{\boldsymbol{\tau}_p^{(n)} \cdot \mathbf{G}_{ip}^{(n)}}{\rho_0} + \mathbf{F}_p^{(n)} S_{ip}^{(n)} \right) \quad (5.18)$$

Because  $j = a$  or  $b$  only here, and assuming particle stresses remain in synch, we have total nodal force as:

$$\mathbf{f}_i = \mathbf{f}_{i,a} + \mathbf{f}_{i,b} \quad \text{but} \quad \mathbf{a}_i = \frac{\mathbf{f}_i}{m_i} \neq \mathbf{a}_{i,a} \neq \mathbf{a}_{i,b} \quad (5.19)$$

After the force extrapolations are done, the following calculations complete the force calculations:

1. Restore momentum to the grid, but at multimaterial nodes, it will restore  $\mathbf{p}_{i,j}$ , which are momenta after contact laws changes done above, but before boundary conditions.
2. Add traction loads on particles.

$$\mathbf{f}_{i,T,j} = \int_{S_{T,j}} \mathbf{N}_i(\mathbf{x}) \mathbf{T} dS \quad (5.20)$$

3. Add any body forces (such as gravity) assuming independent of particle state

$$\mathbf{f}_{i,b,j} = \sum_{p \in j} m_p S_{ip}^{(n)} \mathbf{b}_p = m_{i,j} \mathbf{b}_i \quad (5.21)$$

Because  $j = a$  or  $b$  only here, total body force is sum of force on the two materials.

4. For nodes with velocity boundary conditions, set the force such that the momenta restored in step #1 above will update to the specified velocity.

## 5.7 XPIC Task

The key XPIC task is to find

$$\mathbf{v}^* = \sum_{k=2}^m (-1)^k \mathbf{v}_k^* \quad (5.22)$$

where

$$\mathbf{v}_1^* = \mathbf{v} \quad (5.23)$$

and

$$(\mathbf{v}_k^*)_i = \frac{m-k+1}{k} \sum_p \sum_{j \in \Omega_p} \frac{M_p S_{pi} S_{pj}}{m_i^{(n)}} (\mathbf{v}_{k-1}^*)_j \quad \forall i \in \Omega_p \quad (5.24)$$

In multimaterial mode, each material is done separately. But, if the interface and perfect, interface nodes will have  $\mathbf{v}_a = \mathbf{v}_b = \mathbf{v}_i$ . This state means that  $\mathbf{v}_1^* = \mathbf{v}$  in both material domains. But, then separate domains will have:

$$(\mathbf{v}_k^*)_{i,a} = \frac{m-k+1}{k} \sum_{p \in a} \sum_{j \in \Omega_p} \frac{M_p S_{pi} S_{pj}}{m_i^{(n)}} (\mathbf{v}_{k-1}^*)_j \quad \forall i \in \Omega_p \quad (5.25)$$

$$(\mathbf{v}_k^*)_{i,b} = \frac{m-k+1}{k} \sum_{p \in b} \sum_{j \in \Omega_p} \frac{M_p S_{pi} S_{pj}}{m_i^{(n)}} (\mathbf{v}_{k-1}^*)_j \quad \forall i \in \Omega_p \quad (5.26)$$

$$(\mathbf{v}_k^*)_i = (\mathbf{v}_k^*)_{i,a} + (\mathbf{v}_k^*)_{i,b} \quad (5.27)$$

$$\mathbf{v}^* = \sum_{k=2}^m (-1)^k ((\mathbf{v}_k^*)_a + (\mathbf{v}_k^*)_b) = \mathbf{v}_a^* + \mathbf{v}_b^* \quad (5.28)$$

## 5.8 Update Nodal Momenta

This task updates the nodal momenta. In single material mode:

$$\mathbf{p}_i^+ = \mathbf{p}_i + \mathbf{f}_i \Delta t \quad (5.29)$$



In multimaterial mode

$$\mathbf{p}_{i,j}^+ = \mathbf{p}'_{i,j} + \mathbf{f}'_{i,j}\Delta t \quad \text{where} \quad \mathbf{f}'_{i,j} = \mathbf{f}_{i,j} - \frac{\Delta \mathbf{p}_{i,j}}{\Delta t} \quad (5.30)$$

includes  $\mathbf{f}_{i,j}$  as force due to stress, force, traction, and body force extrapolations and the subtracted force is due to earlier contact calculations. Note that we can also write

$$\mathbf{p}_{i,j}^+ = \mathbf{p}_{i,j} + \mathbf{f}_{i,j}\Delta t \quad (5.31)$$

For any interface we will recover

$$\mathbf{p}_i^+ = \mathbf{p}_{i,a}^+ + \mathbf{p}_{i,b}^+ \quad (5.32)$$

including at velocity boundary conditions.

