

GEC ICP Reactor, Argon Chemistry

The GEC cell was introduced by NIST (National Institute of Standards and Technology) in order to provide a standardized platform for experimental and modeling studies of discharges in different laboratories. The plasma is sustained via inductive heating. The Reference Cell operates as an inductively-coupled plasma in this model.

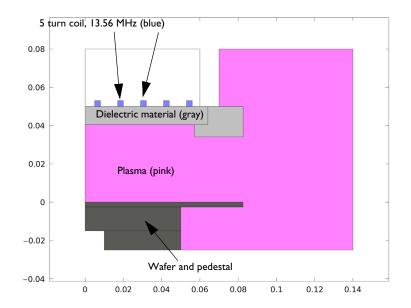


Figure 1: GEC ICP reactor geometry consisting of a 5 turn copper coil, plasma volume, dielectrics, and wafer with pedestal.

Note: This application requires the Plasma Module and the AC/DC Module.

Model Definition

Inductively coupled discharges typically operate at low pressures (<10 Pa) and high charge density (>10¹⁷ m⁻³). High density plasma sources are popular because low pressure ion bombardment can provide a greater degree of anisotropy on the surface of the wafer.

The electron density and mean electron energy are computed by solving a pair of driftdiffusion equations for the electron density and mean electron energy. For detailed

information on electron transport, see *Theory for the Drift Diffusion Interface* in the *Plasma Module User's Guide*.

$$\frac{\partial}{\partial t}(n_e) + \nabla \cdot [-\,n_e(\boldsymbol{\mu}_e \bullet \mathbf{E}) - \mathbf{D}_e \bullet \nabla n_e] \, = \, R_e$$

$$\frac{\partial}{\partial t}(n_{\varepsilon}) + \nabla \cdot [-n_{\varepsilon}(\mu_{\varepsilon} \bullet \mathbf{E}) - \mathbf{D}_{\varepsilon} \bullet \nabla n_{\varepsilon}] + \mathbf{E} \cdot \Gamma_{e} = R_{\varepsilon}$$

The electron source R_e and the energy loss due to inelastic collisions R_{ϵ} are defined later. The electron diffusivity, energy mobility, and energy diffusivity are computed from the electron mobility using:

$$\mathbf{D}_{e} = \mu_{e} T_{e}, \, \mu_{\varepsilon} = \left(\frac{5}{3}\right) \mu_{e}, \, \mathbf{D}_{\varepsilon} = \mu_{\varepsilon} T_{e}$$

The source coefficients in the above equations are determined by the plasma chemistry using rate coefficients. In the case of rate coefficients, the electron source term is given by:

$$R_e = \sum_{j=1}^{M} x_j k_j N_n n_e$$

where x_j is the mole fraction of the target species for reaction j, k_j is the rate coefficient for reaction j (SI unit: m^3/s), and N_n is the total neutral number density (SI unit: $1/m^3$). The electron energy loss is obtained by summing the collisional energy loss over all reactions:

$$R_{\varepsilon} = \sum_{i=1}^{P} x_{j} k_{j} N_{n} n_{e} \Delta \varepsilon_{j}$$

where $\Delta \epsilon_j$ is the energy loss from reaction j (SI unit: V). The rate coefficients can be computed from cross section data by the following integral:

$$k_k = \gamma \int_0^\infty \varepsilon \sigma_k(\varepsilon) f(\varepsilon) d\varepsilon$$

where $\gamma = (2q/m_e)^{1/2}$ (SI unit: $C^{1/2}/kg^{1/2}$), m_e is the electron mass (SI unit: kg), ε is energy (SI unit: V), σ_k is the collision cross section (SI unit: m²), and f is the electron energy distribution function.

For nonelectron species, the following equation is solved for the mass fraction of each species. For detailed information on the transport of the nonelectron species, see *Theory* for the Heavy Species Transport Interface in the Plasma Module User's Guide.

$$\rho \frac{\partial}{\partial t}(w_k) + \rho(\mathbf{u} \cdot \nabla)w_k = \nabla \cdot \mathbf{j}_k + R_k$$

The electrostatic field is computed using the following equation:

$$-\nabla \cdot \varepsilon_0 \varepsilon_r \nabla V = \rho$$

The space charge density ρ is automatically computed based on the plasma chemistry specified in the model using the formula:

$$\rho = q \left(\sum_{k=1}^{N} Z_k n_k - n_e \right)$$

For detailed information about electrostatics see Theory for the Electrostatics Interface in the Plasma Module User's Guide.

