

# Heterogeneous Electrode Geometry Generation

## Introduction

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This example demonstrates how to create a representative 3D geometry of a Nickel-Manganese-Cobalt (NMC) electrode, based on X-ray tomography data. The generated geometry may be used for subsequent heterogeneous electrode modeling.

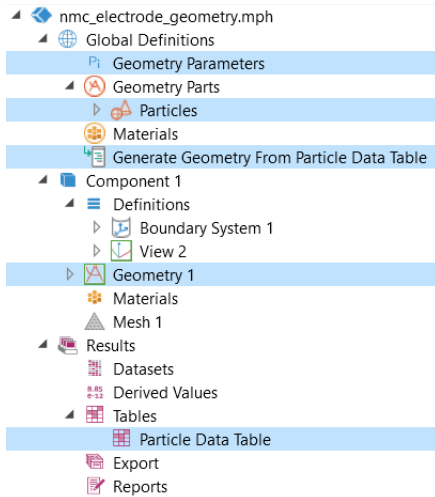
The X-ray data differentiates between the active NMC particles, and a surrounding mix of electron-conductive carbon additives, binder and ion-conducting electrolyte.

A model method is used to loop through tabulated data on particle volumes and center coordinates.

The geometry generated is used in the [Heterogeneous NMC Electrode](#) and [Homogenizing a Heterogeneous Electrode Model](#) tutorials in the Battery Design Module Application Library.

## Model Definition

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*Figure 1: Model tree setup. All nodes involved in the geometry generation are highlighted in blue.*

[Figure 1](#) indicates the nodes in the model tree used for creating the geometry.

#### *Global Definitions>Geometry Parameters*

This node contains parameters related to the generation of the geometry, such as tolerance factors and bounding box.

#### *Results>Tables>Particle Data Table*

This table contains the particle center coordinates and the particle volumes.

The data, available as open source, was downloaded from <http://dx.doi.org/10.5905/ethz-iis-1>. See [Ref. 1](#) for additional information about the data.

#### *Global Definitions>Geometry Parts>Particles*

This geometry part contains the individual particle geometric objects (spheres and cylinders) generated from **Particle Data Table**.

#### *Global Definitions>Generate Geometry From Particle Data Table*

This node calls the corresponding model method for generating the **Geometry Parts>Particles** geometric objects, based on the **Particle Data Table**.

Each entry in the **Particle Data Table** generates a **Sphere** object in **Particles**, given that the particle is located within the bounding box defined in **Geometry Parameters**.

**Cylinder** objects are also added in **Geometry Parts>Particles** in order to avoid sharp angles between overlapping spheres, or to join spheres together when the mutual distance is lower than the tolerance factor.

The model method also calculates and updates the electrode thickness parameter in **Geometry Parameters**.

The model method is written in Java, and maybe edited by switching to the **Application Builder** desktop. Note however that the model method editor is only available in the COMSOL desktop when using Windows. The model method node created in the model is however available and may be executed on all operating systems.

#### *Component 1>Geometry*

The final geometry sequence imports the **Geometry Parts>Particles**, and also includes block geometry objects to define the separator and porous conductive binder domains, based on the settings in **Geometry Parameters**. Geometry selections are also included in order to facility physics setup.

## Results and Discussion

Figure 2 shows the generated model geometry.

