
Advanced MOSFETs and Novel Devices

Dr.-Ing. Josef Biba

7. Tutorial & Exercise

Device Simulation

Device Simulation – Overview

SILVACO

ENGINEERED EXCELLENCE

TCAD

<http://www.silvaco.com/products/tcad/index.html>

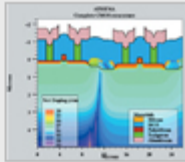
Silvaco provides different **T**echnology **C**omputer **A**ided **D**esign (TCAD)

software packages for modeling semiconductor processes, devices and circuit.

ATHENA

Process Simulation Framework.

Integrates several process simulation modules in a user-friendly environment provided by Silvaco TCAD interactive tools.



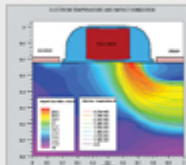
[More...](#)

➤ **Athena** is the process simulator

- Simulates all processes e. g. implantation, oxidation, diffusion, etching etc.

ATLAS

Device simulation framework that enables simulation of the electrical, optical, and thermal behavior of semiconductor devices.



[More...](#)

➤ **Atlas** is the device simulator

- Simulates electrical characteristics e. g. IV, CV of devices

Generating Atlas Input File:

➤ Initial Definitions

- Mesh
- Structure
- Electrodes
- Doping
- Regridding

➤ Models

➤ Solutions

- Initial Guess
- Bias steps
- Bias sweeps

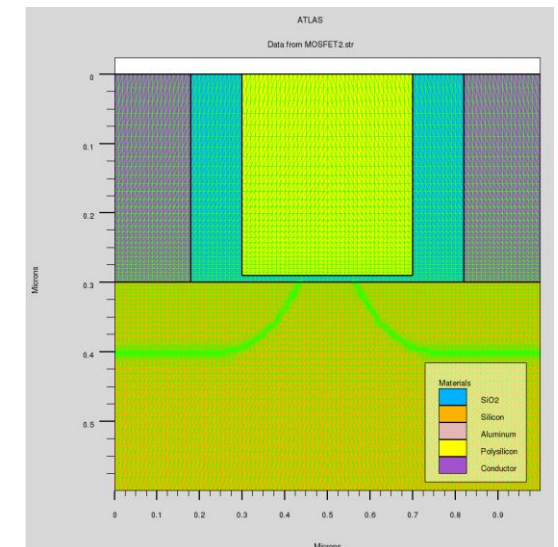
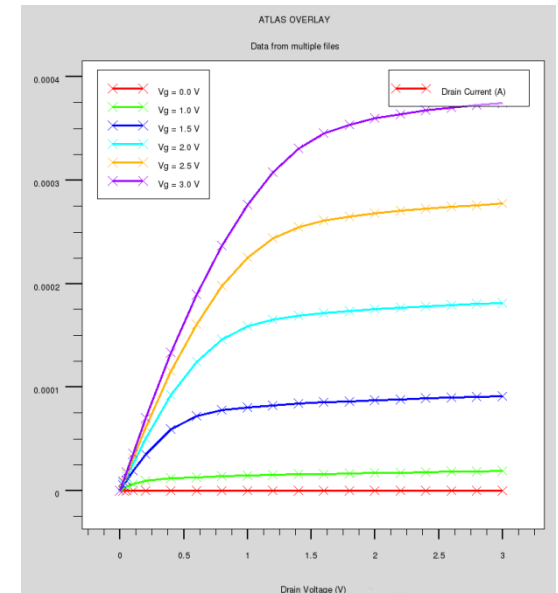
Results:

➤ 1D Plots

- IV Characteristic
- Profiles

➤ 2D Plots

- Structure
- Contour Plots

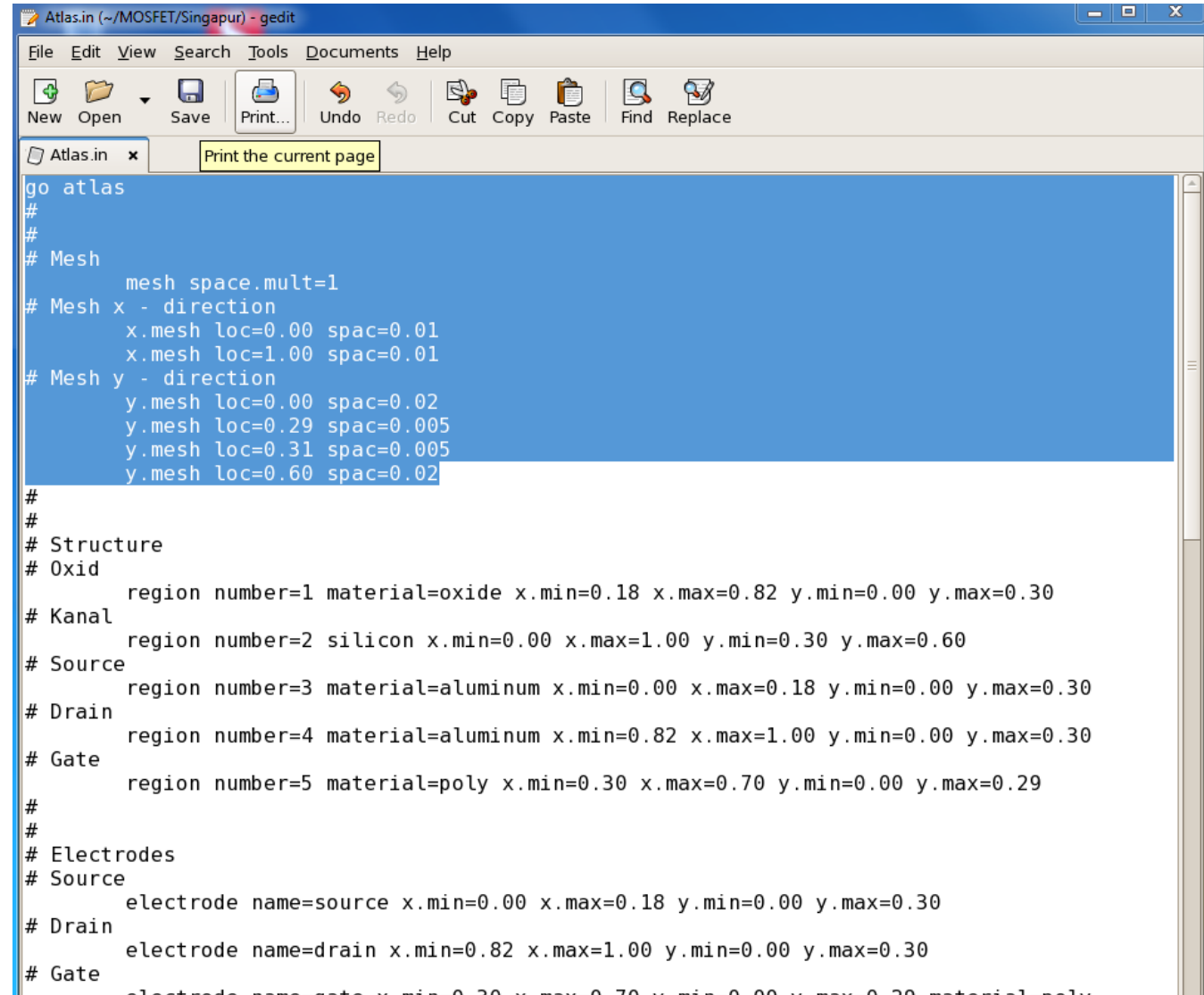


The input file has the following structure:

- The **physical structure** to be simulated
 - Structure from the process simulator
 - Self defined structure
- The **physical models** to be used
 - Choose models for e. g. recombination, tunneling breakdown etc.
- The **numerical methods** for the simulation
- The **bias conditions** for the electrical characteristics
 - IV characteristic, e. g. transfer characteristic
 - CV characteristic

Defining a mesh for Atlas:

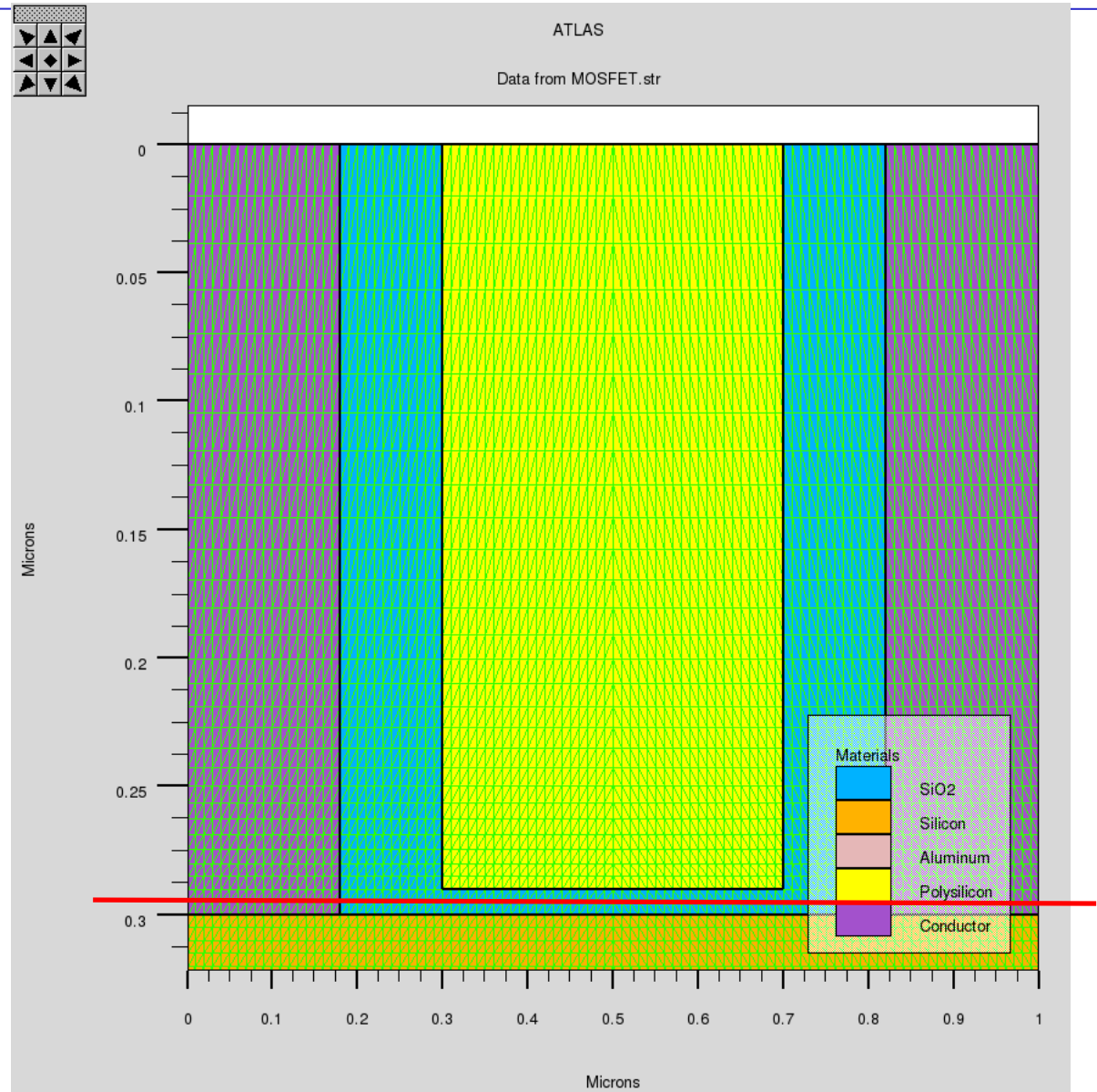
- **Go Atlas** says to use Atlas simulator
- **Mesh space** is a scaling factor for the mesh
- **X.Mesh** defines the mesh in x – direction
- **Y.Mesh** defines the mesh in y - direction



```
Atlas.in (~/.MOSFET/Singapur) - gedit
File Edit View Search Tools Documents Help
New Open Save Print... Undo Redo Cut Copy Paste Find Replace
Atlas.in x Print the current page
go atlas
#
#
# Mesh
    mesh space.mult=1
# Mesh x - direction
    x.mesh loc=0.00 spac=0.01
    x.mesh loc=1.00 spac=0.01
# Mesh y - direction
    y.mesh loc=0.00 spac=0.02
    y.mesh loc=0.29 spac=0.005
    y.mesh loc=0.31 spac=0.005
    y.mesh loc=0.60 spac=0.02
#
#
# Structure
# Oxid
    region number=1 material=oxide x.min=0.18 x.max=0.82 y.min=0.00 y.max=0.30
# Kanal
    region number=2 silicon x.min=0.00 x.max=1.00 y.min=0.30 y.max=0.60
# Source
    region number=3 material=aluminum x.min=0.00 x.max=0.18 y.min=0.00 y.max=0.30
# Drain
    region number=4 material=aluminum x.min=0.82 x.max=1.00 y.min=0.00 y.max=0.30
# Gate
    region number=5 material=poly x.min=0.30 x.max=0.70 y.min=0.00 y.max=0.29
#
#
# Electrodes
# Source
    electrode name=source x.min=0.00 x.max=0.18 y.min=0.00 y.max=0.30
# Drain
    electrode name=drain x.min=0.82 x.max=1.00 y.min=0.00 y.max=0.30
# Gate
    electrode name=gate x.min=0.30 x.max=0.70 y.min=0.00 y.max=0.29 material=poly
```

