Documentation

ContaminationFlow on Linux and Windows

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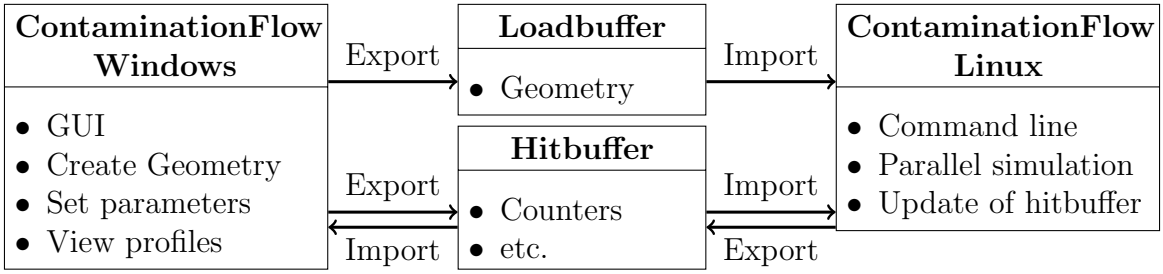
Contents

1. General Structure



General structure for ContaminationFlow simulation

- Code adapted from Molflow
- ContaminationFlow Windows primary used to create Geometry and to view simulation results through the GUI
- ContaminationFlow Linux primary used for simulation and calculation of counters, profiles, etc.
- Loadbuffer contains information of geometry
- Hitbuffer contains information such as hit counters, profiles, etc.
- Import and export of buffer files for communication between ContaminationFlow Windows and ContaminationFlow Linux
- Export of simulationHistory for ContaminationFlow Linux

2. ContaminationFlow Linux

- Parallel simulation on several sub processes
- Processing and control of data in main process
- Update and accumulation of hit counters and other information such as profiles
- SimulationHistory, final hitbuffer and used parameters exported to results folder

