

Langmuir Project Design

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The Langmuir project is designed for ...

The Langmuir project file structure can be seen as below,

Langmuir Project File Structure

packages/Lib/

- | - Poisson1D_Solver.py # 1D Poisson solver
- | - Poisson2D_Solver.py # 2D Poisson solver

packages/Mesh/

- | - Geom2D.py # 2D geometry generator
- | - Mesh2D.py # 2D mesh generator

packages/Model/Common

- | - Particle.cpp # class for single particle, 3D by default
- | - Multi_Particle.py # class for multi-particles, supporting vector operations
- | - Particle_Mover.py # functions to move particle
- | - Field.py # class for field info and solver

packages/Model/Feature2D

- | - Feature2D_main.py # 'main' func for feature model
- | - Feature2D_mesh.py # class for feature mesh, inherited from mesh2d.py
- | - Feature2D_rct.py # class for reaction
- | - Feature2D_rflct.py # class for reflection

packages/Model/Sheath2D

- | - Sheath2D_main.py # 'main' func for sheath model
- | - Sheath2D_coll.py # class for collision

packages/Model/Reactor2D

- | - Reactor2D_main.py # 'main' func for reactor model
- | - Reactor2D_transp.py # solver for plasma transport
- | - Reactor2D_eerg.py # solver for electron energy equation
- | - Reactor2D_edistb.py # solver for electron energy distribution
- | - Reactor2D_field.py # solver for field
- | - Reactor2D_rct.py # class for reaction

```

database/
| - Species.csv # database for species info
| - Cross_Section.csv # database for species info

database/valid
| - reaction_xxx.yaml # validated reaction set for xxx

utest/
| - Particle_test.py # test for Particle.py
| - Multi_Particle_test.py # test for Multi_Particle.py
| - Particle_Mover_test.py # test for Particle_Mover.py

stest/
| - xxx_test.py # test for xxx

run/Feature2D
| - Feature2D_run.py # prepare input and call main()
| - mesh.npz # mesh for feature model
| - feature.yaml # input for feature model
| - reaction.csv # reaction set

run/Sheath2D
| - Sheath2D_run.py # prepare input and call main()
| - sheath.yaml # input for sheath model
| - collision.py # collision set

doc/
| - Langmuir_Doc.lyx # lyx for Langmuir pyhsics
| - Langmuir_Doc.pdf # pdf for Langmuir pyhsics
| - Langmuir_Doc.tex # LaTeX for Langmuir pyhsics
| - Langmuir_Structure.lyx # lyx for Langmuir code structure
| - Langmuir_Structure.pdf # pdf for Langmuir code structure
| - Langmuir_Structure.tex # LaTeX for Langmuir code structure

```

This project design will focus on the modules.

The project design is based on the modeling of the physics and cannot go beyond it. In order to well design the project structure, the models need to be analyzed in details.

Let's see the particle-based Monte Carlo models first.

The most frequently called function in a Monte Carlo model is random number generator. In order to repeat the simulation identically, the random number generator should be able to produce identical pseudo-random numbers for each run. The random number generator will be placed in /basics.

```

Function Poisson2d_Solver(input)
    return solution

```

```

# input:
# solution:

```

Another core module for particle-based models is, as its name states, particle module, which determines the properties and actions of a particle.

```

Class Particle()
-- attribute --
    self.name: str
    self.type: str, one of('Eon', 'Ion', 'Neut')
    self.charge: float, unit in C
    self.mass: float, unit in AMU
    self.position: float array(3), unit in m
    self.speed: float, unit in m/s
    or self.erg: float, unit in eV
    or self.vel: float array(3) unit in m/s
    self.uvec: float array(3), unitless, normalized velocity
    or self.ang: float array(2), theta and phi
    self.accl: float array(3), unit in m/s**2, acceleration
    self.isAlive: bool, state of particle
-- method --
    self.initParticle()
    self.initPosition()
    self.initSpeed(imode)
    self.initEnergy(imode)
        # imode: str, determines the distribution such as
        # Uniform, Normal and Cosine.
    self.move(dL, imode='Space')
        # dL: float, unit in m, spacestep for a move
        # imode: str
        # move in a spacestep, with no field,
        # that is self.accl = (0.0, 0.0, 0.0), used in Feature Model.
    or self.move(dt, EF, imode='Normal')
        # dt: float, unit in s, timestep for a move
        # EF: float array(3), unit in V/m, E-Field
        # move in timestep, with E-field, 1st order accuracy.
        # used in EEDF Model, or Feature Model with E-field on.
    or self.move(dt, EF, imode='leapfrog')
        # dt: float, unit in s, timestep for a move
        # EF: float array(3), unit in V/m, E-Field
        # move in timestep, with time-varying E-field or B-field,
        # 2nd order accuracy, able to track the oscillating motion.
    self.checkBndy(domain, imode):
        # domain: float array(6)
        # imode: str, 'Rflective' or 'Periodic'

```

self.move() is the key method for a particle. Depending on the field, best algorithm needs to be chosen for required accuracy and speed.

self.move() can be designed as a interface, where the real move function is called from /basics or /common. Feature Model needs intensive information for mesh, while Sheath Model and EEDF Model do not.