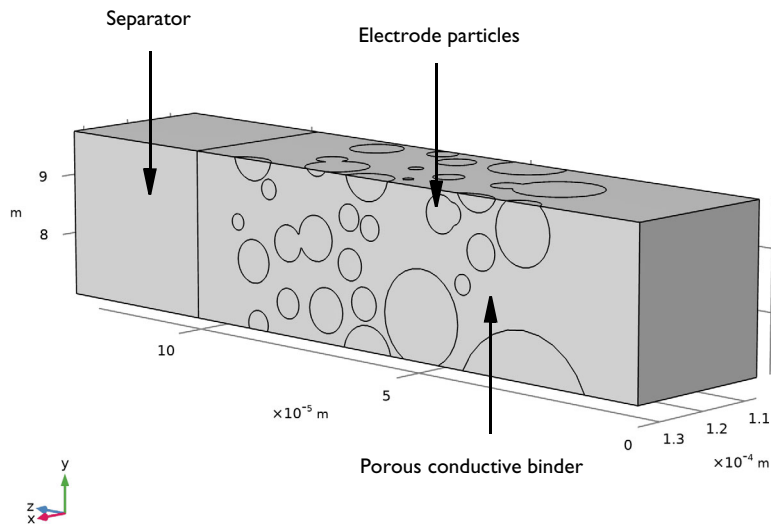


Figure 2: Generated model geometry.

## Reference


1. M.Ebner, F. Geldmacher, F. Marone, M. Stampanoni, and V. Wood, "X-Ray Tomography of Porous, Transition Metal Oxide Based Lithium Ion Battery Electrodes," *Adv. Energy Mater.*, vol. 3, pp. 845–850, 2013. See also supporting information at <https://onlinelibrary.wiley.com/doi/abs/10.1002/aenm.201200932>

**Application Library path:** Battery\_Design\_Module/Batteries,\_Heterogeneous/nmc\_electrode\_geometry

## Modeling Instructions

From the **File** menu, choose **New**.


## NEW

In the **New** window, click  **Blank Model**.

## GLOBAL DEFINITIONS

### *Geometry Parameters*

Load some geometry parameters from a text file.



- 1 In the **Model Builder** window, under **Global Definitions** click **Parameters 1**.
- 2 In the **Settings** window for **Parameters**, type Geometry Parameters in the **Label** text field.
- 3 Locate the **Parameters** section. Click  **Load from File**.
- 4 Browse to the model's Application Libraries folder and double-click the file `nmc_electrode_geometry_parameters.txt`.

## PARTICLES

The particle data is placed in **Geometry Parts**, which is then imported into the model geometry.

- 1 In the **Model Builder** window, right-click **Global Definitions** and choose **Geometry Parts> 3D Part**.
- 2 In the **Settings** window for **Part**, type Particles in the **Label** text field.  
Create a dummy sphere in the part. This will later be replaced by a collection of objects, generated from a model method.

### *Sphere 1 (sph1)*

- 1 In the **Geometry** toolbar, click  **Sphere**.
- 2 In the **Model Builder** window, expand the **Particles** node, then click **Sphere 1 (sph1)**.
- 3 In the **Settings** window for **Sphere**, locate the **Size** section.
- 4 In the **Radius** text field, type  $2\text{e-}6$ .
- 5 Locate the **Position** section. In the **x** text field, type  $11\text{e-}5$ .
- 6 In the **y** text field, type  $7\text{e-}5$ .
- 7 In the **z** text field, type  $4\text{e-}6$ .
- 8 Click  **Build Selected**.



## ADD COMPONENT

In the **Home** toolbar, click  **Add Component** and choose **3D**.


## GEOMETRY I

Now set up the model geometry sequence, making use of the **Particles** geometry part.



### *Particles Part*

- 1 In the **Geometry** toolbar, click  **Parts** and choose **Particles**.
- 2 In the **Settings** window for **Part Instance**, type **Particles Part** in the **Label** text field.
- 3 Click  **Build Selected**.



### *Particles*

- 1 In the **Geometry** toolbar, click  **Booleans and Partitions** and choose **Union**.
- 2 In the **Settings** window for **Union**, type **Particles** in the **Label** text field.
- 3 Select the object **pi1** only.
- 4 Locate the **Union** section. Clear the **Keep interior boundaries** check box.
- 5 Locate the **Selections of Resulting Entities** section. Select the **Resulting objects selection** check box.