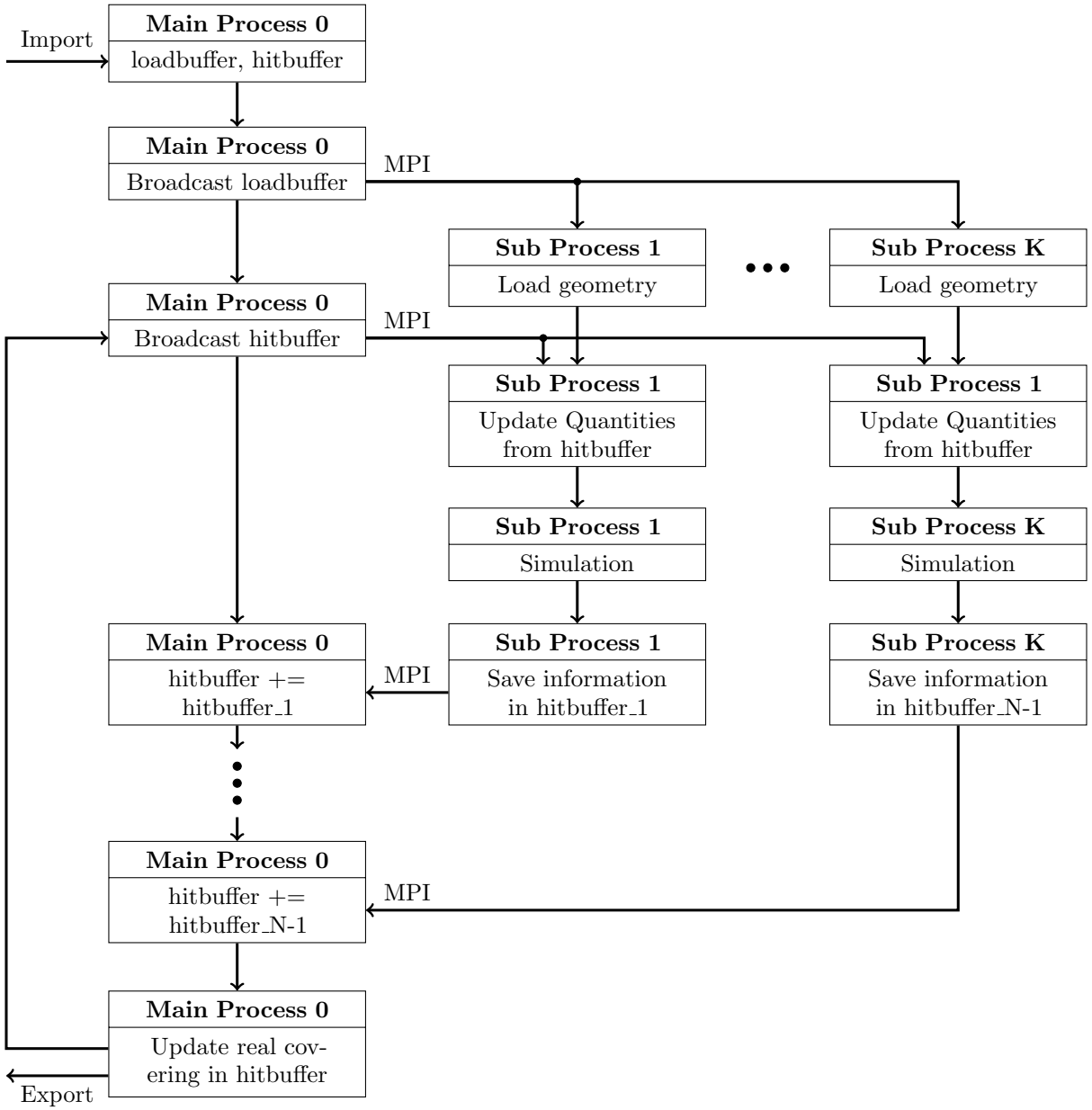


Figure 2.1.: Processing of data in main and sub processes

2.1. Call of Application from Command line

New class ProblemDef

- Defines parameters used for simulation
- Possible adaptation of default paramaters through input file
- Creates result folder for simulation if desired
  - Final resultbuffer
  - Final covering, input file and console output as text files

Application with custom parameters using input file

Call of ContaminationFlow Linux application in the command line:

```
$ module load mpi
$ mpirun -n N MolflowLinux inputfile save
```

with the following command line parameters:

- `N` : desired number of worker processes; simulation on K=N-1 worker processes
- `MolflowLinux` : path to application, e.g. `~/MolflowLinux/Debug/MolflowLinux`
- `inputfile` : path to file that defines simulation parameters
- `save` : determines whether result directory is created (1: true, 0:false); default: `1`

and the input file defining the following parameters:

- $loadbufferPath$: path to loadbuffer file, contains geometry, e.g. `~/loadbuffer`
- $hitbufferPath$: path to hitbuffer file, contains counters, etc., e.g. `~/hitbuffer`
- $simulationTime$: simulation time per iteration step; default: `10.0`
- $unit$: simulation time unit; default: `s`
- $maxTime$: maximum simulation time; default: `10.0`
- $maxUnit$: maximum simulation time unit; default: `y`
- $iterationNumber$: number of iterations; default: 43200
- $particleDia$: diameter of particles; default: `2·76E-12`
- E_{de} : binding energy of a particle on pure substrate; default: `1E-21`
- H_{vap} : vaporization enthalpy of a particle in case of multilayer contamination; default: `0.8E-19`
- W_{tr} : transition width between monolayer and multilayer properties; default: `1`
- $sticking$: constant sticking coefficient for all facets, set to zero, not used at the moment; default: `0`
- $targetPaticles$: minimum number of desorbed particles per iter.; default: `1000`
- $targetError$: average statistical uncertainty (error) to be achieved for each iteration, calculated as the average (weighted with the facets area) of the normalized standard deviation of events per facet; default: `0.001`
- $hitRatioLimit$: Ratio at which hits are ignored, default: `1E-5`
- $Tmin$ or t_min : minimum time for step size; default: `1E-4`
- $maxStepSize$ or t_max : maximum time for step size; default: `max`
- $maxSimPerIt$: maximum simulation steps per iteration; default: `max`
- $coveringMinThresh$: minimum covering (through multiplication); default: `10000`

- *histsize* : Size of history lists; default: `max`
- *vipFacets* : very important facets: facets with have their own target error. input in inputfile as alternating sequence of facet numbers and respective target errors seperated via blanks; default: `[]`

Terminology

- Simulation time: desired computation time until check if target is reached for iteration
- Simulated time: physical time in the simulated system, e.g. flight time or residence time of a particle
- Maximum simulation time: desired total simulated time
- Step size: desired simulated time per particle for iteration