For a nonmagnetized, nonpolarized plasma, the induction currents are computed in the frequency domain using the following equation:

$$(j\omega\boldsymbol{\sigma} - \omega^2\boldsymbol{\varepsilon}_0)\mathbf{A} + \nabla\times(\boldsymbol{\mu}_0^{-1}\nabla\times\mathbf{A}) \;=\; \mathbf{J}^{\mathrm{e}}$$

The electromagnetic wave "sees" a plasma defined by the plasma conductivity in the cold plasma approximation that is set in the Plasma Conductivity Coupling multiphysics feature:

$$\sigma = \frac{n_e q^2}{m_e (v_e + j\omega)}$$

where n_e is the electron density, q is the electron charge, m_e is the electron mass, v_e is the collision frequency, and ω is the angular frequency. The Joule heating term that is responsible to heat the electrons is set in the **Electron Heat Source** multiphysics feature.

BOUNDARY CONDITIONS

Electrons are lost to the wall due to random motion within a few mean free paths of the wall and gained due to secondary emission effects, resulting in the following boundary condition for the electron flux:

$$\mathbf{n} \cdot \Gamma_e = \left(\frac{1}{2} \mathbf{v}_{e, \, \text{th}} n_e\right)$$

and the electron energy flux:

$$\mathbf{n} \cdot \Gamma_{\varepsilon} = \left(\frac{5}{6} v_{e, \, \text{th}} n_{\varepsilon}\right)$$

For the heavy species, ions are lost to the wall due to surface reactions and the fact that the electric field is directed toward the wall:

$$\mathbf{n} \cdot \mathbf{j}_k = M_w R_k + M_w c_k Z \mu_k (\mathbf{E} \cdot \mathbf{n}) [Z_k \mu_k (\mathbf{E} \cdot \mathbf{n}) > 0]$$

The walls of the reactor are grounded.

PLASMA CHEMISTRY

Because the physics occurring in an inductively coupled plasma is rather complex, it is always best to start a modeling project with a simple chemical mechanism. Argon is one of the simplest mechanisms to implement at low pressures. The electronically excited states can be lumped into a single species, which results in a chemical mechanism consisting of only 3 species and 7 reactions (electron impact cross sections are obtained from Ref. 3):

TABLE I: TABLE OF COLLISIONS AND REACTIONS MODELED.

REACTION	FORMULA	ТҮРЕ	Δε(eV)
1	e+Ar=>e+Ar	Elastic	0
2	e+Ar=>e+Ars	Excitation	11.5
3	e+Ars=>e+Ar	Superelastic	-11.5
4	e+Ar=>2e+Ar+	Ionization	15.8
5	e+Ars=>2e+Ar+	Ionization	4.24
6	Ars+Ars=>e+Ar+Ar+	Penning ionization	-
7	Ars+Ar=>Ar+Ar	Metastable quenching	-

Stepwise ionization (reaction 5) can play an important role in sustaining low pressure argon discharges. Excited argon atoms are consumed via superelastic collisions with electrons, quenching with neutral argon atoms, ionization or Penning ionization where two metastable argon atoms react to form a neutral argon atom, an argon ion and an

electron. In addition to volumetric reactions, the following surface reactions are implemented:

TABLE 2: TABLE OF SURFACE REACTIONS.

REACTION	FORMULA	STICKING COEFFICIENT
1	Ars=>Ar	1
2	Ar+=>Ar	I

When a metastable argon atom makes contact with the wall, it reverts to the ground state argon atom with some probability (the sticking coefficient).

ELECTRICAL EXCITATION

From an electrical point of view, the GEC reactor behaves as a transformer. A current is applied to the driving coil (the primary) and this induces a current in the plasma (the secondary). The plasma then induces an opposing current back in the coil, increasing its resistance. The current flowing in the plasma depends on the current applied to the coil and the reaction kinetics. The total plasma current can vary from no current (plasma not sustained) to the same current as the primary which corresponds to perfect coupling between the coil and the plasma.

In this example a fixed power of 1500 W is applied to the coil. Some of this power is dissipated in the coil, some is deposited into the plasma.

The peak electron density occurs at the center of the reactor, underneath the RF coil. The electron density in this case is high enough to cause some shielding of the azimuthal electric field.

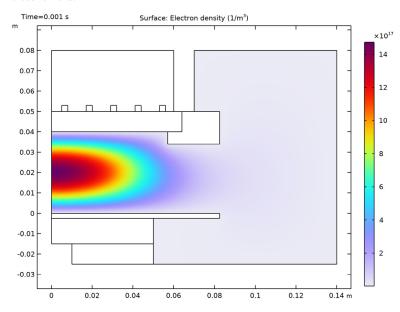


Figure 2: Plot of the electron density inside the GEC ICP reactor.

The electron "temperature" is highest directly underneath the coil, which is where the bulk of the power deposition occurs.

