

# PRISMS-Plasticity

## Crystal Plasticity

### Simple tension example -FCC Copper

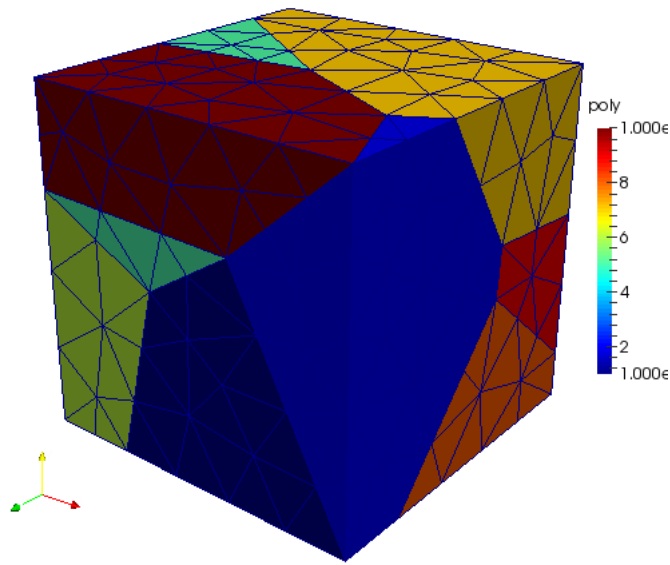


Figure 1: Virtual microstructure (Neper [2])

This is an illustrative example of a simple tension deformation problem. A virtual fcc microstructure generated by Neper[2] was tested with the material parameters of Copper which were obtained from [1]

#### Input Parameters

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```
/*FE parameters*/
#define feOrder 1 // Basis function interpolation order (1-linear)
#define quadOrder 2 // Quadrature point order n^3 (2->8 quadrature points)

/*Mesh parameters*/
//Set the length of the domain in all three dimensions
//Each axes spans from zero to the specified length
#define spanX 1.0
#define spanY 1.0
#define spanZ 1.0
// The number of elements in each direction is 2^(refineFactor) * subdivisions
```

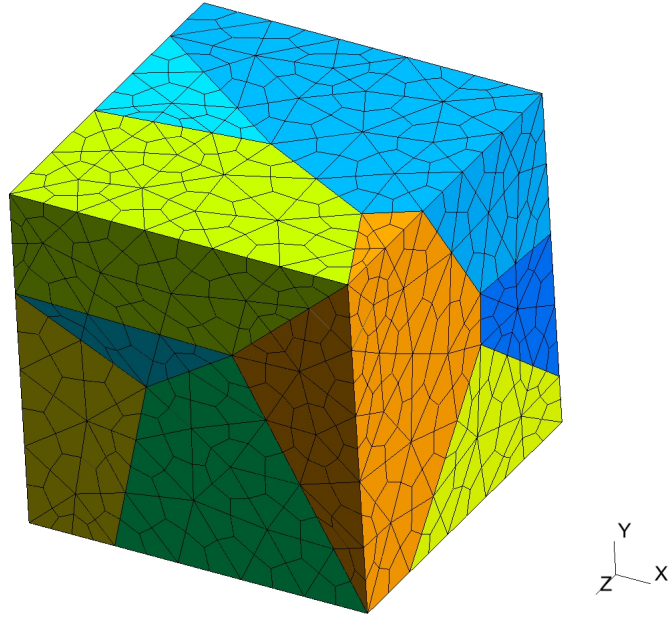


Figure 2: Input microstructure (Hex mesh)

```
// For optimal performance, use meshRefineFactor primarily to determine the
// element size
#define subdivisionsX 1
#define subdivisionsY 1
#define subdivisionsZ 1
#define meshRefineFactor 3 // 2^n*2^n*2^n elements(3->8*8*8 =512 elements)
#define writeMeshToEPS true //Only written for serial runs and if number of
// elements < 10000
#define readExternalMeshes true

/*Solution output parameters*/
#define writeOutput true // flag to write output vtu and pvtu files
#define outputDirectory "."
#define skipOutputSteps 0
#define output_Eqv_strain true
#define output_Eqv_stress true
#define output_Grain_ID true