2.2. Communication

Import and export of buffer files

- New Databuff struct that replaces Dataport struct from MolFlow Windows

```
typedef unsigned char BYTE;
typedef struct {
  signed int size;
  BYTE *buff;
} Databuff;
```

- New functions `importBuff(·)` and `exportBuff(·)` for import of buffer files and export of Databuff struct
- New functions `checkReadable(·)` and `checkWriteable(·)` to check if file is readable or writeable

Communication between worker processes via MPI

- Main process 0 sends Databuff struct containing loadbuffer and Databuff struct containing hitbuffer and required simulationHistory values to sub processes using `MPI_Bcast(·)`
- Sub processes send updated Databuff struct containing hitbuffer and required simulationHistory values to main process 0 using `MPI_Send(·)` and `MPI_Recv(·)`

2.3. Usage of *boost* Library

Multiprecision

- Increase precision for variables if required
- Avoid overflow for integer and underflow for floating point numbers

2.4. New Quantities

New counter `covering`

- Number of carbon equivalent particles on facet
- Increases with adsorption, decreases with desorption
- Extracted from new hitbuffer counter from `Simulationcalc.cpp` file in `getCovering(·)`

Coverage

- Number of monolayers of adsorbed particles
- Calculated from covering, particle diamieter (previously gas mass) and facet area
- Coverage computed from `Simulationcalc.cpp` file in `calcCoverage(·)`

Sticking factor

- Ratio adsorbed particles to impinging particles
- Set to 0, can be adapted for all facets through input file

Binding energy

- Calculated from E_{de} , H_{vap} and W_{tr}
- Energy computed from `Simulationcalc.cpp` file in `calcEnergy(·)`

Desorption [1/s]

- number of particles desorbing per second
- Calculated from binding energy, covering and temperature
- Desorption computed from `Simulationcalc.cpp` file in `calcDesorption(·)`

Desorption Rate [Pa m^3/s]

- Calculated from desorption, gas mass and facet area before each iteration
- Used to determine starting point for new particle
- Desorption rate computed from `Simulationcalc.cpp` file in `calcDesorptionRate(·)`

Outgassing Rate

- Calculated from facet outgassing and temerature defined in sHandle
- Outgassing rate computed from `Worker.cpp` file in `CalcTotalOutgassingWorker()`

$K_{real/virtual}$

- Number of real particles represented by test particles
- Calculated from desorption & outgassing rate and number of desorbed molecules
- $K_{real/virtual}$ computed from `Simulationcalc.cpp` file in `GetMoleculesPerTP(·)`

Statistical Error

- Calculated from hits and desorbed particles (of facet and total)
- Used to determine significance of simulation results of iteration

2.5. Iterative Algorithm

2.5.1. Initialization of simulation

New class to store Simulation History

- SimulationHistory class

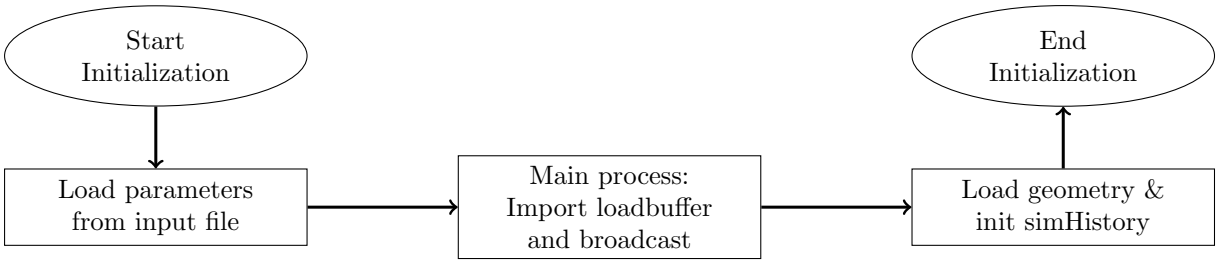


Figure 2.2.: Overview: Initialize simulation

```
class SimulationHistory {
public:
 SimulationHistory();
 SimulationHistory(Databuff *hitbuffer);

 HistoryList<llong> coveringList;
 HistoryList<llong> desorbedList;
 HistoryList<double> hitList;
 HistoryList<double> errorList;

 double lastTime;
 int currentStep;
};
```

```
template <typename T> class HistoryList {
public:
 HistoryList();
 std::vector<std::pair<double,std::vector<T>>> pointInTimeList;
 std::vector<T> currentList;
};
```