Defining a mesh for Atlas:

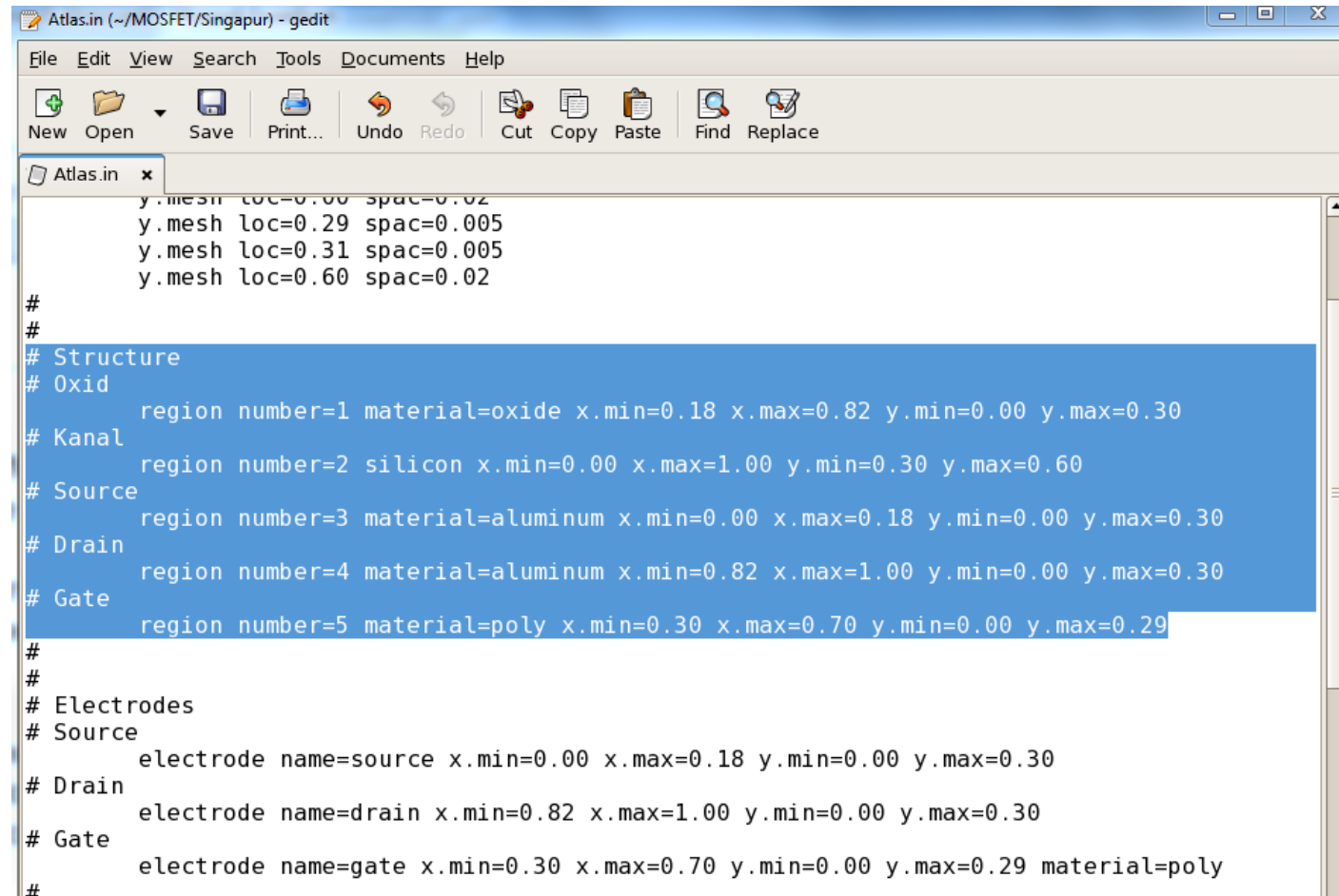
- **Loc** defines the position for a mesh line
- **Spac** defines the space between two mesh lines.
- If different line spacing's were defined, ATLAS adjusts the spacing step by step.
- Use fine mesh where quantities (e. g. doping) changes
- Use coarser mesh where quantities stay constant (e. g. substrate)



Device Simulation – Structure

Defining a structure:

- **Region** defines different layers of the device
- **Number** assigns an identification number to the region
- **Material** defines the material of the region (e. g. silicon, poly, oxide, aluminum etc.)
- **X.min X.max** defines start and end point of the region in x - direction
- **Y.min Y.max** defines start and end point of the region in y - direction



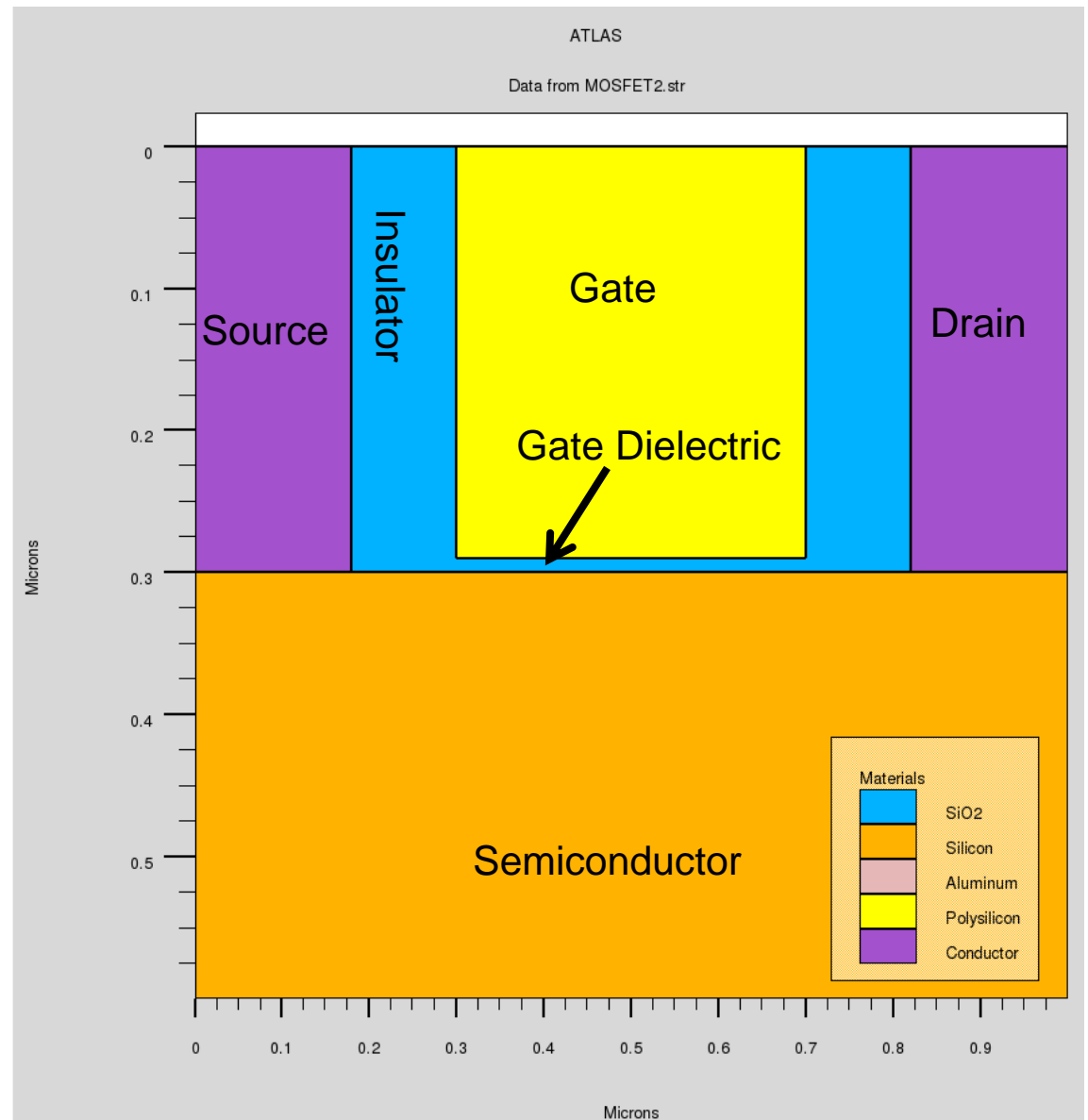
```
Atlas.in (~/.MOSFET/Singapur) - gedit
File Edit View Search Tools Documents Help
New Open Save Print... Undo Redo Cut Copy Paste Find Replace

Atlas.in x
y.mesh loc=0.00 spac=0.02
y.mesh loc=0.29 spac=0.005
y.mesh loc=0.31 spac=0.005
y.mesh loc=0.60 spac=0.02
#
#
# Structure
# Oxid
    region number=1 material=oxide x.min=0.18 x.max=0.82 y.min=0.00 y.max=0.30
# Kanal
    region number=2 silicon x.min=0.00 x.max=1.00 y.min=0.30 y.max=0.60
# Source
    region number=3 material=aluminum x.min=0.00 x.max=0.18 y.min=0.00 y.max=0.30
# Drain
    region number=4 material=aluminum x.min=0.82 x.max=1.00 y.min=0.00 y.max=0.30
# Gate
    region number=5 material=poly x.min=0.30 x.max=0.70 y.min=0.00 y.max=0.29
#
#
# Electrodes
# Source
    electrode name=source x.min=0.00 x.max=0.18 y.min=0.00 y.max=0.30
# Drain
    electrode name=drain x.min=0.82 x.max=1.00 y.min=0.00 y.max=0.30
# Gate
    electrode name=gate x.min=0.30 x.max=0.70 y.min=0.00 y.max=0.29 material=poly
#
```

Device Simulation – Structure

Defining a structure:

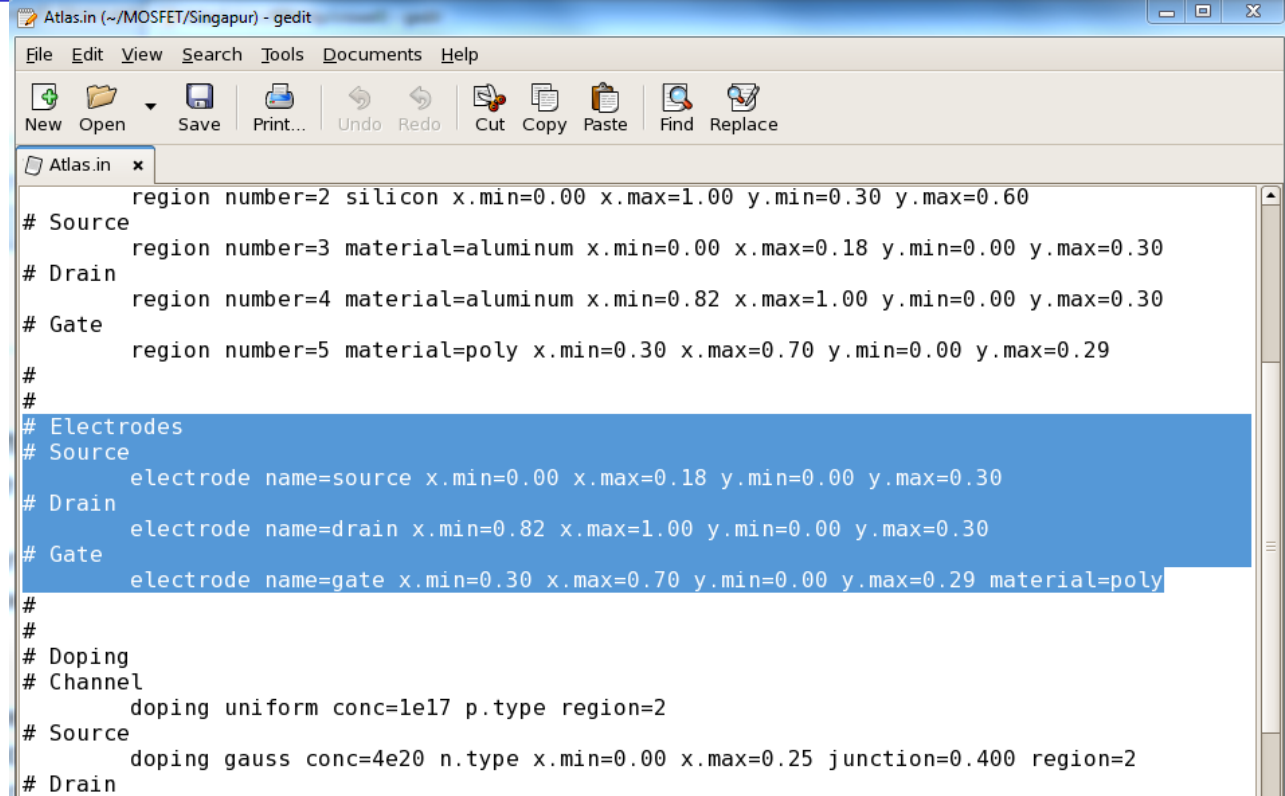
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- **X.min X.max** defines start and end point of the region in x - direction
- **Y.min Y.max** defines start and end point of the region in y - direction



Device Simulation – Electrodes

Defining electrodes:

- **Electrode** defines an electrode for setting bias or current conditions
- **Name** assigns a name to the electrode
- **Material** defines the material of the electrode
- **X.min X.max Y.min Y.max** see region
- All electrical properties are defined in the **contact** statement