### *Block I (blk1)*


- 1 In the **Geometry** toolbar, click  **Block**.
- 2 In the **Settings** window for **Block**, locate the **Size and Shape** section.
- 3 In the **Width** text field, type  $x_{\max} - x_{\min}$ .
- 4 In the **Depth** text field, type  $y_{\max} - y_{\min}$ .
- 5 In the **Height** text field, type  $L_{\text{elec}} + L_{\text{sep}}$ .
- 6 Locate the **Position** section. In the **x** text field, type  $x_{\min}$ .
- 7 In the **y** text field, type  $y_{\min}$ .
- 8 Click  **Build Selected**.

### *Partition Objects I (par1)*




- 1 In the **Geometry** toolbar, click  **Booleans and Partitions** and choose **Partition Objects**.
- 2 Select the object **uni1** only.
- 3 In the **Settings** window for **Partition Objects**, locate the **Partition Objects** section.
- 4 Find the **Tool objects** subsection. Click to select the  **Activate Selection** toggle button.
- 5 Select the object **blk1** only.

### *Box Selection I (boxsel1)*


- 1 In the **Geometry** toolbar, click  **Selections** and choose **Box Selection**.

- 2 In the **Settings** window for **Box Selection**, locate the **Box Limits** section.
- 3 In the **x minimum** text field, type `xmin+geom_tol`.
- 4 In the **x maximum** text field, type `xmax-geom_tol`.
- 5 In the **y minimum** text field, type `ymin+geom_tol`.
- 6 In the **y maximum** text field, type `ymax-geom_tol`.
- 7 Locate the **Resulting Selection** section. From the **Show in physics** list, choose **Off**.
- 8 Click  **Build Selected**.



#### *Complement Selection I (comsell)*

- 1 In the **Geometry** toolbar, click  **Selections** and choose **Complement Selection**.
- 2 In the **Settings** window for **Complement Selection**, locate the **Input Entities** section.
- 3 Click  **Add**.
- 4 In the **Add** dialog box, select **Box Selection I** in the **Selections to invert** list.
- 5 Click **OK**.
- 6 In the **Settings** window for **Complement Selection**, locate the **Resulting Selection** section.
- 7 From the **Show in physics** list, choose **Off**.
- 8 Click  **Build Selected**.




#### *Delete Entities I (delI)*

- 1 In the **Model Builder** window, right-click **Geometry I** and choose **Delete Entities**.
- 2 In the **Settings** window for **Delete Entities**, locate the **Entities or Objects to Delete** section.
- 3 From the **Geometric entity level** list, choose **Domain**.
- 4 From the **Selection** list, choose **Complement Selection I**.
- 5 Click  **Build Selected**.


#### *Electrode*

- 1 In the **Geometry** toolbar, click  **Block**.
- 2 In the **Settings** window for **Block**, type **Electrode** in the **Label** text field.
- 3 Locate the **Size and Shape** section. In the **Width** text field, type `xmax-xmin`.
- 4 In the **Depth** text field, type `ymax-ymin`.
- 5 In the **Height** text field, type `L_elec`.
- 6 Locate the **Position** section. In the **x** text field, type `xmin`.
- 7 In the **y** text field, type `ymin`.
- 8 Click  **Build Selected**.




### *Difference 1 (dif1)*

- 1 In the **Geometry** toolbar, click  **Booleans and Partitions** and choose **Difference**.
- 2 Select the object **blk2** only.
- 3 In the **Settings** window for **Difference**, locate the **Difference** section.
- 4 Find the **Objects to subtract** subsection. Click to select the  **Activate Selection** toggle button.
- 5 Select the object **dif1** only.
- 6 Select the **Keep objects to add** check box.
- 7 Select the **Keep objects to subtract** check box.
- 8 Click  **Build Selected**.