To keep contact correct, revisit contact law calculations in section 5.2, find contact momentum change  $\Delta \mathbf{p}_{i,j}^+$  and update force to match with:

$$\Delta \mathbf{f}_{i,j} = \frac{\Delta \mathbf{p}_{i,j}^+}{\Delta t} \quad (5.33)$$

For this special case of perfect interface, material  $a$  will have

$$\Delta \mathbf{p}_{i,a}^+ = \frac{m_{i,a}}{m_i} \mathbf{p}_i^+ - \mathbf{p}_{i,a}^+ \quad (5.34)$$

$$\mathbf{p}_{i,a}^{+'} = \frac{m_{i,a}}{m_i} \mathbf{p}_i^+ \quad (5.35)$$

The revised material  $a$  (and similar for  $b$ ) will be

$$\mathbf{p}_{i,j}^{+'} = \mathbf{p}'_{i,j} + \mathbf{f}_{i,j}''\Delta t = \mathbf{p}'_{i,j} + \left( \mathbf{f}_{i,a} + \frac{\Delta \mathbf{p}_{i,a}^+ - \Delta \mathbf{p}_{i,a}}{\Delta t} \right) \Delta t \quad (5.36)$$

The velocity in both modes are the same ( $\mathbf{v}_i = \mathbf{v}_{i,j}$ ). The acceleration should be based on  $\mathbf{f}_{i,j}''/m_{i,j}$ . For material  $a$  and perfect interface

$$\mathbf{a}_{i,a} = \frac{1}{m_{i,a}} \left( \mathbf{f}_{i,a} + \frac{\Delta \mathbf{p}_{i,a}^+ - \Delta \mathbf{p}_{i,a}}{\Delta t} \right) \quad (5.37)$$

$$= \frac{1}{m_{i,a}} \left( \mathbf{f}_{i,a} + \frac{m_{i,a}}{m_i} \left( \frac{\mathbf{p}_i}{\Delta t} + \mathbf{f}_i \right) - \left( \frac{\mathbf{p}_{i,a}}{\Delta t} + \mathbf{f}_{i,a} \right) - \frac{m_{i,a}}{m_i} \frac{\mathbf{p}_i}{\Delta t} + \frac{\mathbf{p}_{i,a}}{\Delta t} \right) \quad (5.38)$$

$$= \frac{\mathbf{f}_i}{m_i} \quad (5.39)$$

which is correct result for a perfect interface

These contact calculations might disrupt velocity boundary conditions so they need to be reimposed. If momentum is changed, it should change force as well. Actually force started this step in synch with boundary conditions, but contact calculations might add  $\Delta \mathbf{f}_{i,j}$  that disrupts it, so it needs to be brought back in line.

## 5.9 Update the Particles

Update particle position and velocity using nodal velocity (from  $\mathbf{v}_{i,j} = \mathbf{p}_{i,j}^+/m_{i,j}$ ) and acceleration (from  $\mathbf{a}_{i,j} = \mathbf{f}_{i,j}/m_{i,j}$ ) found in material velocity field `matfld` (for type of particle). For a perfect interface, we should have (within round-off error)  $\mathbf{v}_{i,j} = \mathbf{v}_i$  and  $\mathbf{a}_{i,j} = \mathbf{a}_i$ , meaning the particle update should be same for single and multimaterial modes.

## 5.10 Optional Update Strain Last

Optionally the calculations can update strain after the momentum update. This calculation is done for methods  $\text{USAVG}\pm$  and for  $\text{USL}\pm$ . For a perfect interface, these should continue to see single an multimaterial mode have the same velocity and therefore results should be identical.

# Bibliography

- [1] C. C. Hammerquist and J. A. Nairn. Modeling nanoindentation using the material point method. *J. Materials Research*, submitted, 2017.
- [2] J. A. Nairn. Numerical implementation of imperfect interfaces. *Computational Materials Science*, 40:525–536, 2007.
- [3] J. A. Nairn, S. G. Bardenhagen, and G. S. Smith. Generalized contact and improved frictional heating in the material point method. *Computational Particle Mechanics*, in press, 2017.