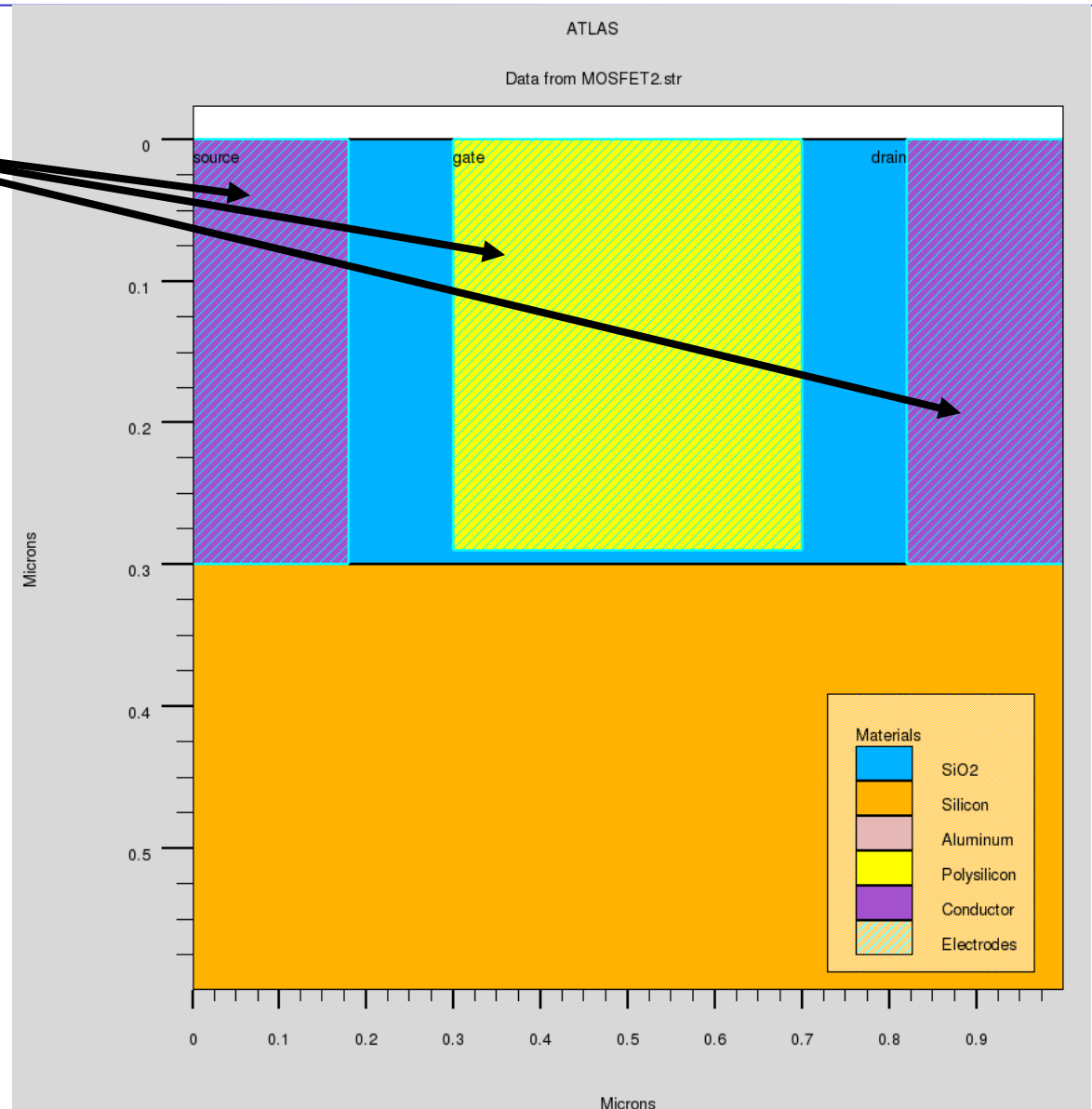


```
Atlas.in (~\MOSFET/Singapur) - gedit
File Edit View Search Tools Documents Help
New Open Save Print... Undo Redo Cut Copy Paste Find Replace
Atlas.in x
region number=2 silicon x.min=0.00 x.max=1.00 y.min=0.30 y.max=0.60
# Source
region number=3 material=aluminum x.min=0.00 x.max=0.18 y.min=0.00 y.max=0.30
# Drain
region number=4 material=aluminum x.min=0.82 x.max=1.00 y.min=0.00 y.max=0.30
# Gate
region number=5 material=poly x.min=0.30 x.max=0.70 y.min=0.00 y.max=0.29
#
#
# Electrodes
# Source
electrode name=source x.min=0.00 x.max=0.18 y.min=0.00 y.max=0.30
# Drain
electrode name=drain x.min=0.82 x.max=1.00 y.min=0.00 y.max=0.30
# Gate
electrode name=gate x.min=0.30 x.max=0.70 y.min=0.00 y.max=0.29 material=poly
#
#
# Doping
# Channel
doping uniform conc=1e17 p.type region=2
# Source
doping gauss conc=4e20 n.type x.min=0.00 x.max=0.25 junction=0.400 region=2
# Drain
```

Device Simulation – Electrodes

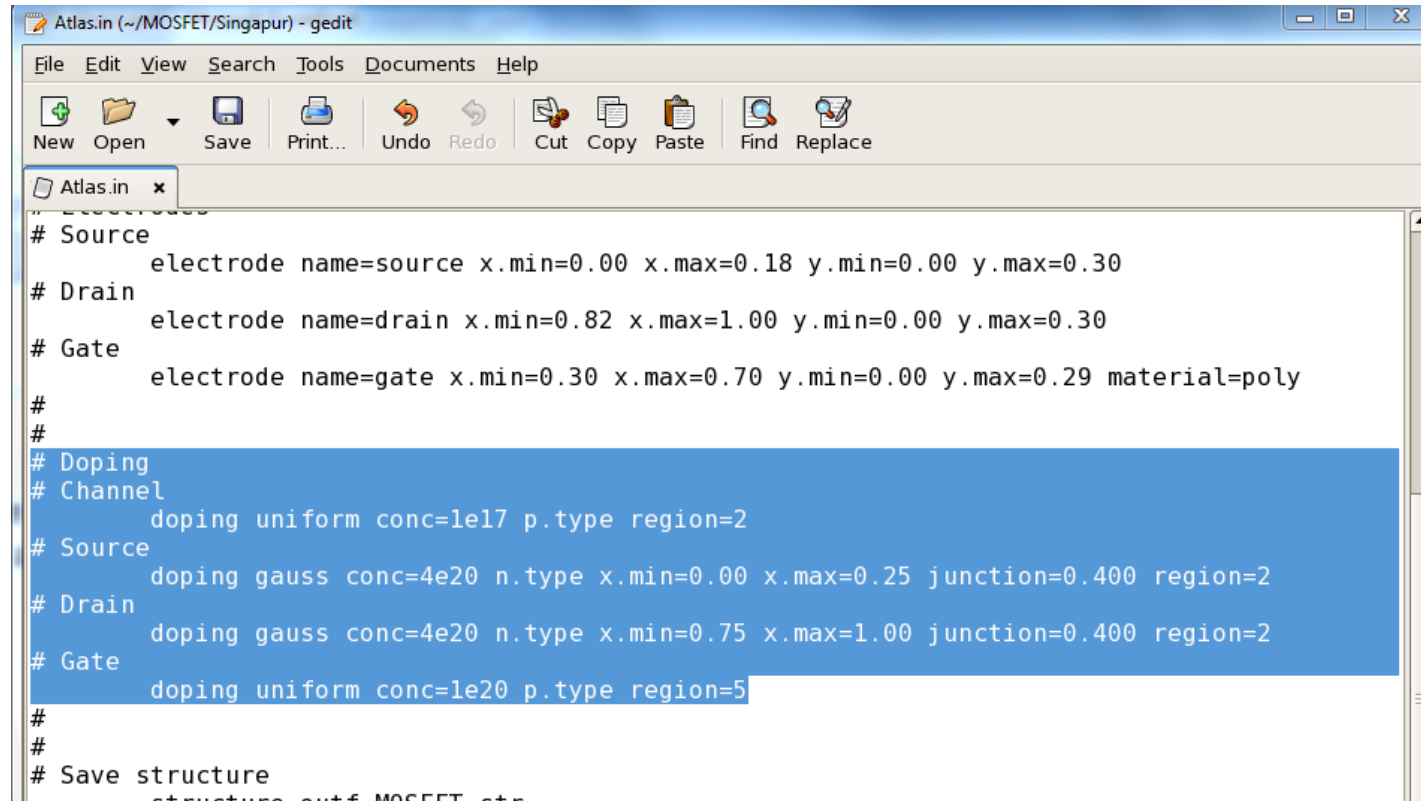
Defining electrodes:

- **Electrode** defines an electrode for setting bias or current conditions
- **Name** assigns a name to the electrode
- **Material** defines the material of the electrode
- **X.min X.max Y.min Y.max** see region
- All electrical properties are defined in the **contact** statement



Defining Doping:

- **Doping** defines the doping of different areas of the device
- **Uniform** makes a constant doping
- **Gauss** defines a doping with Gaussian shape
- **Junction** defines the depth of a Gaussian doping
- **Region** assigns the doping to the previous defined region.

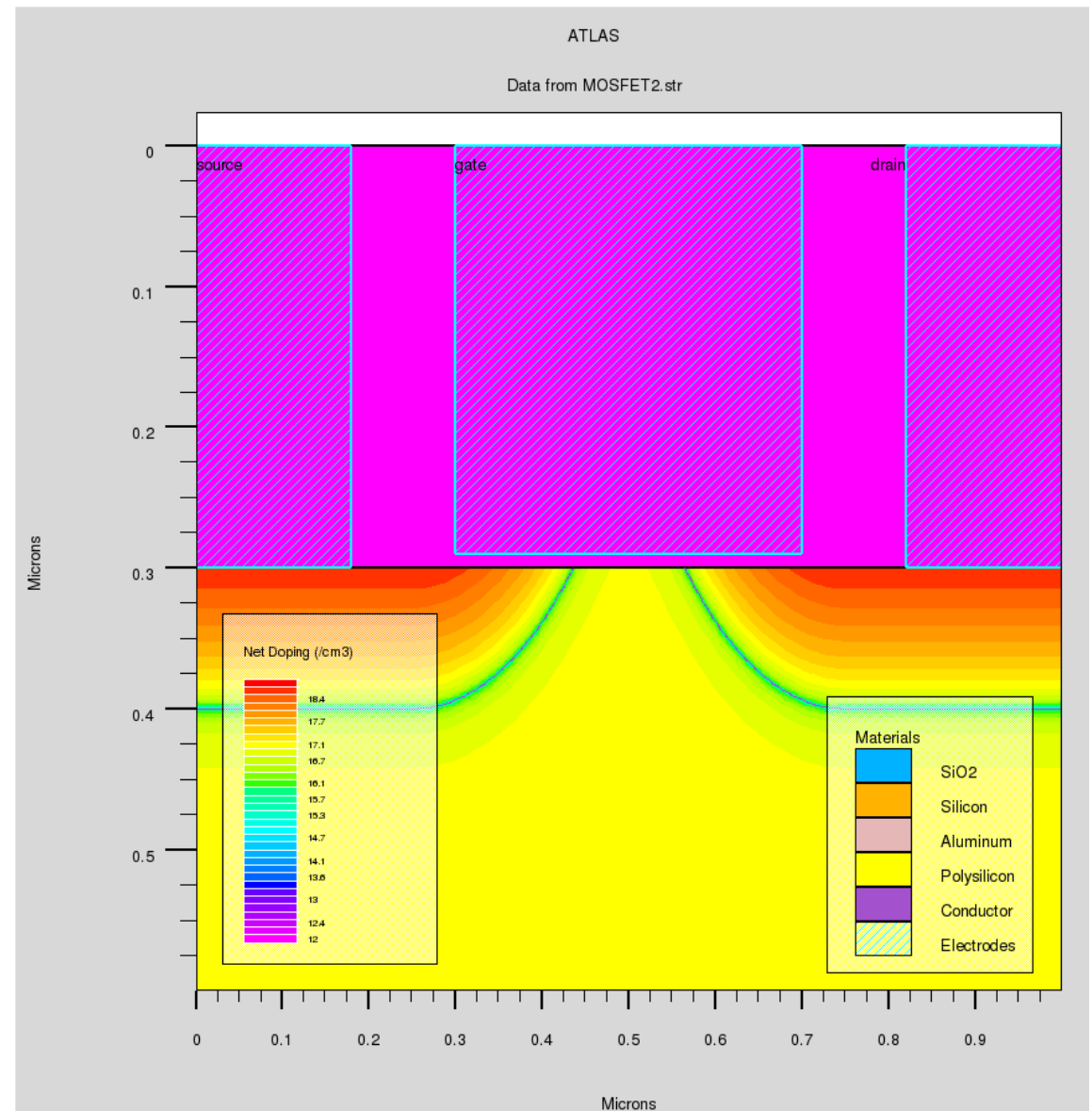
A screenshot of a gedit window titled 'Atlas.in (~/.MOSFET/Singapur) - gedit'. The window shows a text file with configuration parameters for a MOSFET simulation. The parameters are organized into sections: Source, Drain, Gate, Doping, and Save structure. The Doping section is highlighted with a blue background. The text content is as follows:

```
# Source
electrode name=source x.min=0.00 x.max=0.18 y.min=0.00 y.max=0.30
# Drain
electrode name=drain x.min=0.82 x.max=1.00 y.min=0.00 y.max=0.30
# Gate
electrode name=gate x.min=0.30 x.max=0.70 y.min=0.00 y.max=0.29 material=poly
#
#
# Doping
# Channel
doping uniform conc=1e17 p.type region=2
# Source
doping gauss conc=4e20 n.type x.min=0.00 x.max=0.25 junction=0.400 region=2
# Drain
doping gauss conc=4e20 n.type x.min=0.75 x.max=1.00 junction=0.400 region=2
# Gate
doping uniform conc=1e20 p.type region=5
#
#
# Save structure
structure out f=MOSFET.ctr
```