### *Delete Entities 2 (del2)*

- 1 Right-click **Geometry 1** and choose **Delete Entities**.
- 2 In the **Settings** window for **Delete Entities**, locate the **Entities or Objects to Delete** section.
- 3 From the **Geometric entity level** list, choose **Object**.
- 4 Select the object **dif1** only.
- 5 Click  **Clear Selection**.
- 6 Select the object **blk2** only.

### *Separator*

- 1 In the **Geometry** toolbar, click  **Block**.
- 2 In the **Settings** window for **Block**, type Separator in the **Label** text field.
- 3 Locate the **Size and Shape** section. In the **Width** text field, type  $x_{max} - x_{min}$ .
- 4 In the **Depth** text field, type  $y_{max} - y_{min}$ .
- 5 In the **Height** text field, type  $L_{sep}$ .
- 6 Locate the **Position** section. In the **x** text field, type  $x_{min}$ .
- 7 In the **y** text field, type  $y_{min}$ .
- 8 In the **z** text field, type  $L_{elec}$ .
- 9 Locate the **Selections of Resulting Entities** section. Select the **Resulting objects selection** check box.
- 10 Click  **Build Selected**.
- 11 Click the  **Zoom Extents** button in the **Graphics** toolbar.



### *Form Union (fin)*

- 1 In the **Model Builder** window, click **Form Union (fin)**.




- 2 In the **Settings** window for **Form Union/Assembly**, click  **Build Selected**.


#### *Porous Conductive Binder*

- 1 In the **Geometry** toolbar, click  **Selections** and choose **Complement Selection**.
- 2 In the **Settings** window for **Complement Selection**, type Porous Conductive Binder in the **Label** text field.
- 3 Locate the **Input Entities** section. Click  **Add**.
- 4 In the **Add** dialog box, in the **Selections to invert** list, choose **Particles** and **Separator**.
- 5 Click **OK**.



#### *Lithium Foil*

- 1 In the **Geometry** toolbar, click  **Selections** and choose **Box Selection**.
- 2 In the **Settings** window for **Box Selection**, type Lithium Foil in the **Label** text field.
- 3 Locate the **Geometric Entity Level** section. From the **Level** list, choose **Boundary**.
- 4 Locate the **Box Limits** section. In the **z minimum** text field, type  $L_{elec} + L_{sep}/2$ .
- 5 Locate the **Output Entities** section. From the **Include entity if** list, choose **Entity inside box**.



#### *NMC Current Collector*

- 1 In the **Geometry** toolbar, click  **Selections** and choose **Box Selection**.
- 2 In the **Settings** window for **Box Selection**, type NMC Current Collector in the **Label** text field.
- 3 Locate the **Geometric Entity Level** section. From the **Level** list, choose **Boundary**.
- 4 Locate the **Box Limits** section. In the **z maximum** text field, type 0.
- 5 Locate the **Output Entities** section. From the **Include entity if** list, choose **Entity inside box**.



#### *Particle Boundaries*

- 1 In the **Geometry** toolbar, click  **Selections** and choose **Adjacent Selection**.
- 2 In the **Settings** window for **Adjacent Selection**, type Particle Boundaries in the **Label** text field.
- 3 Locate the **Input Entities** section. Click  **Add**.
- 4 In the **Add** dialog box, select **Particles** in the **Input selections** list.
- 5 Click **OK**.
- 6 In the **Settings** window for **Adjacent Selection**, locate the **Resulting Selection** section.
- 7 From the **Show in physics** list, choose **Off**.



### *Binder Boundaries*

- 1 In the **Geometry** toolbar, click  **Selections** and choose **Adjacent Selection**.
- 2 In the **Settings** window for **Adjacent Selection**, type **Binder Boundaries** in the **Label** text field.
- 3 Locate the **Input Entities** section. Click  **Add**.
- 4 In the **Add** dialog box, select **Porous Conductive Binder** in the **Input selections** list.
- 5 Click **OK**.
- 6 In the **Settings** window for **Adjacent Selection**, locate the **Resulting Selection** section.
- 7 From the **Show in physics** list, choose **Off**.