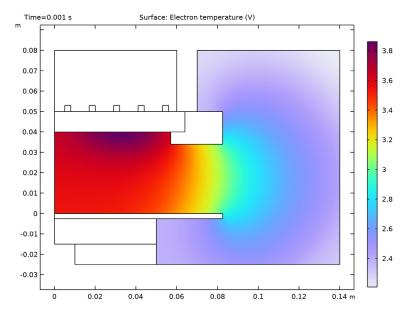


Figure 3: Plot of the electron "temperature" inside the GEC ICP reactor.

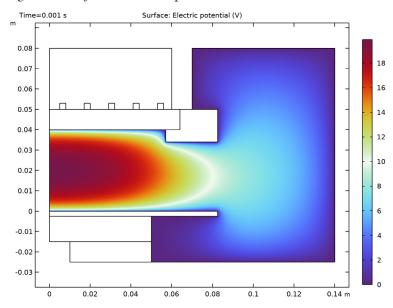


Figure 4: Plot of the electric potential inside the GEC ICP reactor.

From an electrical standpoint, the quantities of interest are total power deposition, coil resistance and inductance, and reactor efficiency. These "global" parameters are relatively easy to measure when the plasma is on or off, so such quantities provide an easy route for comparison with experimental data, without the need for expensive optical emission spectroscopy equipment or Langmuir probes.

The resistance of the coil increases by a little less than a factor of 4 when the plasma is on. When the plasma is on, there is a substantial opposing current induced back into the coil from the plasma. The electric potential applied across the coil needs to increase in order to maintain the same total current.

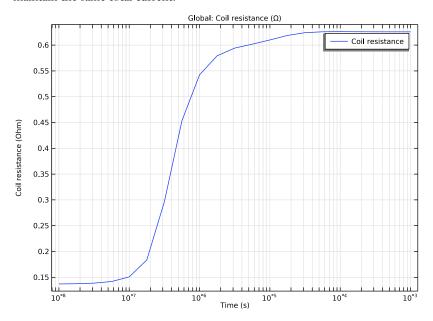


Figure 5: Plot of coil resistance versus time in GEC ICP reactor.

Initially the power dissipated is all dissipated in the coil (\sim 1400 W). After about 1 microsecond, the plasma ignition begins and as the neutral gas atoms split into electrons and ions, the electrons begin to absorb more and more power. Over a period of 2 microseconds, the plasma goes from absorbing no power to absorbing around 1200 W.

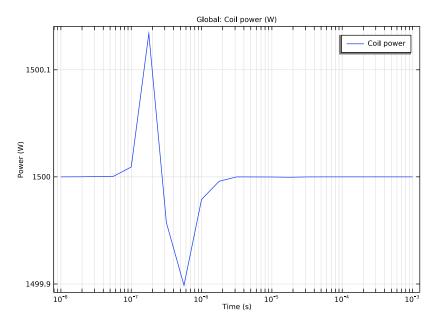


Figure 6: Plot of total power versus time in GEC ICP reactor.

The ion density is exactly the same except from in a thin region close to the walls. In this region, the ion density dominates the electron density which leads to a positive potential in the plasma bulk with respect to the walls. The positive potential increases the flux of ions and reduces the flux of electrons to the wall.

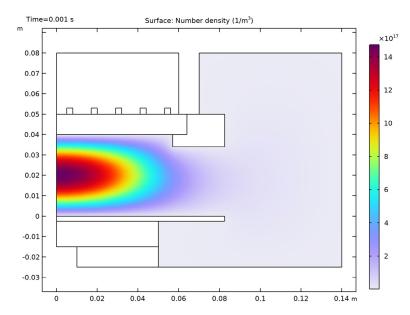


Figure 7: Plot of number density of argon ions in the GEC ICP reactor.

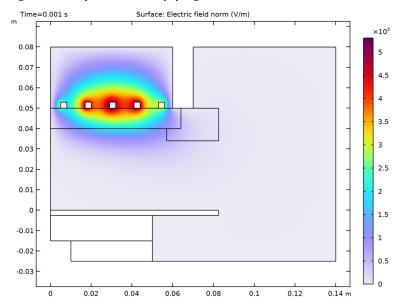


Figure 8: Plot of the norm of the electric field due to the induction currents.

The number density of excited species is also greatest in the center of the reactor. Unlike the charged species, there is no rapid drop off in number density close to the walls. The physics of the excited species is relatively simple: they are formed in the center of the reactor by high energy electrons and are lost to the via either stepwise ionization or diffusion to the wall. Because the excited argon atoms are not susceptible to migration due to the electric field, they can exist in much higher quantities than ions. The peak number density of excited argon atoms represents a mass fraction of around 0.02.