Device Simulation – Doping

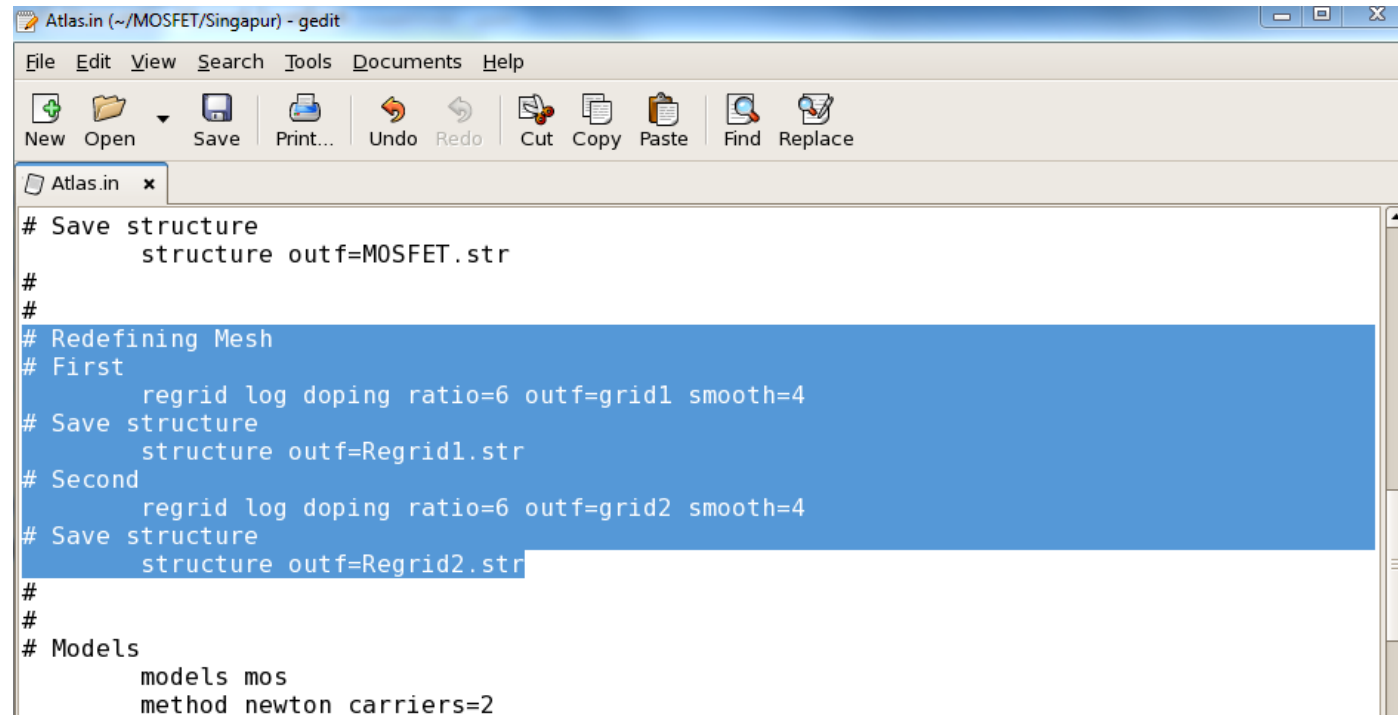
Defining Doping:

- **Doping** defines the doping of different areas of the device
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- **Junction** defines the depth of a Gaussian doping
- **Region** assigns the doping to the previous defined region.



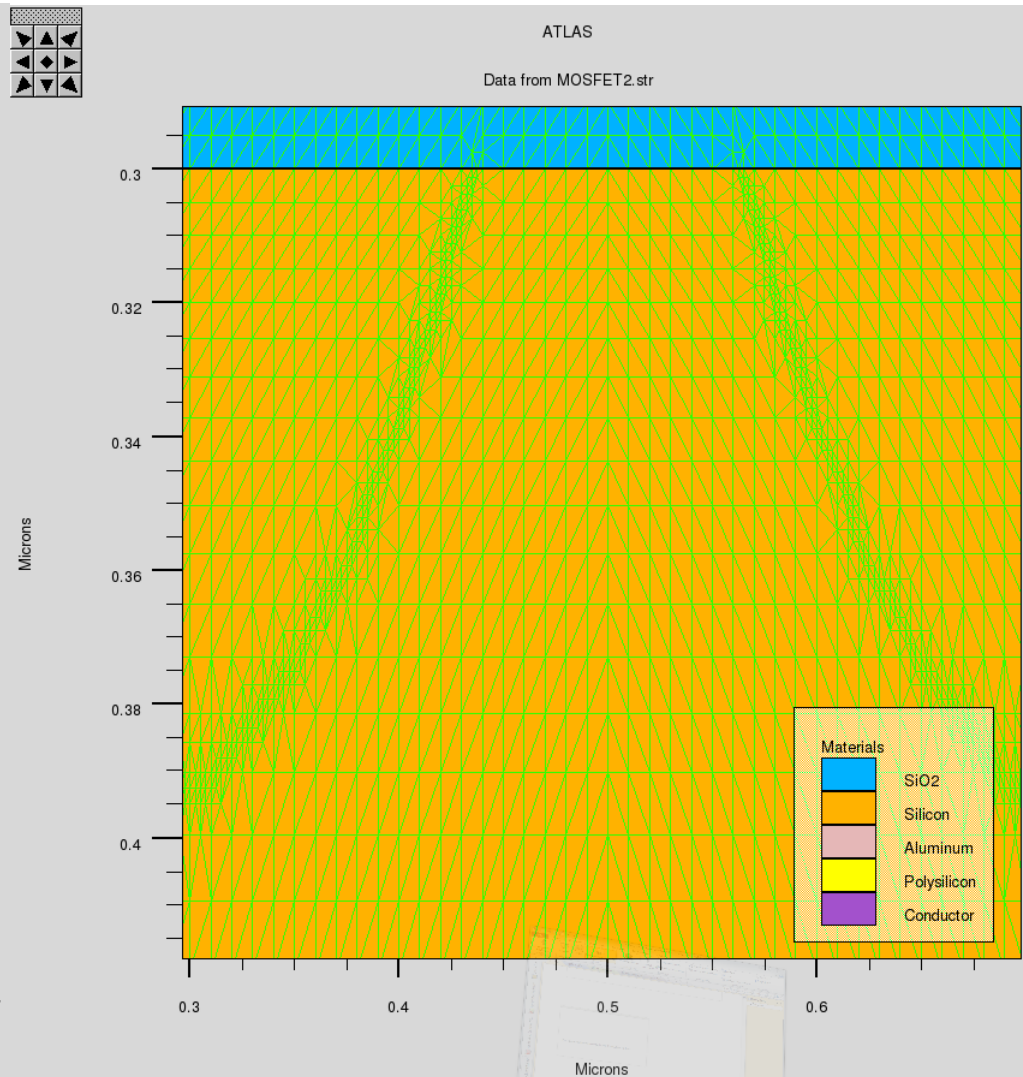
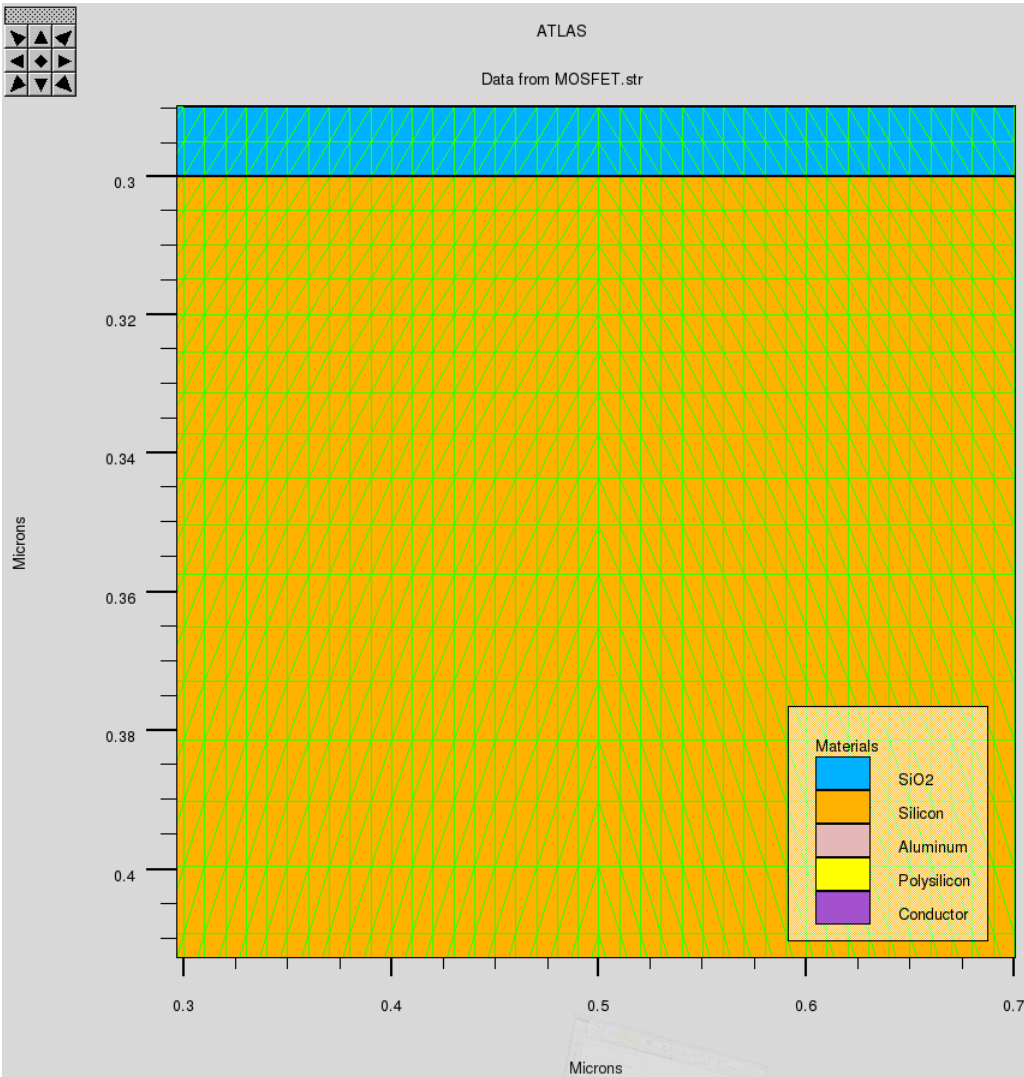
Defining Doping:

- **Regrid** redefines the mesh according to a specific quantity
- **Log** uses a logarithmic refinement scale
- **Doping** sets doping as refinement quantity
- **Ratio** defines the maximum allowed variance across one element
- **Smooth** reduces number of obtuse triangles