- In `SimulationLinux.h` and `SimulationLinux.cpp` file
- SimulationHistory updated after each iteration in `UpdateCovering(·)` from `UpdateMainProcess.` file
- Currently recorded quantities: covering and error for each facet for each iteration, total hits and desorbed particles for each facet
- lastTime: simulated time (accumulated time steps) instead of computation time

2.5.2. Simulation on subprocesses

Calculate Step Size

- Use `simHistory→currentStep` to calculate logarithmic step size
- Calculation in `UpdateMainProcess.cpp` file in `getStepSize()`

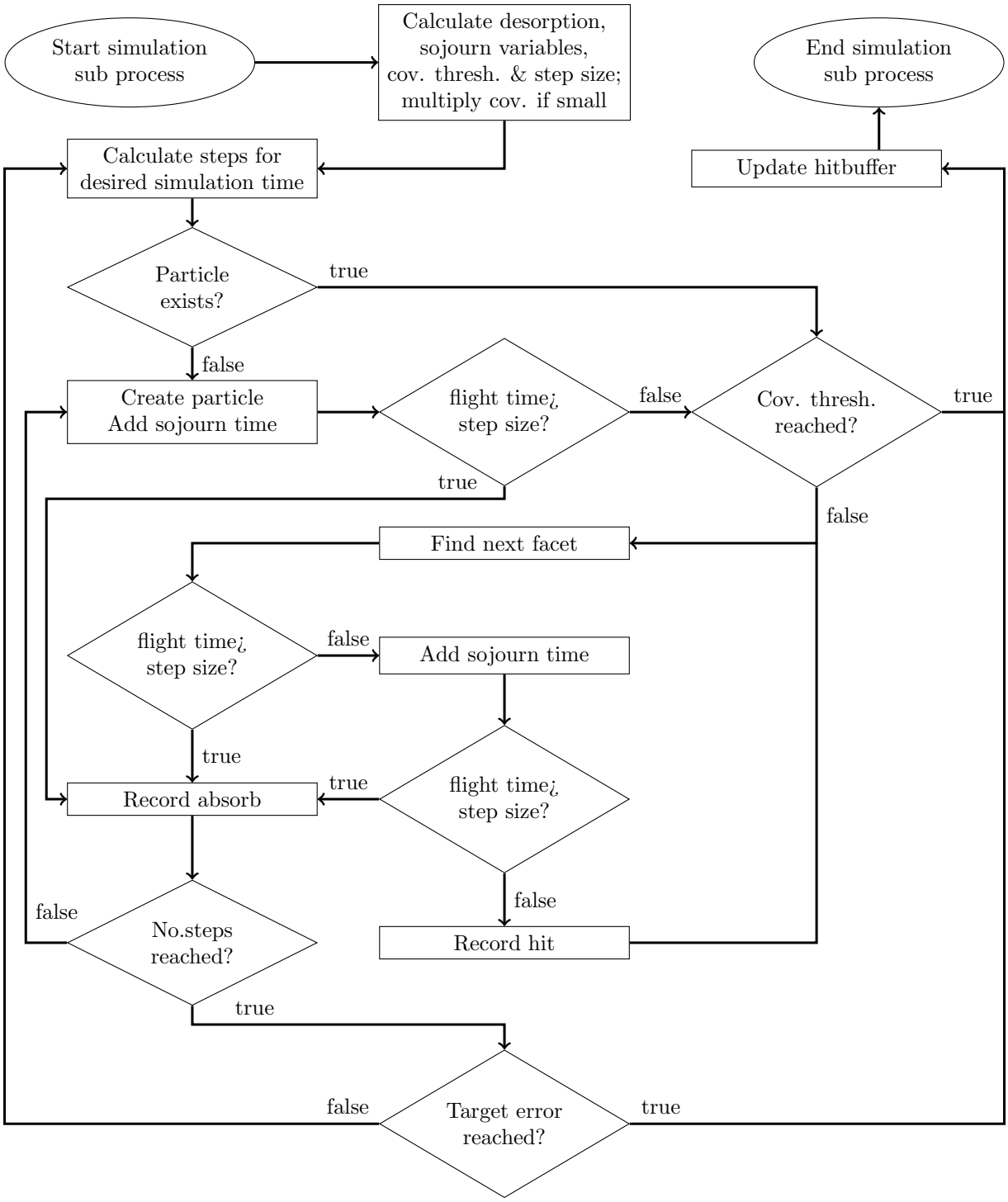


Figure 2.3.: Overview: simulation on sub processes

Calculate Covering Threshold

- Set lower threshold for covering for each facet to prevent covering getting negative
- Stop simulation once threshold is reached
- Threshold set in `setCoveringThreshold(·)` from `Iteration.cpp` file

Multiply small covering

- Multiply covering so that smallest covering $\geq$ `ProblemDef::coveringThreshMin`
- Multiply covering threshold with same factor
- Adapt covering threshold and `simHistory→coveringList`
- Calculation in `checkSmallCovering(·)` from `SimulationLinux.cpp` file

Calculate desorption

- Desorption rate calculated from current covering values
- Calculation in `UpdateDesorptionRate(·)` from `UpdateSubProcess.cpp` file

Calculate residence time and binding energy

- residence time of a particle on a surface calculated from binding energy, coverage and thermal oscillation frequency
- binding energy calculated using binding energy on pure substrate, vaporization enthalpy and coverage
- thermal oscillation frequency calculated using temperature, Boltzmann constant and Planck's constant
- Calculation in `UpdateSojourn(·)` from `UpdateSubProcess.cpp` file

Target error reached?

- Calculate statistical error in `UpdateError()` from `UpdateSubProcess.cpp` file
- Check if vip facets reached their own target error
- Check if normal facets total error reached target error
- Total error calculated from summing facet error weighted with facet area
- Set error of facets that reached `ProblemDef::hitRatioLimit` to inf
- Set error of facets with no hits and desorption to inf
- Facets with error=inf are not considered
  - if vip facet: target automatically reached
  - if normal facet: facet error and area not used for calculation
- Check in `checkErrorSub(·)` from `UpdateSubProcess.cpp` file

2.6. Update main buffer

Multiply small covering before summation of subprocesses

- Only if covering was multiplied in subprocesses
- Multiply covering in `hitbuffer` and `simHistory` of main process before simulation using same factor

Divide small covering after summation of subprocesses

- Only if covering was multiplied in subprocesses
- Divide covering in `hitbuffer` and `simHistory` of main process before simulation using same factor
- Adapt hitbuffer and `simHistory→coveringList`