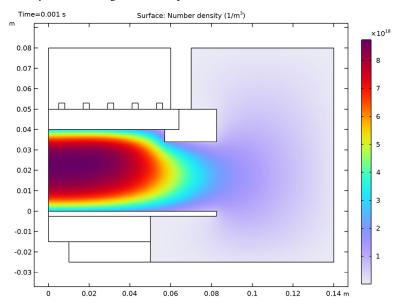


Figure 9: Plot of the number density of excited argon atoms in the GEC ICP reactor.

The skin depth of the plasma is on the order of 1cm which prevents the electric field from penetrating into the core of the plasma. The skin depth is defined as:

$$\delta = \sqrt{\frac{2}{\mu\omega\sigma}}$$

where μ is the permeability, σ is the plasma conductivity, and ω is the angular frequency. This tells us that increasing the driving frequency does not necessarily couple more power into the plasma. As the frequency increases, the plasma tends to shield the region over which power is deposited into a thin layer close to the upper wall.

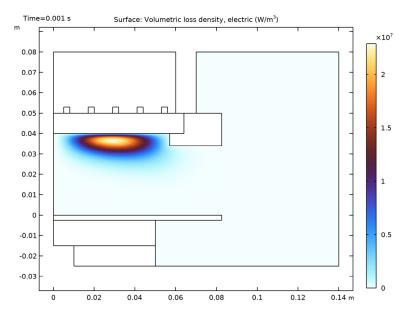


Figure 10: Plot of the power deposition into the plasma in the GEC ICP reactor. The region over which power is deposited to the plasma is governed by the plasma skin depth.

References

- 1. G.J.M. Hagelaar and L.C. Pitchford, "Solving the Boltzmann Equation to Obtain Electron Transport Coefficients and Rate Coefficients for Fluid Models," *Plasma Sources Sci. Technol.*, vol. 14, pp. 722–733, 2005.
- 2. D.P. Lymberopoulos and D.J. Economou, "Two-Dimensional Self-Consistent Radio Frequency Plasma Simulations Relevant to the Gaseous Electronics Conference RF Reference Cell," *J. Res. Natl. Inst. Stand. Technol.*, vol. 100, pp. 473–494, 1995.
- 3. Phelps database, www.lxcat.net, retrieved 2017.

Application Library path: Plasma_Module/Inductively_Coupled_Plasmas/argon_gec_icp

From the File menu, choose New.

NEW

In the New window, click Model Wizard.

MODEL WIZARD

- I In the Model Wizard window, click 2D Axisymmetric.
- 2 In the Select Physics tree, select Plasma>Inductively Coupled Plasma.
- 3 Click Add.
- 4 Click Study.
- 5 In the Select Study tree, select Preset Studies for Selected Multiphysics>Frequency-Transient.
- 6 Click **Done**.

GEOMETRY I

Insert the prepared geometry sequence from file. You can read the instructions for creating the geometry in the appendix.

- I In the Geometry toolbar, click Insert Sequence and choose Insert Sequence.
- 2 Browse to the model's Application Libraries folder and double-click the file argon gec icp geom sequence.mph.

Add some predefined selections for the geometric entities which will be referenced later on.

DEFINITIONS

Walls

- I In the **Definitions** toolbar, click **\(\frac{1}{3} \) Explicit**.
- 2 In the Settings window for Explicit, type Walls in the Label text field.
- 3 Locate the Input Entities section. From the Geometric entity level list, choose Boundary.
- **4** Select Boundaries 6, 8, 35–38, 44, 45, and 51–56 only.

Coils

- I In the **Definitions** toolbar, click **\(\frac{1}{3} \) Explicit**.
- 2 In the Settings window for Explicit, type Coils in the Label text field.
- 3 Select Domains 6 and 8–11 only.

Coil Boundaries

- I In the **Definitions** toolbar, click **\(\bigcap_{\bigcap} \) Explicit**.
- 2 In the Settings window for Explicit, type Coil Boundaries in the Label text field.
- **3** Select Domains 6 and 8–11 only.
- **4** Locate the **Output Entities** section. From the **Output entities** list, choose **Adjacent boundaries**.

GLOBAL DEFINITIONS

Parameters 1

- I In the Model Builder window, under Global Definitions click Parameters I.
- 2 In the Settings window for Parameters, locate the Parameters section.
- **3** In the table, enter the following settings:

Name	Expression	Value	Description
Psp	1500[W]	1500 W	Power input
mueN	4E24[1/(m*V*s)]	4E24 I/(V·m·s)	Reduced electron mobility
ТО	300[K]	300 K	Gas temperature
p0	0.02[torr]	2.6664 Pa	Gas pressure

PLASMA (PLAS)

- I In the Model Builder window, under Component I (compl) click Plasma (plas).
- 2 Select Domain 3 only.