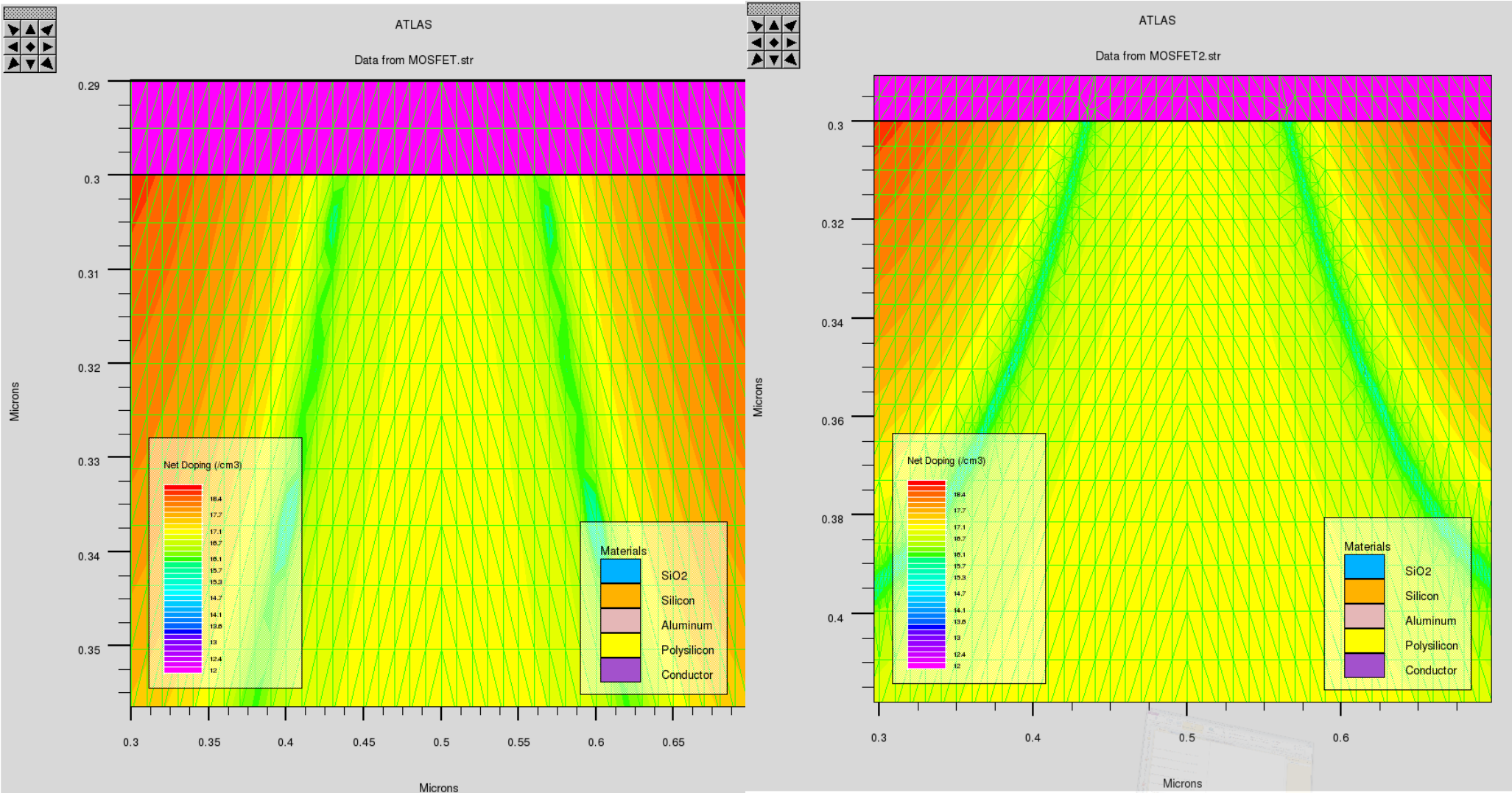


```
Atlas.in (~/.MOSFET/Singapur) - gedit
File Edit View Search Tools Documents Help
New Open Save Print... Undo Redo Cut Copy Paste Find Replace
Atlas.in x
# Save structure
    structure outf=MOSFET.str
#
#
# Redefining Mesh
# First
    regrid log doping ratio=6 outf=grid1 smooth=4
# Save structure
    structure outf=Regrid1.str
# Second
    regrid log doping ratio=6 outf=grid2 smooth=4
# Save structure
    structure outf=Regrid2.str
#
#
# Models
    models mos
    method newton carriers=2
```

Device Simulation – Regrid



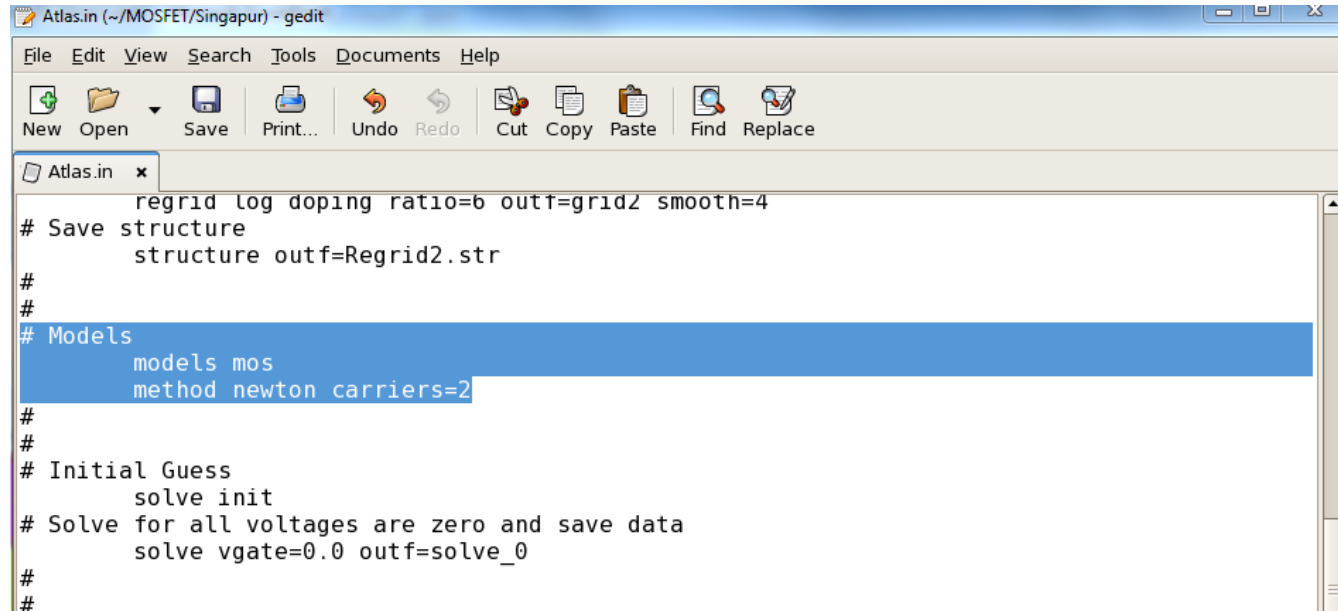
Device Simulation – Regrid



Device Simulation – Models / Methods

Selecting Models / Method:

- **Models** selecting physical models for carrier statistics, mobility, recombination, impact ionization, tunneling
- **Mos** selects a group of predefined models for a special case
- **Method** selects the numerical technique for calculating solutions
- **Carriers** defines if electrons holes, both or none are considered

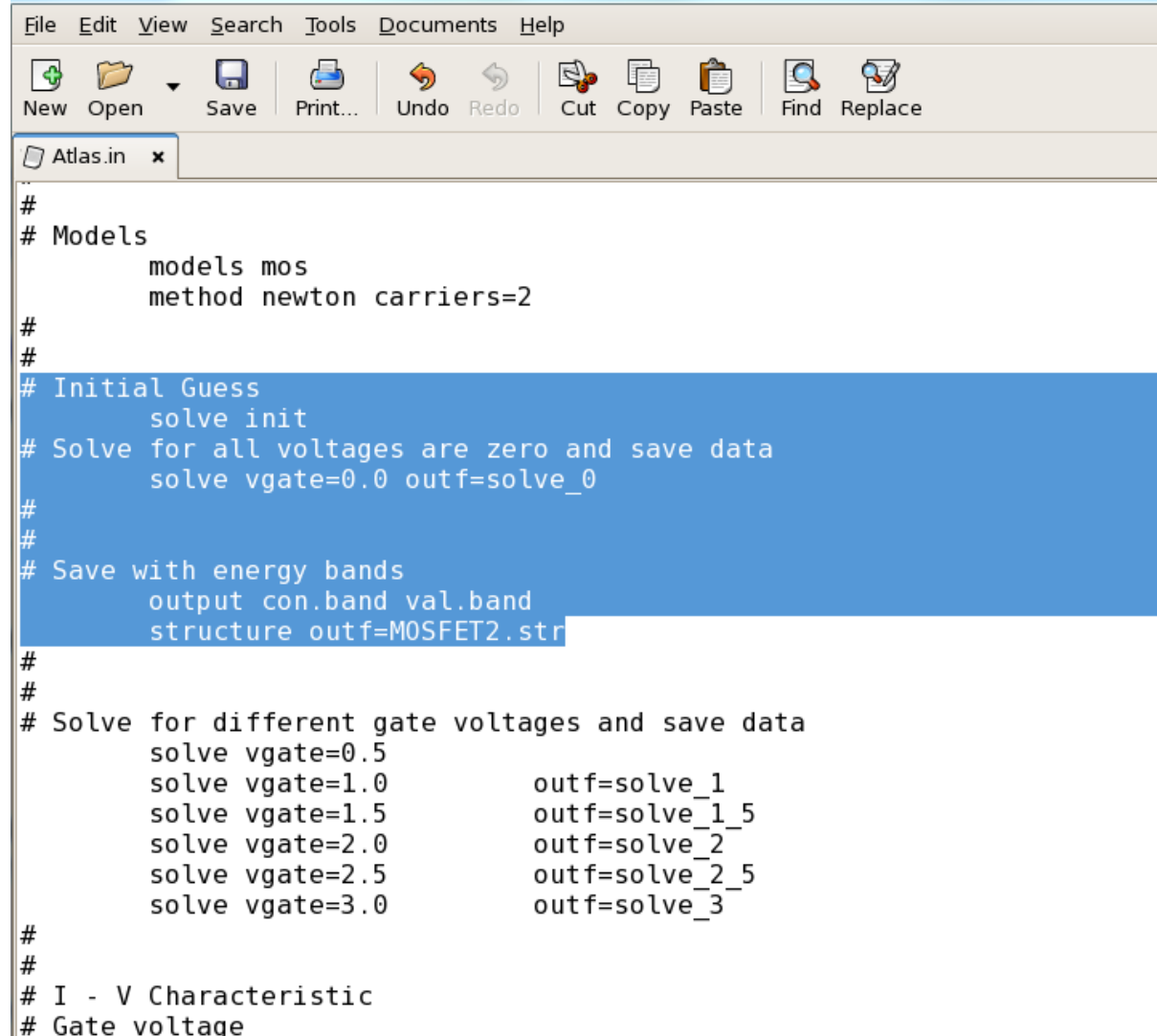


```
Atlas.in (~/.MOSFET/Singapur) - gedit
File Edit View Search Tools Documents Help
New Open Save Print... Undo Redo Cut Copy Paste Find Replace
Atlas.in x
    regrid log doping ratio=6 outf=grid2 smooth=4
# Save structure
    structure outf=Regrid2.str
#
#
# Models
    models mos
    method newton carriers=2
#
#
# Initial Guess
    solve init
# Solve for all voltages are zero and save data
    solve vgate=0.0 outf=solve_0
#
#
```


Device Simulation – Initial Guess / Save Structure

Initial Guess and save structure:

- **Solve** performs a solution for a given bias point
- **Init** sets all electrodes to 0 V
- **Vgate** defines bias for the former defined electrode “gate”
- **Outf** saves data to file
- **Output** specifies data stored in a structure file
- **Structure** writes mesh and solution informations

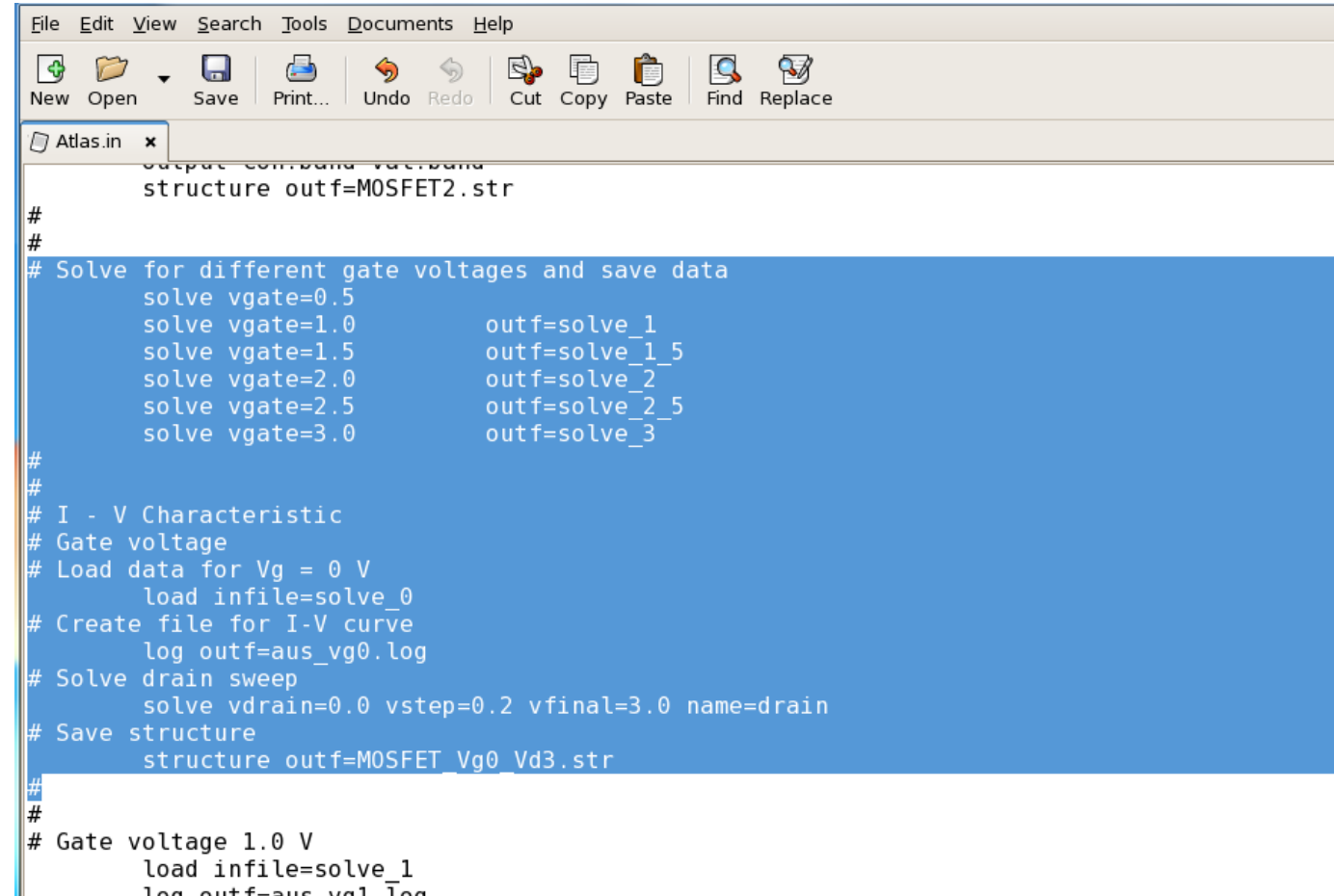


```
File Edit View Search Tools Documents Help
New Open Save Print... Undo Redo Cut Copy Paste Find Replace
Atlas.in x
#
# Models
    models mos
    method newton carriers=2
#
#
# Initial Guess
    solve init
# Solve for all voltages are zero and save data
    solve vgate=0.0 outf=solve_0
#
#
# Save with energy bands
    output con.band val.band
    structure outf=MOSFET2.str
#
#
# Solve for different gate voltages and save data
    solve vgate=0.5
    solve vgate=1.0      outf=solve_1
    solve vgate=1.5      outf=solve_1_5
    solve vgate=2.0      outf=solve_2
    solve vgate=2.5      outf=solve_2_5
    solve vgate=3.0      outf=solve_3
#
#
# I - V Characteristic
# Gate voltage
```