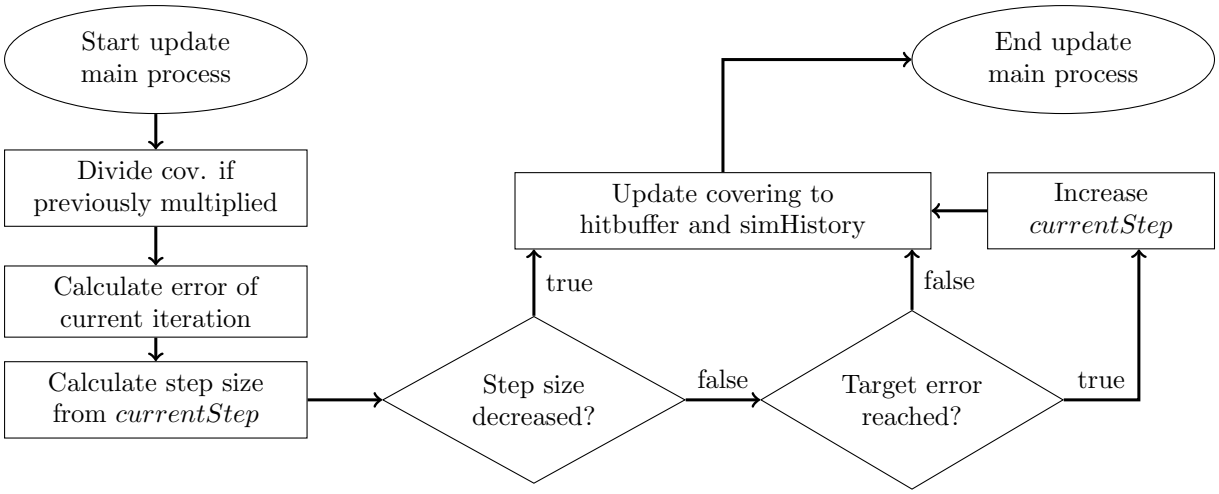


Figure 2.4.: Overview: update of covering in hitbuffer

- Calculation in `UndoSmallCovering(·)` from `SimulationLinux.cpp`

Error Calculation

- Calculate statistical error of normal facets
- Facet error calculated from hits and desorbed particles, weighted with opacity
- Total error calculated from summing facet error weighted with facet area
- Set error of facets that reached `ProblemDef::hitRatioLimit` to inf
- Set error of facets with no hits and desorption to inf
- Facets with error=inf are not used for calculation of total error
  - if vip facet: target automatically reached
  - if normal facet: facet error and area not used for calculation
- Save error in `simHistory→errorList`
- Management in `UpdateErrorMain(·)` from `UpdateMainProcess.cpp` file

Management of Step Size

- Adapt step size if desorption × time step was larger than covering
- Increase `simHistory→currentStep` if no adaptation & target errors reached
- Management in `manageStepSize(·)` from `UpdateMainProcess.cpp` file

Update Covering

- Use step size and $K_{real/virt}$ to calculate new covering
- Calculation in `UpdateCovering(·)` from `UpdateMainProcess.cpp` file

2.7. Summary

General Pipeline

- Initialize MPI, `ProblemDef p` and `SimulationHistory simHistory`
- Load geometry into `Simulation sHandle` using `LoadSimulation()`