Cross Section Import 1

- I In the Physics toolbar, click A Global and choose Cross Section Import.
- 2 In the Settings window for Cross Section Import, locate the Cross Section Import section.
- 3 Click **Browse**.
- **4** Browse to the model's Application Libraries folder and double-click the file Ar xsecs.txt.
- 5 Click | Import.
- 6 In the Model Builder window, click Plasma (plas).
- 7 In the Settings window for Plasma, locate the Plasma Properties section.

8 Select the Use reduced electron transport properties check box.

Now you add two more regular reactions which describe how electronically excited argon atoms are consumed on the volumetric level. The rate coefficients for these reactions are taken from the literature.

Reaction 1

- I In the Physics toolbar, click **Domains** and choose **Reaction**.
- 2 In the Settings window for Reaction, locate the Reaction Formula section.
- 3 In the Formula text field, type Ars+Ars=>e+Ar+Ar+.
- **4** Locate the **Reaction Parameters** section. In the k^{f} text field, type 3.734E8.

Reaction 2

- I In the Physics toolbar, click **Domains** and choose **Reaction**.
- 2 In the Settings window for Reaction, locate the Reaction Formula section.
- 3 In the Formula text field, type Ars+Ar=>Ar+Ar.
- **4** Locate the **Reaction Parameters** section. In the k^{f} text field, type 1807.

When solving any type of reacting flow problem there always needs to be one species which is selected to fulfill the mass constraint. This should be taken as the species with the largest mass fraction.

Species: Ar

- I In the Model Builder window, click Species: Ar.
- 2 In the Settings window for Species, locate the Species Formula section.
- 3 Select the From mass constraint check box.

When solving a plasma problem the plasma must be initially charge neutral. COMSOL automatically computes the initial concentration of a selected ionic species such that the initial electroneutrality constraint is satisfied. Once the simulation begins to time step, the plasma need not be charge neutral. In fact, the separation of space charge between the ions and electrons close to the wall is a critical component in sustaining the discharge.

Species: Ar+

- I In the Model Builder window, click Species: Ar+.
- 2 In the Settings window for Species, locate the Species Formula section.
- 3 Select the Initial value from electroneutrality constraint check box.

Initial conditions for the electron number density and mean electron energy are critical for any plasma model. If the initial electron density is too low then the plasma may not

be able to sustain itself and may self extinguish. If the initial electron density is too high then convergence problems may occur during initial time steps.

Initial Values 1

- I In the Model Builder window, click Initial Values I.
- 2 In the Settings window for Initial Values, locate the Initial Values section.
- **3** In the $n_{e,0}$ text field, type 1E15.
- **4** In the ε_0 text field, type 5.

The Magnetic Fields are computed everywhere except the wafer and the wafer pedestal.

MAGNETIC FIELDS (MF)

- I In the Model Builder window, under Component I (compl) click Magnetic Fields (mf).
- 2 Select Domains 3-6 and 8-12 only.

The **Coil** feature is used to electrically excite the system. The coil operates with a fixed total power of 1500 watts.

Coil I

- I In the Physics toolbar, click **Domains** and choose Coil.
- 2 In the Settings window for Coil, locate the Domain Selection section.
- 3 From the Selection list, choose Coils.
- 4 Locate the Coil section. From the Coil excitation list, choose Power.
- **5** Select the **Coil group** check box.
- **6** In the P_{coil} text field, type Psp.

PLASMA (PLAS)

Plasma Model I

- I In the Model Builder window, under Component I (compl)>Plasma (plas) click
 Plasma Model I
- 2 In the Settings window for Plasma Model, locate the Model Inputs section.
- **3** In the *T* text field, type T0.
- **4** In the p_A text field, type p0.
- 5 Locate the **Electron Density and Energy** section. In the $\mu_e N_n$ text field, type mueN. Next, define the material properties. There is no need to define the material properties in the plasma domain, as these are defined by the **Plasma Conductivity Coupling** feature.

MATERIALS

Material I (mat I)

- I In the Model Builder window, under Component I (compl) right-click Materials and choose Blank Material.
- 2 In the Settings window for Material, locate the Geometric Entity Selection section.
- **3** From the **Selection** list, choose **Coils**.
- **4** Locate the **Material Contents** section. In the table, enter the following settings:

Property	Variable	Value	Unit	Property group
Electrical conductivity	sigma_iso; sigmaii = sigma_iso, sigmaij = 0	6e7	S/m	Basic
Relative permeability	mur_iso; murii = mur_iso, murij = 0	1	I	Basic
Relative permittivity	epsilonr_iso; epsilonrii = epsilonr_iso, epsilonrij = 0	1	I	Basic