Feature Model mesh class is designed as

```

Class Mesh()
-- attribute --
    self.name: str
    self.domain: float array(6), unit in m, (top, bottom, left, right, front, back)
    self.res: float array(3), unit in m, resolution in structured mesh
    self.ngrid: int array(3), num of nodes in structured mesh
    self.x, self.z, self.y: float array(m,n), unit in m, coordinates in axis x, z,
    self.mat: int array(m,n), material number, vacuum = 0 by default
    self.matDict: dictionary to map material number to material name
    self.surf: int array(m,n), indicator for surface nodes
        # 0: non-surf node
        # 1: surf node in material
        # -1: surf node in vacuum
-- method --
    self.initMesh()
    self.readMesh()
    self.addShape(ishape='Rectangle', 'Triangle' or 'Circle')
    or self.addRectangle(), self.addTriangle() and self.addCircle()
    self.find_surf(): assign values to self.surf
    self.saveMesh()

```

This mesh class can be shared by all mesh-based models, such as Feature Model and Fluid Model.

This mesh class will support read-in function so that it enables restart capability for Sheath Model.

In Feature Model, mesh evolution is the key so that mesh needs more methods. Let's create a FeatMesh class to inheritate generic Mesh class.

```

Class FeatMesh(Mesh): FeatMesh inherits all attributes and methods from Mesh
-- attribute --
    self.surfNorm: list of float array(3) store all surface normal along the surface
-- method --
    self.checkHit(Particle.position)
        # the position of the particle is passed to checkHit()
        # position in continuous space is mapped to meshgrid
        # position --> index
        # check the material of the index
        # if self.mat[index] != 0, it is a hit
    self.calcSurfNorm(global or local)
        # calculate the surface normal either globally or for a given node
    self.updateMat(index, newMat)

```

```

        # change the mat at index to newMat
        # to protect the self.mat, we define a method to explicitly
        # change the material
self.findFloatChell()
        # find the floating cells which are detached from surface
        # the algorithm is not determined yet
self.dropFloatCell(imode='Remove' or 'Drop')
        # process floating cells
        # imode='Remove': Remove floating cells, change mat to zero
        # imode='Drop': Drop floating cells to bottom, like deposition

```

checkHit() is placed under FeatMesh() class instead of Particle() because checkHit() is more associated to Mesh than Particle.

self.calcSurfNorm() and self.findFloatCell() can be designed as a interface, where the real functions are called from /basics or /common

There are two classes specific to Feature Model, Reflection and Reaction.

When a hit occurs, the program first check the reaction. If a reaction occurs, the mesh material will be updated; otherwise, the particle is going to reflect.

Reaction class can be seen as blow.

```

Class Reaction()
-- attribute --
    self.index: int array(3), index of hit node
-- method --
    self.readReaction(fname)
        # fname: str, filename for chemical reaction
        # Read in reaction files and
        # create a reaction list with associated probability
self.getParticle(Particle)
        # get information from Particle
        # name, erg, uvec
self.getMesh(FeatMesh)
        # get information from FeatMesh
        # mat, matDict, surfNorm
self.determineReaction(erg, angle, surfNorm)
        # explanied with an example
        # Cl+ + Si_ --> Etch : p1(erg, angle, surfNorm)
        # Cl+ + Si_ --> SiCl_ : p2(erg, angle, surfNorm)
        # Cl+ + Si_ --> SiCl : p3(erg, angle, surfNorm)
        # Cl+ + Si_ --> Reflect : p4(erg, angle, surfNorm)
        # normalize pi as pi/(p1 + p2 + p3 + p4)
        # roll a dice and determine which reaction to happen
        # return the serial number of the reaction
self.makeReaction(index)
        # index: int, the serial number of the reaction in reaction list

```

```

        # change Particle.isAlive to False
        # call updateMat()
        # if the reaction is a type of "Byproduct": initParticle()

```

Reflection class can be seen as below.

```

Class Reflection()
-- attribute --
    self.index: int array(3), index of hit node
-- method --
    self.getParticle(Particle)
        # get information from Particle
        # name, erg, uvec
    self.getMesh(FeatMesh)
        # get information from FeatMesh
        # mat, matDict, surfNorm          self.reflect()
    self.updateAngle(diffThld, specThld)
        # diffThld: threshold of diffusive reflection
        # specThld: threshold of specular reflection
        # return the anlge or uvec after reflection and
        #         the energy after reflection
        # if erg < diffThld: diffusive reflection
        # if diffThld < erg < specThld: mixed reflection
        #         anlge is determined by the combination of
        #         diffusive and specular reflections
        #         angleMixed =
        #         [(erg - diffThld)*angleSpec + (erg - specThld)*angleDiff]
        #         /(specThld - diffThld)
        # if erg > specThld: specular reflection
    self.updateEnergy(thermThld, thermErg, facErgLoss)
        # thermThld: threshold of falling to thermErg
        # thermErg: thermal energy
        #         thermErg must be less than thermThld
        # facErgLoss: factor of energy loss due to each collision
        # if erg < thermThld: Particle.erg = thermErg
        # if erg > thermThld: Particle.erg = facErgLoss*Particle.erg
    self.postMove(numStep):
        # numStep: int, the number of steps to move after reflection
        # if the reflection angle is close to the surface (angle is
        # almost 90 degrees w.r.t. surface normal), the particle
        # travel a long distance within a material before going out.
        # the postMove will move the particle by steps without
        # checkHit() until it gets out of the material

```

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