### *Particle Surfaces*

- 1 In the **Geometry** toolbar, click  **Selections** and choose **Intersection Selection**.
- 2 In the **Settings** window for **Intersection Selection**, type **Particle Surfaces** in the **Label** text field.
- 3 Locate the **Geometric Entity Level** section. From the **Level** list, choose **Boundary**.
- 4 Locate the **Input Entities** section. Click  **Add**.
- 5 In the **Add** dialog box, in the **Selections to intersect** list, choose **Particle Boundaries** and **Binder Boundaries**.
- 6 Click **OK**.

### *Separator + Binder*



- 1 In the **Geometry** toolbar, click  **Selections** and choose **Union Selection**.
- 2 In the **Settings** window for **Union Selection**, type **Separator + Binder** in the **Label** text field.
- 3 Locate the **Input Entities** section. Click  **Add**.
- 4 In the **Add** dialog box, in the **Selections to add** list, choose **Separator** and **Porous Conductive Binder**.
- 5 Click **OK**.

### *All Domains*

- 1 In the **Geometry** toolbar, click  **Selections** and choose **Box Selection**.
- 2 In the **Settings** window for **Box Selection**, type **All Domains** in the **Label** text field.
- 3 Locate the **Box Limits** section. In the **x minimum** text field, type **xmin**.
- 4 In the **x maximum** text field, type **xmax**.
- 5 In the **y minimum** text field, type **ymin**.

- 6 In the **y maximum** text field, type **y<sub>max</sub>**.
- 7 Locate the **Output Entities** section. From the **Include entity if** list, choose **All vertices inside box**.
- 8 Locate the **Resulting Selection** section. From the **Show in physics** list, choose **Off**.

#### *All External Boundaries*

- 1 In the **Geometry** toolbar, click  **Selections** and choose **Adjacent Selection**.
- 2 In the **Settings** window for **Adjacent Selection**, type All External Boundaries in the **Label** text field.
- 3 Locate the **Input Entities** section. Click  **Add**.
- 4 In the **Add** dialog box, select **All Domains** in the **Input selections** list.
- 5 Click **OK**.

### **RESULTS**


Import particle data in the form of a table.

#### *Particle Data Table*

- 1 In the **Model Builder** window, expand the **Results** node.
- 2 Right-click **Results>Tables** and choose **Table**.
- 3 In the **Settings** window for **Table**, type Particle Data Table in the **Label** text field.  
Download the data from the entry Particle Statistics (NMC\_94wt\_0bar) from <http://dx.doi.org/10.5905/ethz-iis-1> and import it as a text file to the table.
- 4 Locate the **Data** section. Click **Import**.

For generating the **Particles** geometry part based on the table data, we will use a model method. Note that the method editor is only available in the Windows® version of the COMSOL Desktop.

### **APPLICATION BUILDER**

In the **Home** toolbar, click  **Application Builder**.

### **NEW METHOD**

- 1 In the **Application Builder** window, right-click **Methods** and choose **New Method**.
- 2 In the **New Method** dialog box, type GenerateGeometryFromTable in the **Name** text field.
- 3 Click **OK**.