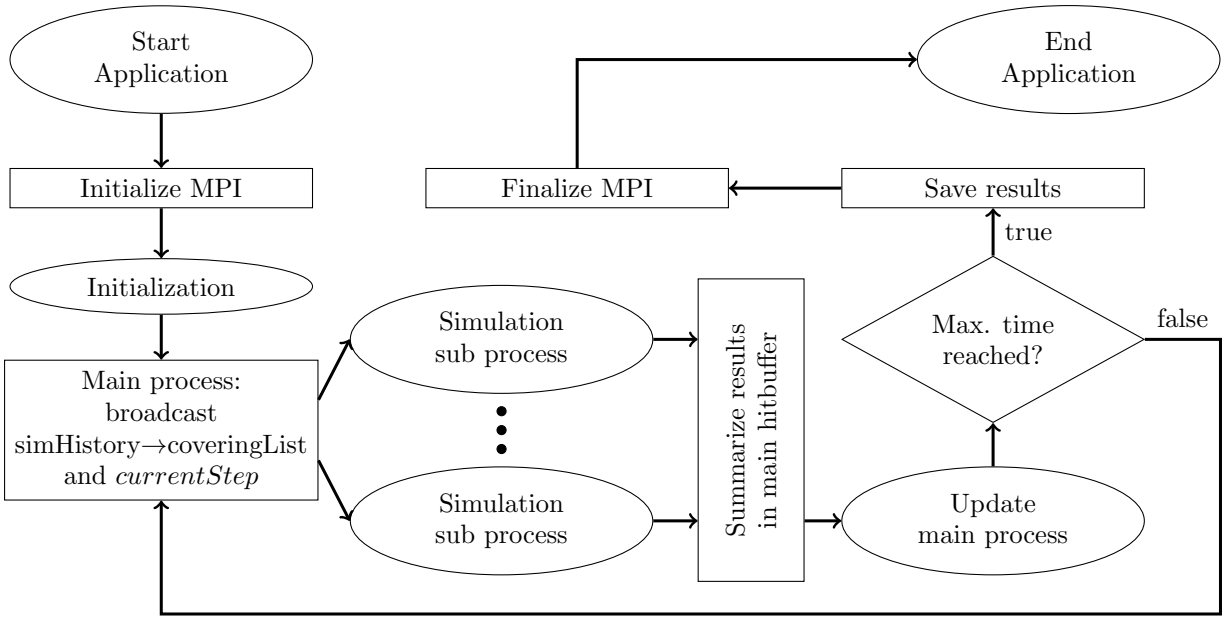


Figure 2.5.: Overview: ContaminationFlow application

- Iteration until desired maximum simulation time is reached:
  - Reset hitbuffer counters using `initbufftotero(·)`
  - Broadcast `simHistory→coveringList` using `MPI_Bcast(·)`
  - Simulation in sub processes
    - Simulate until `targetParticles` and `targetError` or `covthresh` reached
    - Update hitbuffers of sub processes from `sHandle` using `UpdateSubHits(·)` from `UpdateSubProcess.cpp`
  - Update Main process:
    - Send hitbuffer to main process using `MPI_Send(·)` and `MPI_Recv(·)`
    - Update of hitbuffer in `UpdateMainHits(·)` from `UpdateMainProcess.cpp`
  - Update error of iteration using `UpdateErrorMain(·)` from `UpdateMainProcess.cpp`
  - Calculate real covering in main process using $K_{real/virtual}$ and simulated step size in `UpdateCovering(·)` from `UpdateMainProcess.cpp`, save in `simHistory`
  - Update real covering in hitbuffer of main process in `UpdateCoveringphys(·)` from `UpdateMainProcess.cpp`
- Export final results (hitbuffer and simulationHistory) to results folder
- Close MPI

3. ContaminationFlow Windows

- Create Geometry and set parameters such as initial coverage and temperature

3.1. Graphical User Interface

Add screenshot of GUI

New GUI elements

- "Particles out" renamed to Contamination level
  - Text field for covering
  - Text field for coverage
- New facet properties
  - Effective surface factor
  - Facet depth and facet volume
  - Diffusion coefficient
  - Concentration and gas mass
- Window for CoveringHistory (reworked to SimulationHistory in ContaminationFlow Linux)
- PressureEvolution window expanded
  - Added list that contains information of graph
  - Option to show only selected facets or all
  - List exportable

3.2. Communication

Import and export of buffer files via GUI

- New Databuff struct

```
typedef unsigned char BYTE;
typedef struct
  signed int size;
  BYTE *buff;
Databuff;
```

- New functions `importBuff(·)` and `exportBuff(·)` for import and export of buffer files/Databuff struct