Material 2 (mat2)

- I Right-click Materials and choose Blank Material.
- **2** Select Domain 5 only.
- 3 In the Settings window for Material, locate the Material Contents section.
- **4** In the table, enter the following settings:

Property	Variable	Value	Unit	Property group
Relative permeability	mur_iso; murii = mur_iso, murij = 0	1	I	Basic

Property	Variable	Value	Unit	Property group
Electrical conductivity	sigma_iso; sigmaii = sigma_iso, sigmaij = 0	0	S/m	Basic
Relative permittivity	epsilonr_iso; epsilonrii = epsilonr_iso, epsilonrij = 0	1	I	Basic

Material 3 (mat3)

- I Right-click Materials and choose Blank Material.
- **2** Select Domains 4 and 12 only.
- 3 In the Settings window for Material, locate the Material Contents section.
- **4** In the table, enter the following settings:

Property	Variable	Value	Unit	Property group
Relative permeability	mur_iso; murii = mur_iso, murij = 0	1	I	Basic
Electrical conductivity	sigma_iso; sigmaii = sigma_iso, sigmaij = 0	0	S/m	Basic
Relative permittivity	epsilonr_iso; epsilonrii = epsilonr_iso, epsilonrij = 0	4.2	I	Basic

PLASMA (PLAS)

Surface reactions must always be included in a plasma model since they describe how ionic, excited and radical species interact with the wall.

Surface Reaction 1

- I In the Physics toolbar, click Boundaries and choose Surface Reaction.
- 2 In the Settings window for Surface Reaction, locate the Reaction Formula section.
- 3 In the Formula text field, type Ars=>Ar.
- 4 Locate the Boundary Selection section. From the Selection list, choose Walls.

Surface Reaction 2

- I In the Physics toolbar, click Boundaries and choose Surface Reaction.
- 2 In the Settings window for Surface Reaction, locate the Reaction Formula section.
- 3 In the Formula text field, type Ar+=>Ar.
- 4 Locate the Boundary Selection section. From the Selection list, choose Walls. Now, add boundary conditions to describe how the electrons interact with the wall.

Wall I

- I In the Physics toolbar, click Boundaries and choose Wall.
- 2 In the Settings window for Wall, locate the General Wall Settings section.
- **3** In the r_e text field, type 0.2.
- 4 Locate the Boundary Selection section. From the Selection list, choose Walls.

Ground I

- I In the Physics toolbar, click Boundaries and choose Ground.
- 2 In the Settings window for Ground, locate the Boundary Selection section.
- 3 From the Selection list, choose Walls.

MESH I

Meshing is a critical step in any plasma model. Boundary layer meshing on the reactor walls is nearly always necessary. This is needed to capture the separation of space charge between the electrons and ions close to the wall. You also add a fine mesh in the coil domains since the skin depth needs to be resolved.

- I In the Model Builder window, under Component I (compl) click Mesh I.
- 2 In the Settings window for Mesh, locate the Physics-Controlled Mesh section.
- **3** From the **Element size** list, choose **Finer**.

Edge 1

- I In the Mesh toolbar, click A Edge.
- **2** Select Boundaries 6, 8, 44, 45, and 54 only.

Size 1

- I Right-click Edge I and choose Size.
- 2 In the Settings window for Size, locate the Geometric Entity Selection section.
- 3 From the Geometric entity level list, choose Entire geometry.
- **4** Locate the **Element Size** section. Click the **Custom** button.

- 5 Locate the Element Size Parameters section.
- 6 Select the Maximum element size check box. In the associated text field, type 1E-3.

Free Triangular 1

- I In the Mesh toolbar, click Free Triangular.
- 2 In the Settings window for Free Triangular, locate the Domain Selection section.
- 3 From the Geometric entity level list, choose Domain.
- 4 Select Domain 3 only.

Size 1

- I Right-click Free Triangular I and choose Size.
- 2 In the Settings window for Size, locate the Element Size section.
- 3 From the Predefined list, choose Extra fine.

Boundary Layers 1

- I In the Mesh toolbar, click Boundary Layers.
- 2 In the Settings window for Boundary Layers, locate the Domain Selection section.
- 3 From the Geometric entity level list, choose Domain.
- 4 Select Domain 3 only.

Boundary Layer Properties

- I In the Model Builder window, click Boundary Layer Properties.
- 2 In the Settings window for Boundary Layer Properties, locate the Boundary Selection section.
- 3 From the Selection list, choose Walls.
- 4 Locate the Layers section. In the Number of layers text field, type 5.
- 5 In the Stretching factor text field, type 1.4.