### *GenerateGeometryFromTable*

1 In the **Application Builder** window, under **Methods** click **GenerateGeometryFromTable**.

2 Copy the following code into the **GenerateGeometryFromTable** window:

```
// Read data from Results->Table in model builder tree
String tableData[][] = model.result().table().get("tbl1").getTableData(false);

// Read some geometry parameter values from Parameters
Double voxel_size = model.param().evaluate("vox_size");
Double geom_tol = model.param().evaluate("geom_tol");
Double xmin = model.param().evaluate("xmin")+geom_tol;
Double xmax = model.param().evaluate("xmax")+geom_tol;
Double ymin = model.param().evaluate("ymin")+geom_tol;
Double ymax = model.param().evaluate("ymax")+geom_tol;

// Clear geometry part
model.geom("part1").feature().clear();

// Create matrix for added spheres
Double sphereData[][] = new Double[tableData.length][4];
int i = 0; // Row counter for matrix

// Variable for electrode thickness calculation
Double maxz = new Double(0.0);

// Variables for calculating mass-averaged sphere radius
Double Vsum = new Double(0.0);
Double rVsum = new Double(0.0);

// Loop for all rows in the table
for (String[] row : tableData) {
    Double xminData = Double.parseDouble(row[2])*voxel_size;
    Double yminData = Double.parseDouble(row[3])*voxel_size;
    Double xmaxData = Double.parseDouble(row[5])*voxel_size;
    Double ymaxData = Double.parseDouble(row[6])*voxel_size;

    // Check if within bounding box, and add sphere if radius larger than tolerance
    if (((xminData < xmax && xminData > xmin) || (xmaxData < xmax && xmaxData >
xmin)) &&
        ((yminData < ymax && yminData > ymin) || (ymaxData < ymax && ymaxData > ymin))) {
        Double volume = Double.parseDouble(row[1])*Math.pow(voxel_size, 3.0);
        Double radius = Math.pow(0.75*volume/Math.PI, 0.333333);

        if (radius > geom_tol) {
            Double xpos = Double.parseDouble(row[8])*voxel_size;
            Double ypos = Double.parseDouble(row[9])*voxel_size;
            Double zpos = Double.parseDouble(row[10])*voxel_size;

            // Add data to sphere data matrix
            sphereData[i][0] = xpos;
            sphereData[i][1] = ypos;
            sphereData[i][2] = zpos;
            sphereData[i][3] = radius;
            i++;
        }
    }
}
```

```

// Update electrode thickness variable
maxz = maxz < (zpos+radius) ? zpos+radius : maxz;

// Update mass-averaged radius variables if center of sphere inside bounding box
if (((xpos < xmax && xpos > xmin) && (ypos < ymax && ypos > ymin))) {
Vsum = Vsum+volume;
rVsum = rVsum+radius*volume;
}

// Add sphere to part geometry
GeomFeature sph = model.geom("part1").create("sph"+i, "Sphere");
sph.set("pos", new String[]{Double.toString(xpos)+"[m]",
Double.toString(ypos)+"[m]", Double.toString(zpos)+"[m]"});
sph.label("Particle "+(int) Math.round(Double.parseDouble(row[0])));
sph.set("r", radius+"[m]");
}
}

// Add cylinders to eliminate narrow regions or sharp angles between spheres
double hmin = model.param().evaluate("hmin"); // Minimum mesh element size
int cyl = 1;

// Check distances and gaps between all binary combinations of spheres
for (int j = 0; j < i; j++) {
for (int k = j+1; k < i; k++) {
Double x = sphereData[k][0]-sphereData[j][0];
Double y = sphereData[k][1]-sphereData[j][1];
Double z = sphereData[k][2]-sphereData[j][2];

Double dist = Math.sqrt(x*x+y*y+z*z);

Double gap = dist-(sphereData[k][3]+sphereData[j][3]);
int index1 = sphereData[k][3] < sphereData[j][3] ? k : j;
int index2 = sphereData[k][3] > sphereData[j][3] ? k : j;
Double r_sphere = sphereData[index1][3];

// Add cylinder if spheres are too close or overlapping
if (gap < hmin && gap > -0.667*r_sphere) {

Double x_sphere = sphereData[index1][0];
Double y_sphere = sphereData[index1][1];
Double z_sphere = sphereData[index1][2];

Double direction = sphereData[k][3] < sphereData[j][3] ? -1.0 : 1.0;

Double r_help = r_sphere+gap-hmin;
Double r_cyl = Math.max(Math.sqrt(r_sphere*r_sphere-r_help*r_help), hmin*2);

// Solve root to non-linear equation for cylinder radius (so that sphere-to-
sphere distance = hmin along cylinder surface).
Double r1 = r_sphere;
Double r2 = sphereData[index2][3];
double ta = 2*r2;