/*Solver parameters*/
#define linearSolverType PETScWrappers::SolverCG // Type of linear solver
#define totalNumIncrements 100 // No. of increments
#define maxLinearSolverIterations 50000 // Maximum iterations for linear solver
#define relLinearSolverTolerance 1.0e-10 // Relative linear solver tolerance
#define maxNonLinearIterations 4 // Maximum no. of non-linear iterations
#define absNonLinearTolerance 1.0e-18 // Non-linear solver tolerance
#define relNonLinearTolerance 1.0e-3 // Relative non-linear solver tolerance
```

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#define stopOnConvergenceFailure false // Flag to stop problem if convergence
    fails

/*Adaptive time-stepping parameters*/
#define enableAdaptiveTimeStepping false //Flag to enable adaptive time steps
#define adaptiveLoadStepFactor 0.5 // Load step factor
#define adaptiveLoadIncreaseFactor 1.25
#define successiveIncForIncreasingTimeStep 10

//Elastic Parameters
double elasticStiffness[6][6]={170.0e3, 124.0e3, 124.0e3, 0, 0, 0},
                                {124.0e3, 170.0e3, 124.0e3, 0, 0, 0},
                                {124.0e3, 124.0e3, 170.0e3, 0, 0, 0},
                                {0, 0, 0, 75.0e3, 0, 0},
                                {0, 0, 0, 0, 75.0e3, 0},
                                {0, 0, 0, 0, 0, 75.0e3}}; // Elastic
                                Stiffness Matrix -Voigt Notation (MPa)

//Crystal Plasticity parameters
#define numSlipSystems 12 // generally 12 for FCC
#define latentHardeningRatio 1.4 //q1

double initialSlipResistance[numSlipSystems]= {16.0, 16.0, 16.0, 16.0, 16.0,
    16.0, 16.0, 16.0, 16.0, 16.0, 16.0, 16.0}; //CRSS of the slip sytems
double initialHardeningModulus[numSlipSystems]= {180.0, 180.0, 180.0, 180.0,
    180.0, 180.0, 180.0, 180.0, 180.0, 180.0, 180.0, 180.0}; //Hardening moduli
    of slip systems
double powerLawExponent[numSlipSystems]= {2.25, 2.25, 2.25, 2.25, 2.25, 2.25,
    2.25, 2.25, 2.25, 2.25, 2.25, 2.25}; // Power law coefficient
double saturationStress[numSlipSystems]= {148.0, 148.0, 148.0, 148.0, 148.0,
    148.0, 148.0, 148.0, 148.0, 148.0, 148.0, 148.0}; // Saturation stress

//Slip systems files
#define slipDirectionsFile "slipDirections.txt" // Slip Directions File
#define slipNormalsFile "slipNormals.txt" // Slip Normals File

// Crystal Plasticity Constitutive model tolerances (for advanced users)
#define modelStressTolerance 1.0e-6 // Stress tolerance for the yield surface
    (MPa)
#define modelMaxSlipSearchIterations 20 // Maximum no. of active slip search
    iterations
#define modelMaxSolverIterations 10 // Maximum no. of iterations to achieve
    non-linear convergence
#define modelMaxPlasticSlipL2Norm 0.8 // L2-Norm of plastic slip strain-used for
    load-step adaptivity

//Read Input Microstructure
unsigned int numPts[3]={20, 20, 22}; // No. of voxels in x,y and z directions

```

```

#define grainIDFile "grainID.txt" // Grain ID File
#define headerLinesGrainIDFile 5 // No. of header Lines
#define grainOrientationsFile "orientations.txt" // Slip Normals File

```

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Table 1: FCC Copper Slip Systems

System Number	Slip Direction	Slip Plane
1	$[1\ 1\ 0]$	$(1\ 1\ 1)$
2	$[-1\ 0\ 1]$	$(1\ 1\ 1)$
3	$[0\ 1\ -1]$	$(1\ 1\ 1)$
4	$[1\ 0\ 1]$	$(-1\ 1\ 1)$
5	$[-1\ -1\ 0]$	$(-1\ 1\ 1)$
6	$[0\ 1\ -1]$	$(-1\ 1\ 1)$
7	$[-1\ 0\ 1]$	$(1\ -1\ 1)$
8	$[0\ -1\ -1]$	$(1\ -1\ 1)$
9	$[1\ 1\ 0]$	$(1\ -1\ 1)$
10	$[-1\ 1\ 0]$	$(-1\ -1\ 1)$
11	$[1\ 0\ 1]$	$(-1\ -1\ 1)$
12	$[0\ -1\ -1]$	$(-1\ -1\ 1)$

## References

- [1] Anand, L., and M. Kothari. "A computational procedure for rate-independent crystal plasticity." Journal of the Mechanics and Physics of Solids 44.4 (1996): 525-558.
- [2] Neper homepage [Neper](#)