- New options in file menu: `Export buffer` and `Import buffer`

3.3. New Quantities

New counter `covering`

- Covering computed in `SimulationMC.cpp` file in `updatecovering(·)`

- Added covering counter to hitbuffer
- Added covering to GUI, can be defined through textfield

New facet property `effetiveSurfaceFactor`

- Defines increase of facet area due to texture

New facet property `facetDepth`

- Defines depth of facet

New facet property `diffusionCoefficient`

- Defines diffusion coefficient

New facet property `concentration`

- Defines concentration = mass of particles in volume

Removal of irrelevant quantities

- Sticking factor and pumping speed removed from GUI
- `calcSticking()` and `calcFlow()` in `Molflow.cpp` file not used anymore
- Flow not needed for iterative Algorithm

3.4. Iterative algorithm

New class to store covering for all facets at any time

- In `HistoryWin.cpp` and `HistoryWin.h` file
- `std::vector<std::pair<double,std::vector<double>>> pointintime_list` to store points in time and respective covering for all facets
- New GUI option to add and remove entries for `pointintime_list`
- New GUI option to export or import a complete list

A. Formulas for new Quantities

Constants

$$\begin{aligned} carbondiameter &= 2 \cdot 76\text{E} - 12 \\ K_b &= 1.38\text{E} - 23 \\ h &= 6.626\text{E} - 34 \end{aligned} \tag{A.1}$$

Number of carbon equivalent particles of one monolayer

$$N_{mono} = \frac{\text{Area of Facet } [\text{m}^2]}{carbondiameter^2} \tag{A.2}$$

Carbon equivalent relative mass factor

$$\Delta N_{surf} = \frac{\text{carbon equivalent gas mass}}{12.011} \tag{A.3}$$

Covering θ^*

$$\theta^* = N_{\text{particles on facet}} \tag{A.4}$$

Coverage θ

$$\theta = \frac{\theta^*}{N_{mono}/\Delta N_{surf}} \tag{A.5}$$

Step function $step(x, y_{start}, y_{end}, x_{turningpoint}, w)$

$$E = \frac{y_{start} - y_{end}}{2} \cdot \tanh\Big((x_{turningpoint} - x) \cdot \frac{5.4}{w}\Big) + \frac{y_{start} + y_{end}}{2} \tag{A.6}$$