```

```

double tb = 2*r1;
double tc = Math.sqrt(-((gap-hmin+tb+ta)*(gap-hmin+tb)*(gap-hmin+ta)*(gap-
hmin)));
r_cyl = ((1/(gap-hmin+r1+r2))*tc)/2;

Double dist_corr = Math.sqrt(r_sphere*r_sphere-r_cyl*r_cyl)-2*hmin;
Double posx = x_sphere+x/dist*dist_corr*direction;
Double posy = y_sphere+y/dist*dist_corr*direction;
Double posz = z_sphere+z/dist*dist_corr*direction;

// Create cylinder
GeomFeature cylinder = model.geom("part1").create("cyl"+cyl++, "Cylinder");
with(cylinder);
set("pos", new String[] {Double.toString(posx)+"[m]", Double.toString(posy)+
"[m]", Double.toString(posz)+"[m]"});
set("axistype", "cartesian");
set("ax3", new String[] {Double.toString(x*direction)+"[m]", Double.toString(y*
direction)+"[m]", Double.toString(z*direction)+"[m]"});
set("r", r_cyl+"[m]");
set("h", hmin*5+"[m]");
endwith();
}
}

// Set electrode length in model
model.param().set("L_elec", maxz+geom_tol+"[m]");

// Build geometry in model
model.geom("part1").run();
model.component("comp1").geom("geom1").run();


// Measure and update volume parameter
model.component("comp1").geom("geom1").measure().selection().init(3);
model.component("comp1").geom("geom1").measure().selection().named("uni1");
model.param().set("Vp", model.geom("geom1").measure().getVolume()+"[m^3]");

// Measure and update area parameter
model.component("comp1").geom("geom1").measure().selection().init(2);
model.component("comp1").geom("geom1").measure().selection().named("intsel1");
model.param().set("Ap", model.geom("geom1").measure().getVolume()+"[m^2]");

// Update mass-averaged radius parameter
model.param().set("rp_avg_spheres", rVsum/Vsum+"[m]");


```

## METHODS


In the **Home** toolbar, click  **Model Builder**.

## GLOBAL DEFINITIONS

Make the model method available in the model tree, and run it as follows:

Click  **Method Call** and choose **GenerateGeometryFromTable**.

### *Generate Geometry From Particle Data Table*

- 1 In the **Model Builder** window, under **Global Definitions** click **GenerateGeometryFromTable 1**.
- 2 In the **Settings** window for **Method Call**, type Generate Geometry From Particle Data Table in the **Label** text field.
- 3 Click  **Run**.

### **PARTICLES**



In the **Model Builder** window, collapse the **Global Definitions>Geometry Parts>Particles** node.


### **DEFINITIONS**

Finally, update the view so that the y-axis points upwards, and the z-axis to the left, and the x-axis toward you.

- 1 In the **Model Builder** window, expand the **Component 1 (comp1)>Definitions** node.

#### *Camera*

- 1 In the **Model Builder** window, expand the **Component 1 (comp1)>Definitions>View 2** node, then click **Camera**.
- 2 Select the object only.
- 3 Click the  **Zoom Extents** button in the **Graphics** toolbar.
- 4 In the **Settings** window for **Camera**, locate the **Position** section.
- 5 In the **x** text field, type  $6.1 \times 10^{-4}$ .
- 6 In the **y** text field, type  $2.1 \times 10^{-4}$ .
- 7 In the **z** text field, type  $-4.1 \times 10^{-4}$ .
- 8 Locate the **Up Vector** section. In the **x** text field, type 0.
- 9 In the **y** text field, type 1.
- 10 In the **z** text field, type 0.
- 11 Click  **Update**.

12 Click the  **Zoom Extents** button in the **Graphics** toolbar.