Device Simulation – Bias step / sweep

Bias step / sweep:

- Solve for different gate voltages and save result
- **Load** loads previous saved data specified by **Infile**
- **Log** saves all electrode characteristics to a file
- **Vdrain** sets drain voltage to zero
- **Vstep** defines steps for the voltage sweep
- **Vfinal** sets the final voltage for the sweep

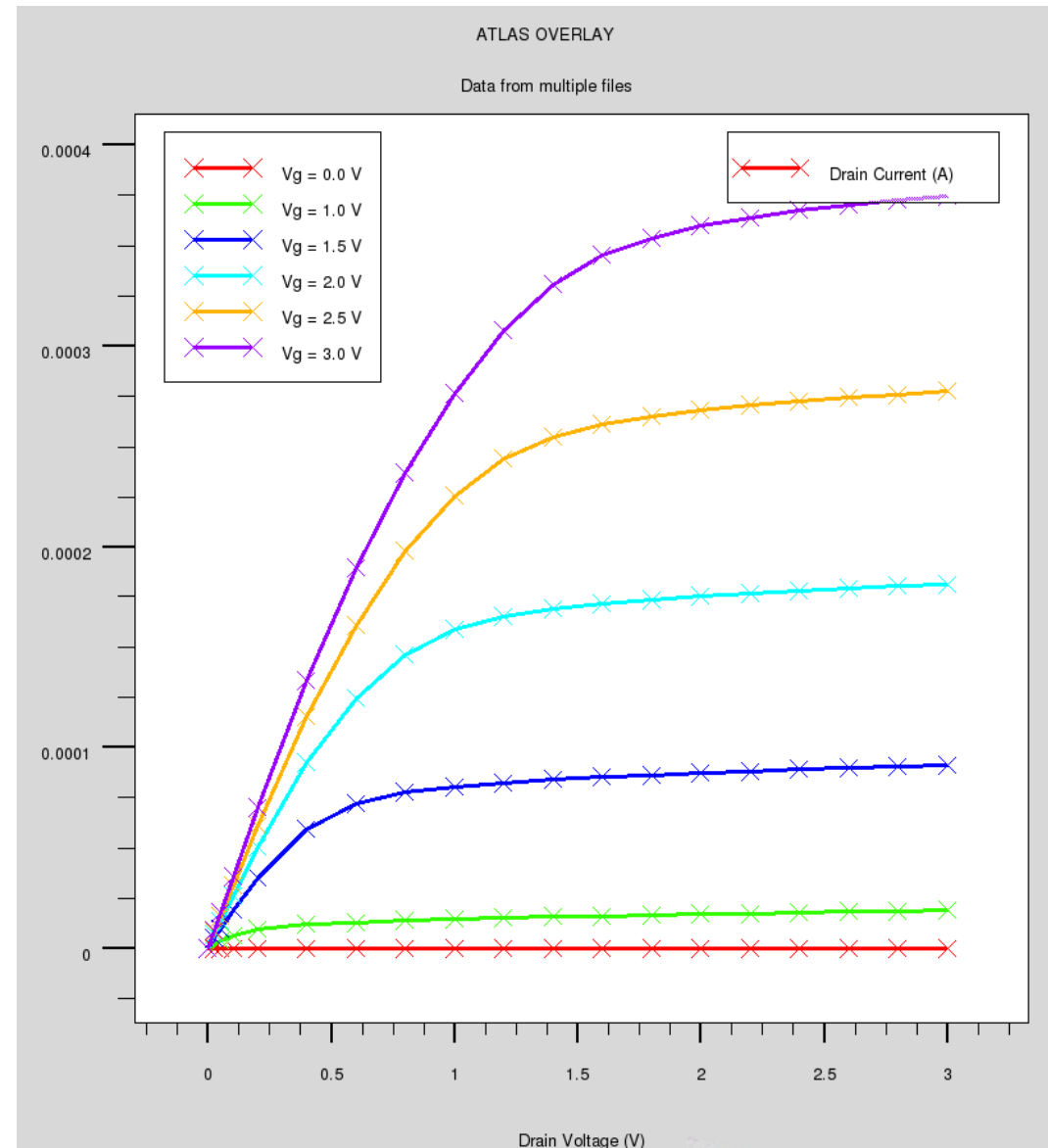


```
File Edit View Search Tools Documents Help
New Open Save Print... Undo Redo Cut Copy Paste Find Replace
Atlas.in x
#
#
# Solve for different gate voltages and save data
solve vgate=0.5
solve vgate=1.0      outf=solve_1
solve vgate=1.5      outf=solve_1_5
solve vgate=2.0      outf=solve_2
solve vgate=2.5      outf=solve_2_5
solve vgate=3.0      outf=solve_3
#
#
# I - V Characteristic
# Gate voltage
# Load data for Vg = 0 V
load infile=solve_0
# Create file for I-V curve
log outf=aus_vg0.log
# Solve drain sweep
solve vdrain=0.0 vstep=0.2 vfinal=3.0 name=drain
# Save structure
structure outf=MOSFET_Vg0_Vd3.str
#
#
# Gate voltage 1.0 V
load infile=solve_1
log outf=aus_vg1.log
```

- **Name** is the name of the swept electrode

1D Plots:

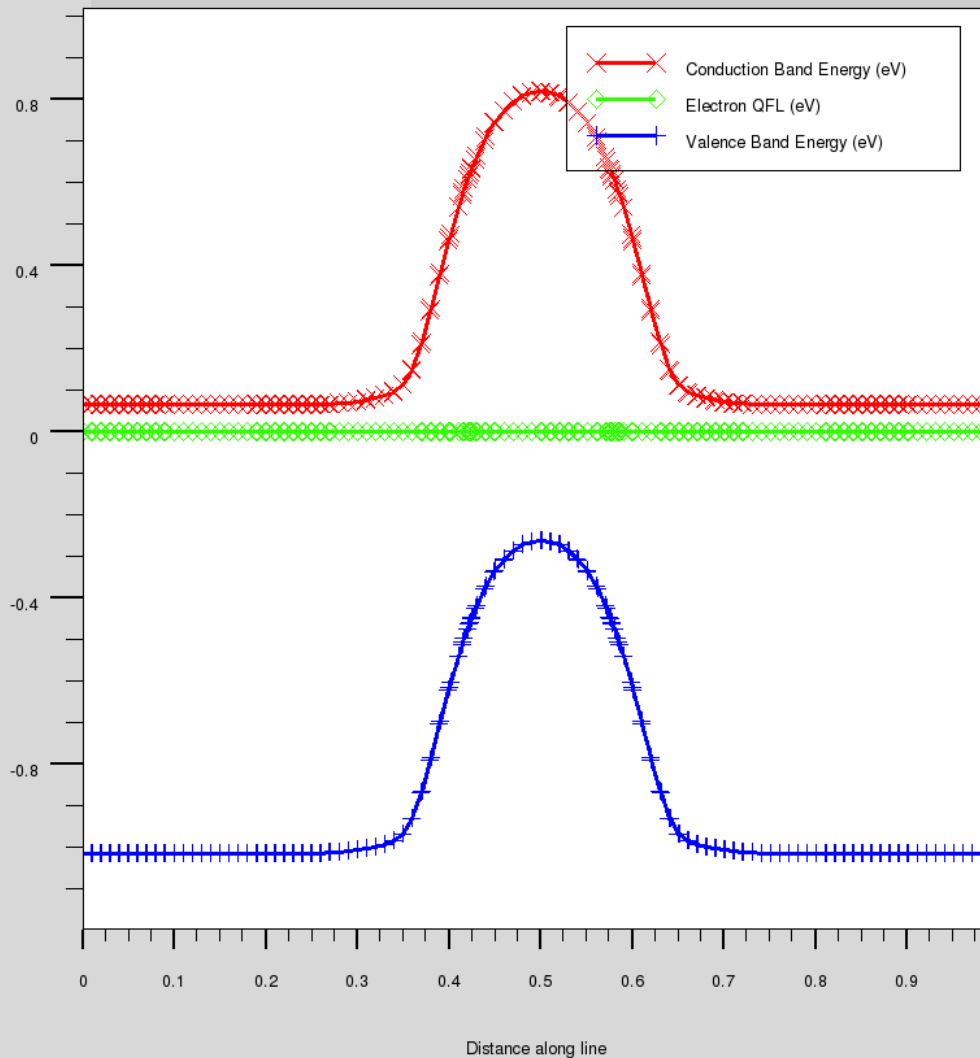
- Simulated output characteristic for different applied gate voltages
- In this case the current is not A! It is A / μm !
- Other 1D plots:
 - Doping
 - Acceptor / Donor concentrations
 - Currents
 - Quasi Fermi Levels
 - Conduction / Valance Bands
 - Electric field