Mapped I

- I In the Mesh toolbar, click Mapped.
- 2 In the Settings window for Mapped, locate the Domain Selection section.
- 3 From the Geometric entity level list, choose Domain.
- 4 From the Selection list, choose Coils.

Distribution I

- I Right-click Mapped I and choose Distribution.
- 2 In the Settings window for Distribution, locate the Boundary Selection section.

- 3 From the Selection list, choose Coil Boundaries.
- 4 Locate the Distribution section. From the Distribution type list, choose Predefined.
- 5 In the Number of elements text field, type 25.
- 6 In the Element ratio text field, type 20.
- 7 From the Growth rate list, choose Exponential.
- 8 Select the Symmetric distribution check box.

Free Triangular 2

- I In the Mesh toolbar, click Free Triangular.
- 2 In the Model Builder window, right-click Mesh I and choose Build All.

STUDY I

Steb 1: Frequency-Transient

- I In the Model Builder window, under Study I click Step I: Frequency-Transient.
- 2 In the Settings window for Frequency-Transient, locate the Study Settings section.
- 3 In the **Output times** text field, type 0 10^{range(-8,5/20,-3)}.
- 4 In the Frequency text field, type 13.56E6.
- 5 In the Home toolbar, click **Compute**.

RESULTS

Electron Density (plas)

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

Electron Temperature (plas)

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

Electric Potential (plas)

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

Now add a global plot for the coil resistance. This is defined as the real part of the total voltage drop over the coil divided by the applied current. The Coil feature creates predefined expressions for the resistance.

Coil Resistance

- I In the Home toolbar, click Add Plot Group and choose ID Plot Group.
- 2 In the Settings window for ID Plot Group, type Coil Resistance in the Label text field.
- 3 Locate the Plot Settings section.

- 4 Select the x-axis label check box. In the associated text field, type Time (s).
- 5 Select the **y-axis label** check box. In the associated text field, type Coil resistance (Ohm).

Global I

- I Right-click Coil Resistance and choose Global.
- 2 In the Settings window for Global, locate the y-Axis Data section.
- **3** In the table, enter the following settings:

Expression	Unit	Description
mf.RCoil_1	Ω	Coil resistance

- 4 Click the x-Axis Log Scale button in the Graphics toolbar.
- 5 In the **Coil Resistance** toolbar, click **Plot**.

 Now verify that 1500 watts is being applied to the system.

Coil Power

- I In the Home toolbar, click Add Plot Group and choose ID Plot Group.
- 2 In the Settings window for ID Plot Group, type Coil Power in the Label text field.
- 3 Locate the Plot Settings section.
- 4 Select the x-axis label check box. In the associated text field, type Time (s).
- 5 Select the y-axis label check box. In the associated text field, type Power (W).

Global I

- I Right-click Coil Power and choose Global.
- 2 In the Settings window for Global, locate the y-Axis Data section.
- **3** In the table, enter the following settings:

Expression	Unit	Description
mf.PCoil_1	W	Coil power

- 4 Click the x-Axis Log Scale button in the Graphics toolbar.
- 5 In the Coil Power toolbar, click Plot.

Ion Number Density

- I In the Home toolbar, click **Add Plot Group** and choose **2D Plot Group**.
- 2 In the Settings window for 2D Plot Group, type Ion Number Density in the Label text field.

Surface I

- I Right-click Ion Number Density and choose Surface.
- 2 In the Settings window for Surface, click Replace Expression in the upper-right corner of the Expression section. From the menu, choose Component I (compl)>Plasma> Number densities>plas.n_wAr_Ip - Number density - I/m3.
- 3 In the lon Number Density toolbar, click Plot.
- 4 Click the **Zoom Extents** button in the **Graphics** toolbar.

High Frequency Electric Field

- I In the Model Builder window, right-click Ion Number Density and choose Duplicate.
- 2 In the Settings window for 2D Plot Group, type High Frequency Electric Field in the Label text field.

Surface I

- I In the Model Builder window, expand the High Frequency Electric Field node, then click Surface 1.
- 2 In the Settings window for Surface, click Replace Expression in the upper-right corner of the Expression section. From the menu, choose Component I (compl)>Magnetic Fields> Electric>mf.normE - Electric field norm - V/m.
- 4 Click the **Zoom Extents** button in the **Graphics** toolbar.

Observe that the electric field is slightly shielded by the plasma. This is due to the skin effect in the plasma. As the electron number density increases, the plasma tends to shield itself from the electric field.

Excited Argon Number Density

- I In the Model Builder window, right-click High Frequency Electric Field and choose Duplicate.
- 2 In the Settings window for 2D Plot Group, type Excited Argon Number Density in the Label text field.