Device Simulation – 1D Plots MOSFET Band Diagram

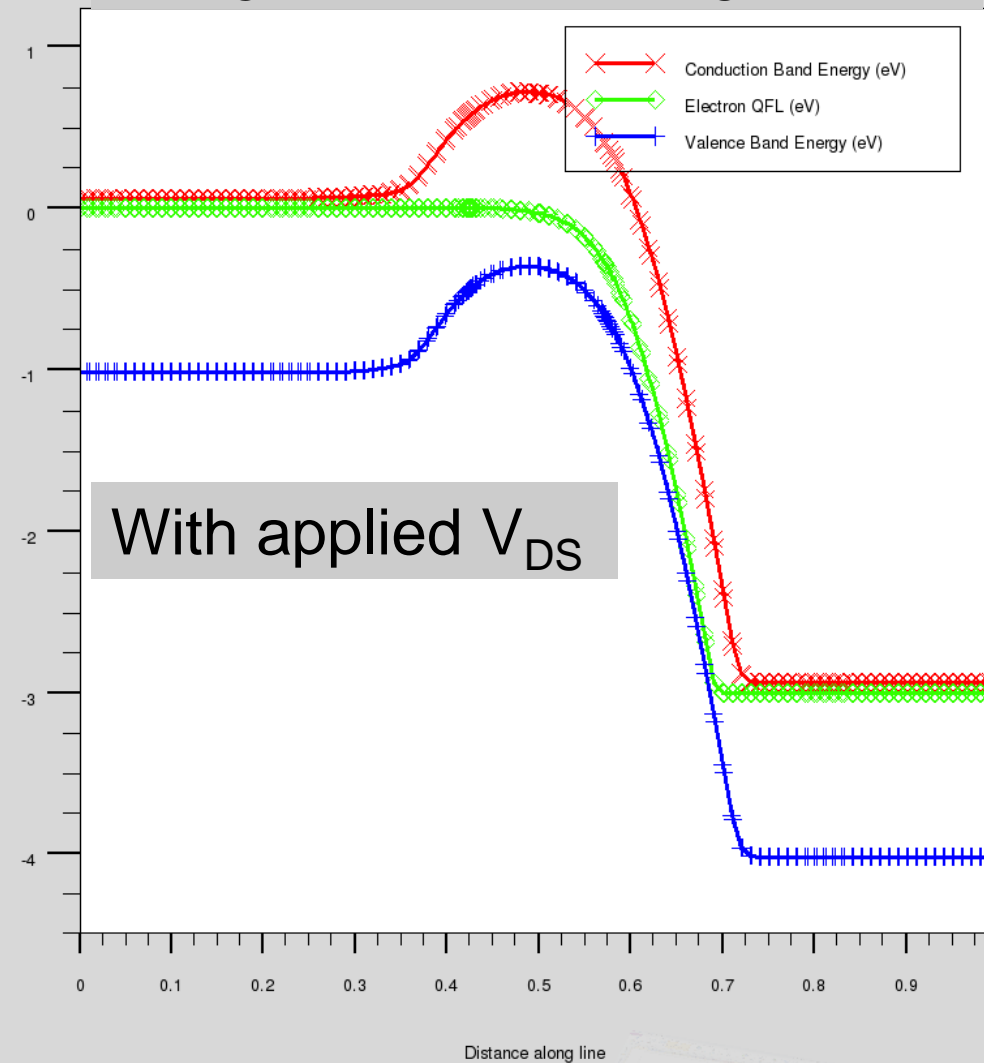
$$V_G = 0V$$

$$V_{DS} = 0V$$

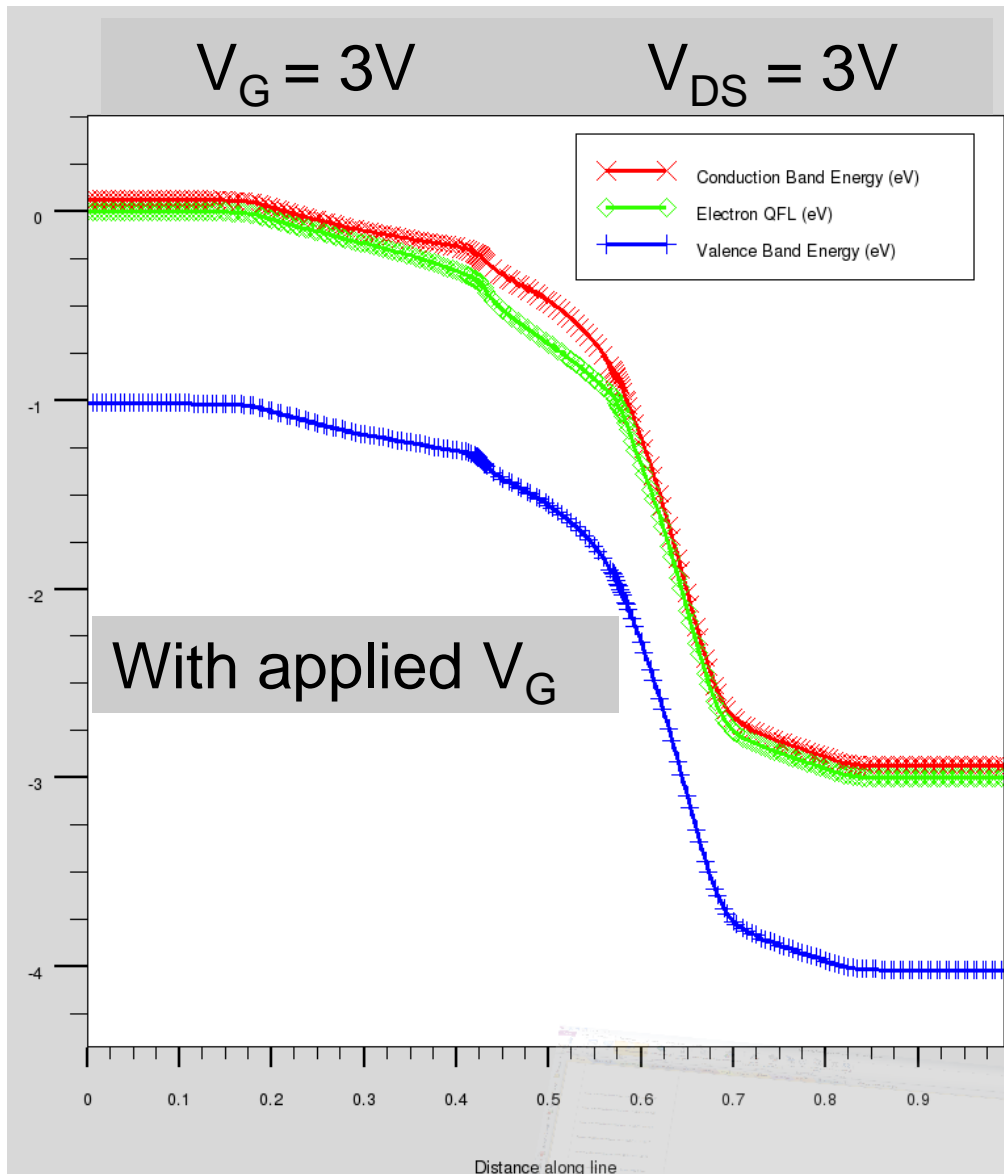


$$V_G = 0V$$

$$V_{DS} = 3V$$



Device Simulation – 1D Plots MOSFET Band Diagram



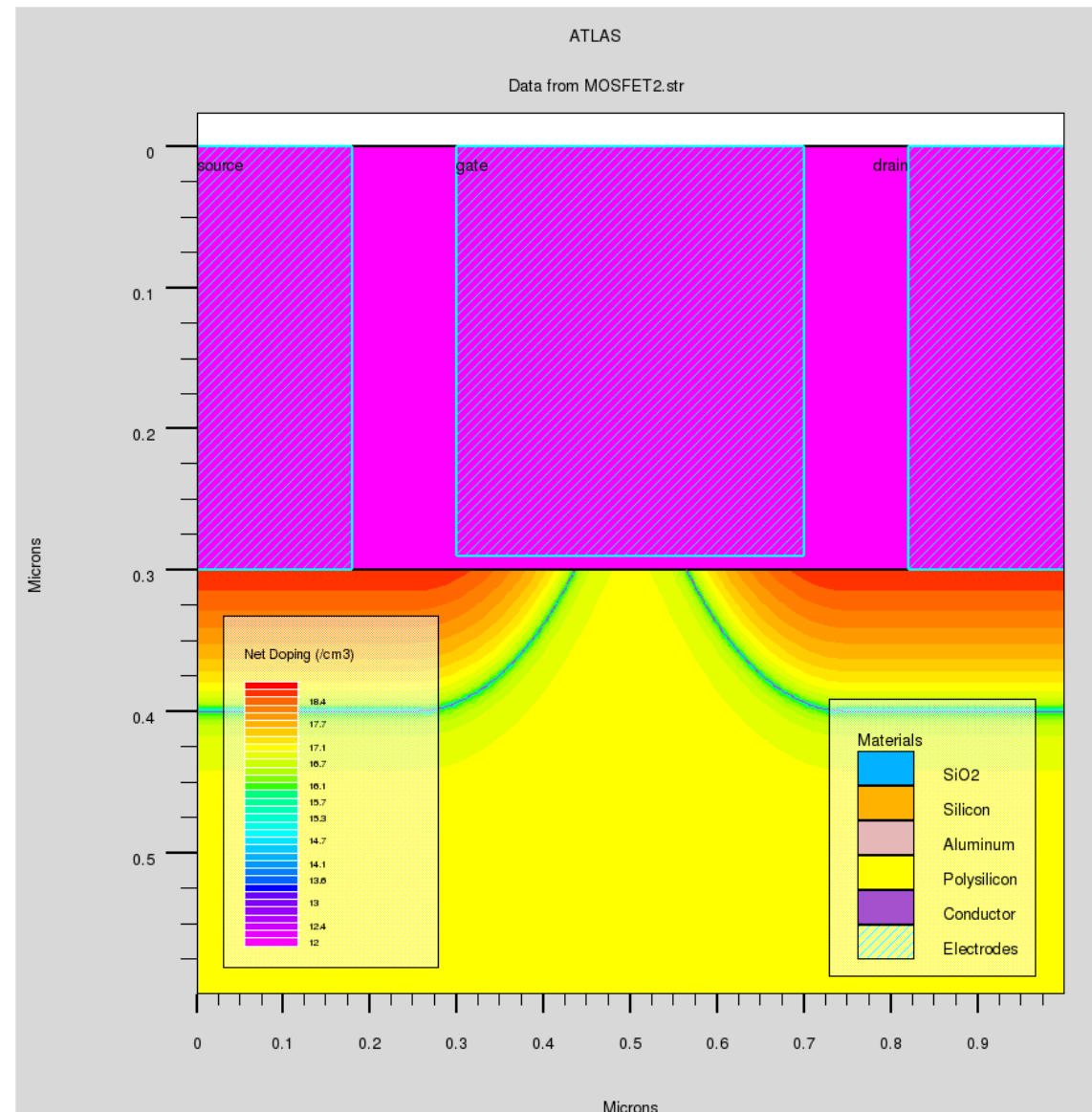
1D Plots:

- Simulated band diagrams
- Barrier height reduced
- Compare to second Tutorial

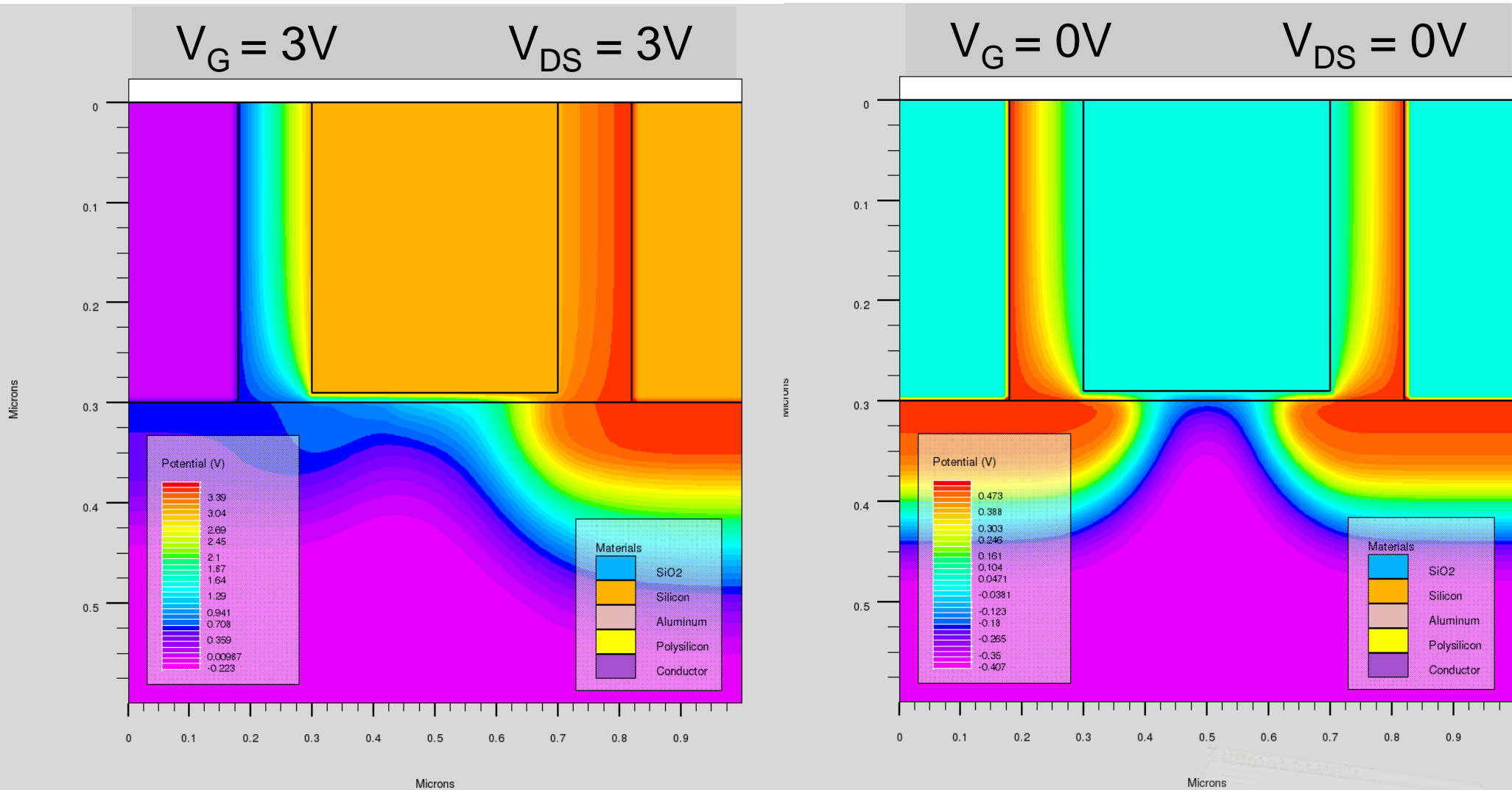
Device Simulation – 2D Plots

2D Plots:

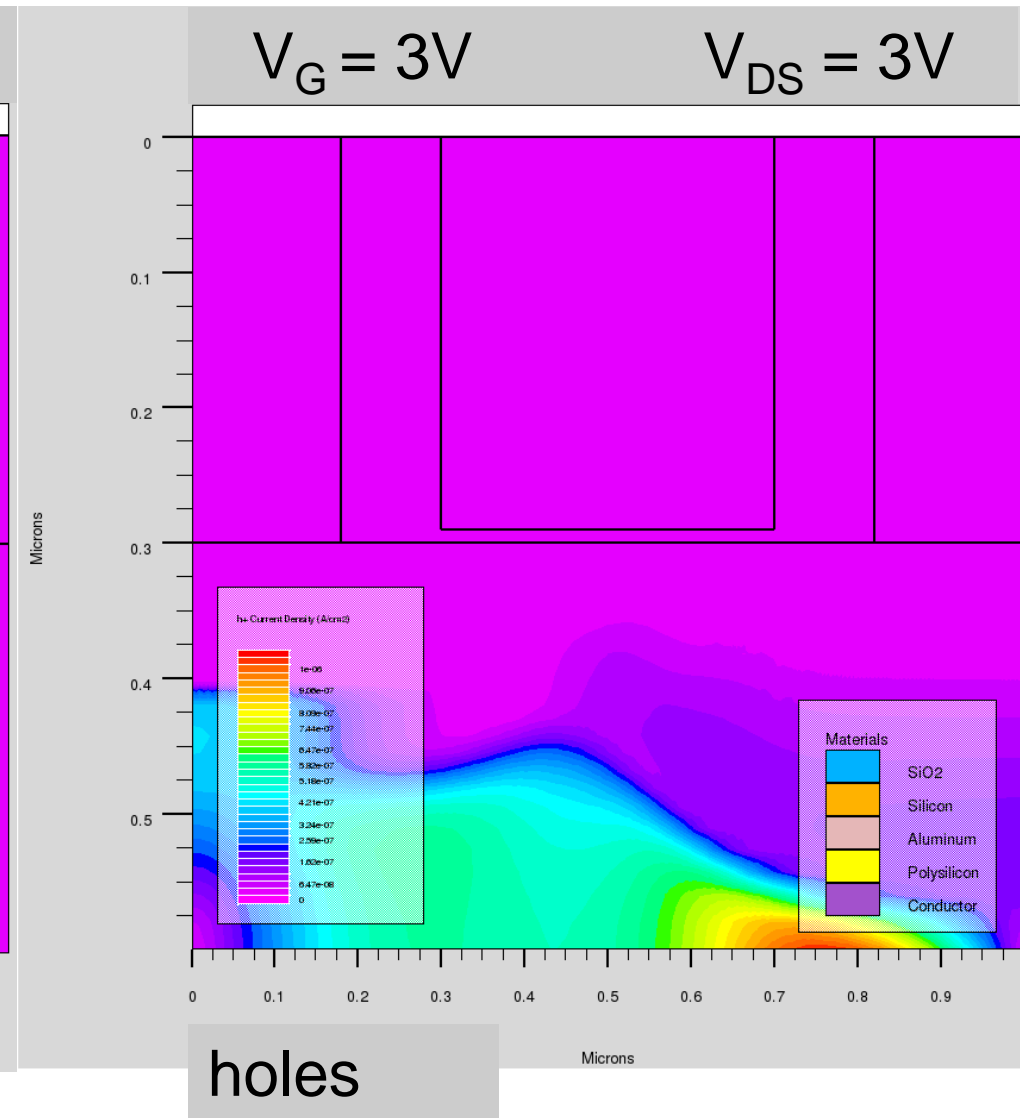
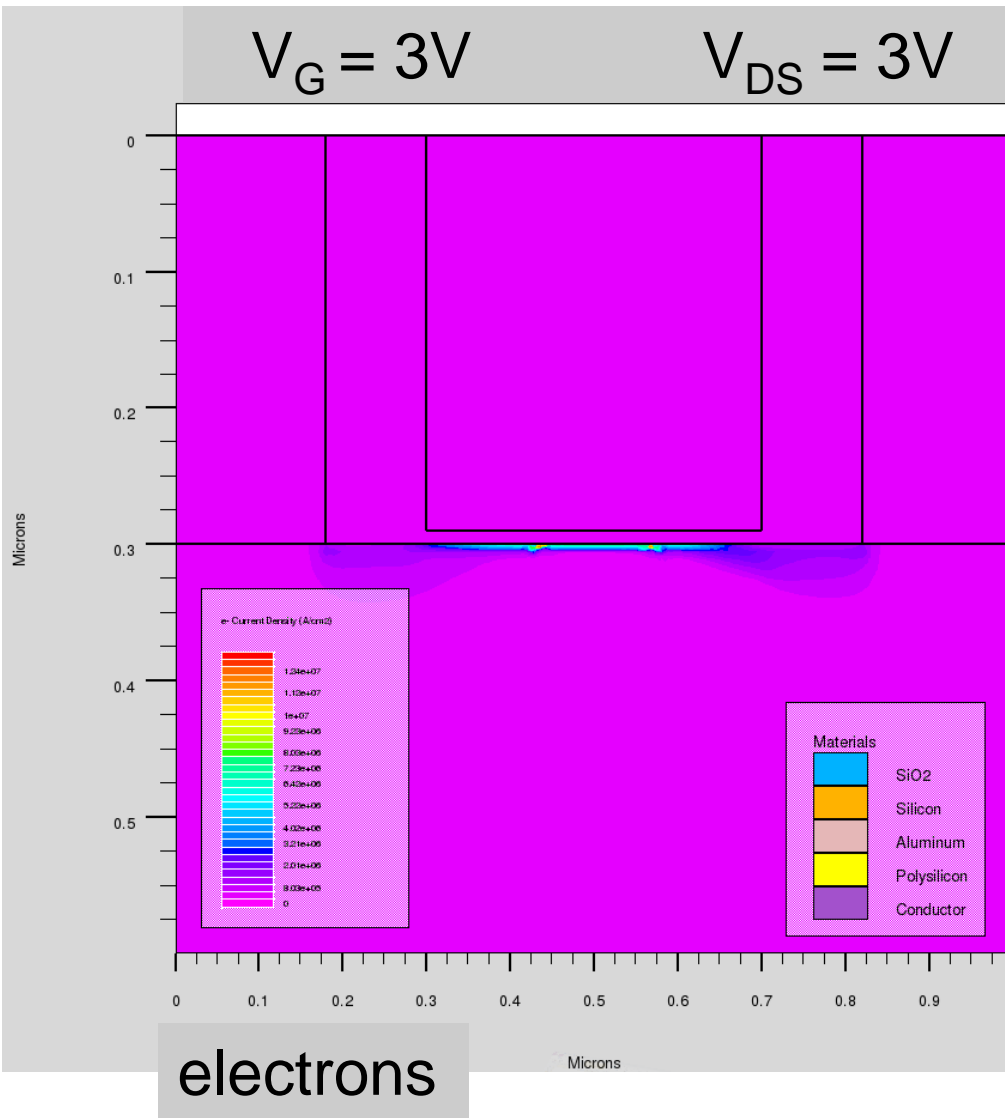
- Simulated doping profile
- Other 2D plots:
 - Doping
 - Acceptor / Donor concentrations
 - Currents
 - Quasi Fermi Levels
 - Conduction / Valance Bands
 - Electric field



Device Simulation – 2D Plots Potential



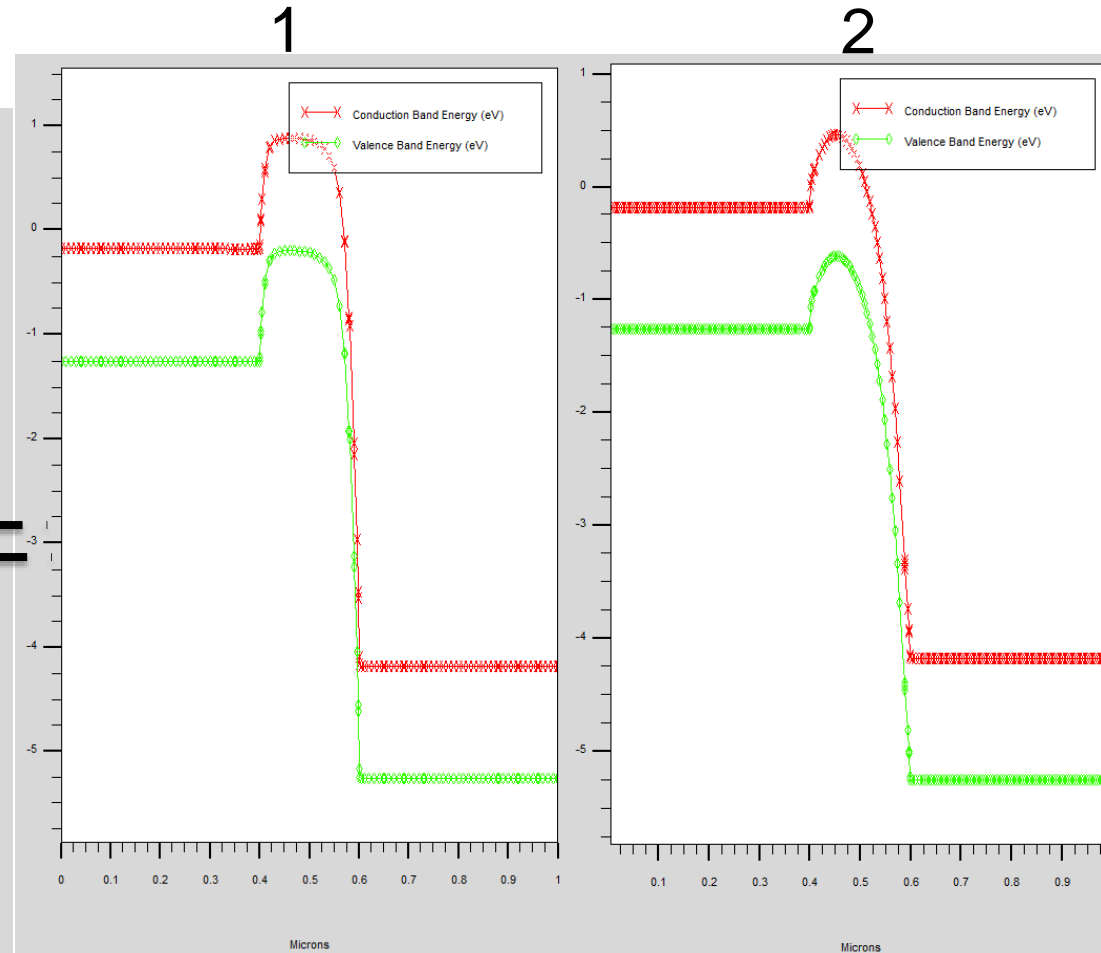
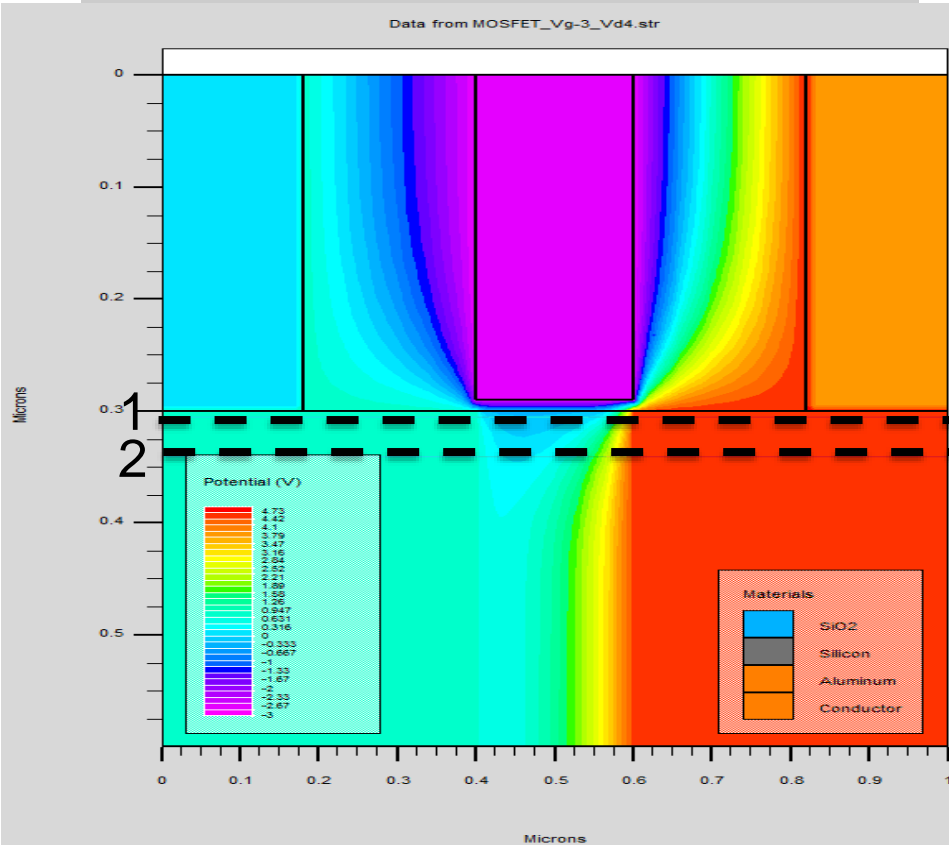
Device Simulation – 2D Plots Current Density



Device Simulation – Short Channel

$$V_G = -3V$$

$$V_{DS} = 3V$$



Transistor switched off

SILVACO

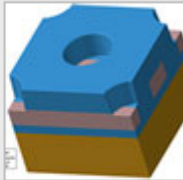
ENGINEERED EXCELLENCE

TCAD

<http://www.silvaco.com/products/tcad/index.html>

Victory Process

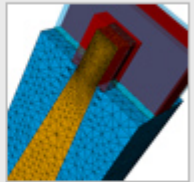
3D Process Simulator. Includes complete process flow core simulator and three advanced simulation modules: Monte Carlo Implant, Advanced Diffusion and Oxidation, and Physical Etch and Deposit.



[More...](#)

Victory Device

General-purpose 3D device simulator using a tetrahedral meshing engine for fast and accurate simulation of complex 3D geometries.



[More...](#)

- **Victory** is the process and device simulator in 3D
 - Simulates all processes e. g. implantation, oxidation, diffusion, etching etc.
 - Simulates electrical characteristics e. g. IV, CV of devices
 - Useful for FinFET calculations

➤ **Important:** Always choose the right models