Surface 1

- I In the Model Builder window, expand the Excited Argon Number Density node, then click Surface 1.
- 2 In the Settings window for Surface, click Replace Expression in the upper-right corner of the Expression section. From the menu, choose Component I (compl)>Plasma> Number densities>plas.n_wArs - Number density - 1/m3.
- 3 In the Excited Argon Number Density toolbar, click **Plot**.

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

Power Deposition

- I In the Home toolbar, click Add Plot Group and choose 2D Plot Group.
- 2 In the Settings window for 2D Plot Group, type Power Deposition in the Label text field.

Surface 1

- I Right-click Power Deposition and choose Surface.
- 2 In the Settings window for Surface, click Replace Expression in the upper-right corner of the Expression section. From the menu, choose Component I (compl)>Magnetic Fields> Heating and losses>mf.Qrh Volumetric loss density, electric W/m³.
- 3 Locate the Coloring and Style section. Click Change Color Table.
- 4 In the Color Table dialog box, select Thermal>ThermalWave in the tree.
- 5 Click OK.

Selection 1

- I Right-click Surface I and choose Selection.
- **2** Select Domain 3 only.
- 3 In the Power Deposition toolbar, click Plot.
- 4 Click the **Zoom Extents** button in the **Graphics** toolbar.

 The effect of the shielding of the electric field due to the skin depth of the plasma is also

The effect of the shielding of the electric field due to the skin depth of the plasma is also apparent when plotting the power deposition.

Appendix — Geometry Instructions

ADD COMPONENT

In the Home toolbar, click Add Component and choose 2D Axisymmetric.

GEOMETRY I

Polygon I (poll)

- I In the Geometry toolbar, click / Polygon.
- 2 In the Settings window for Polygon, locate the Coordinates section.
- **3** From the **Data source** list, choose **Vectors**.
- 4 In the r text field, type 0.01 0.01 0.14 0.14 0.07 0.07 0 0 0.01.

- 5 In the z text field, type -0.015 -0.025 -0.025 0.08 0.08 0.05 0.05 -0.015 -0.015. 6 Click Build All Objects. Rectangle I (rI) I In the Geometry toolbar, click Rectangle. 2 In the Settings window for Rectangle, locate the Size and Shape section. 3 In the Width text field, type 0.064. 4 In the Height text field, type 0.01. **5** Locate the **Position** section. In the **z** text field, type **0.04**. Rectangle 2 (r2) I In the Geometry toolbar, click Rectangle. 2 In the Settings window for Rectangle, locate the Size and Shape section. 3 In the Width text field, type 0.06. 4 In the Height text field, type 0.03. **5** Locate the **Position** section. In the **z** text field, type **0.05**. Rectangle 3 (r3) I In the Geometry toolbar, click Rectangle. 2 In the Settings window for Rectangle, locate the Size and Shape section. **3** In the **Width** text field, type 0.0825. **4** In the **Height** text field, type 0.0025. **5** Locate the **Position** section. In the **z** text field, type -0.0025. Rectangle 4 (r4) I In the Geometry toolbar, click Rectangle. 2 In the Settings window for Rectangle, locate the Size and Shape section. **3** In the **Width** text field, type **0.05**. **4** In the **Height** text field, type 0.015-0.0025. **5** Locate the **Position** section. In the **z** text field, type -0.015.
- Rectangle 5 (r5)
- I In the Geometry toolbar, click Rectangle.
- 2 In the Settings window for Rectangle, locate the Size and Shape section.
- 3 In the Width text field, type 0.04.

- 4 In the Height text field, type 0.01.
- **5** Locate the **Position** section. In the **r** text field, type **0.01**.
- 6 In the z text field, type -0.025.

Polygon 2 (pol2)

- I In the Geometry toolbar, click / Polygon.
- 2 In the Settings window for Polygon, locate the Coordinates section.
- 3 From the Data source list, choose Vectors.
- 4 In the r text field, type 0.057 0.057 0.0825 0.0825 0.064 0.064 0.057.
- 5 In the z text field, type 0.04 0.034 0.034 0.05 0.05 0.04 0.04.
- 6 Click **Build All Objects**.

Rectangle 6 (r6)

- I In the Geometry toolbar, click Rectangle.
- 2 In the Settings window for Rectangle, locate the Size and Shape section.
- 3 In the Width text field, type 0.003.
- 4 In the Height text field, type 0.003.
- **5** Locate the **Position** section. In the **r** text field, type 0.005.
- 6 In the z text field, type 0.05.

Array I (arrI)

- I In the Geometry toolbar, click Transforms and choose Array.
- 2 Select the object r6 only.
- 3 In the Settings window for Array, locate the Size section.
- 4 In the r size text field, type 5.
- **5** Locate the **Displacement** section. In the **r** text field, type **0.012**.
- 6 Click **Build All Objects**.