Binding Energy E

$$\begin{aligned} E &= step(\theta, E_{de}, H_{vap}, 1, W_{tr}) \\ &= \frac{E_{de} - H_{vap}}{2} \cdot \tanh\Big((1 - \theta) \cdot \frac{5.4}{W_{tr}}\Big) + \frac{E_{de} + H_{vap}}{2} \end{aligned} \tag{A.7}$$

Residence

$$\begin{aligned} Frequency &= \frac{K_b T}{h} \\ Energy &= E \end{aligned} \tag{A.8}$$

Residence Time

$$\begin{aligned} A &= \exp\Big(- Energy/(k_b T)\Big) \\ residence\ time &= \frac{-\ln(rnd)}{A \cdot Frequency} \end{aligned} \tag{A.9}$$

Desorption rate des

$$\begin{aligned} \tau &= \frac{h}{K_b T} \\ d &= step(\theta, 1, 0, 1, W_{tr}) \end{aligned} \tag{A.10}$$

$$des = \begin{cases} \frac{1}{\tau}\ \theta^d\ \exp\left(-\frac{E}{K_b T}\right) \cdot \frac{N_{mono}}{\Delta N_{surf}} \cdot K_b T, & \text{if } \theta^* > 0 \\ 0, & \text{otherwise} \end{cases}$$

Outgassing rate out

$$out = \frac{\text{Facet outgassing}}{K_b T} \tag{A.11}$$

$K_{real/virtual}$

$$\mathrm{K}_{\mathrm{real/virtual}} = \frac{\sum\limits_{\text{facets}} \left(out + \frac{des}{K_b T}\right)}{\text{number of total desorbed molecules}} \tag{A.12}$$

Step Size T_{step}

$$\begin{aligned} T_{min} &= \mathrm{Tmin} \\ T_i &= T_{min} \cdot \exp\left(\mathrm{i} \cdot \ln(\text{max. simulation time}/T_{min})/\text{max. \# of steps}\right) \\ T_{step} &= T_{currentStep+1} - T_{currentStep} \end{aligned} \tag{A.13}$$

Error

$$error = \begin{cases} inf \quad \text{, if (hits+desorbed) on facet =0} \\ \left(\frac{1}{\text{(hits + desorbed) on facet}} \cdot \frac{1 - \text{(hits + desorbed) on facet}}{\text{total (hits + desorbed)}}\right)^{0.5} \quad \text{, else} \end{cases} \tag{A.14}$$

B. Datatypes

B.1. Class Members

Name	Datatype	Alias
SimulationHistory::coveringList	boost::multiprecision::uint_128t	covBoost
FacetHitBuffer::covering	llong	covLlong
FacetProperties::desorption	boost::multiprecision::float128	desBoost
Simulation::coveringThreshold	llong	

B.2. Functions

Function	Output Datatype	Relevant Input
getCovering()	boost::multiprecision::float128	covBoost
getCovering()	llong	covLlong
calcCoverage()	boost::multiprecision::float128 or llong	getCovering()
calcDesorption()	boost::multiprecision::float128	calcCoverage()
calcDesorptionRate()	boost::multiprecision::float128	calcDesorption()
calctotalDesorption()	boost::multiprecision::float128	desBoost
GetMoleculesPerTP()	boost::multiprecision::float128	desBoost

C. Overview of new Classes and Functions

C.1. New Classes

SimulationHistory	
coveringList	of class HistoryList, stores covering history
errorList	of class HistoryList, stores error history
hitList	of class HistoryList, stores hits for each facet
desorbedList	of class HistoryList, stores desorbed particles for each facet
startNewParticle	Determines wether to create a new particle for next iteration
numFacet	number of Facets
numSubProcess	number of sub processes used for simulation
nbDesorbed_old	number of total desorbed molecules of previous iteration ⇒ To calculate difference between consecutive iterations
flightTime	Simulated flight time for iteration
nParticles	Simulated particles for iteration
lastTime	Total simulated time = last time in Lists
currentStep	step of logarithmic time step calculation in `getStepSize()`
stepSize	current step size
updateHistory()	Reset and update from hitbuffer
appendList()	Updates coveringList from hitbuffer
print()	Print to terminal
write()	Write to file

HistoryList	
pointInTimeList	list containing history respective facet values
currentList	list containing facet values at current step
currIt	current iteration number
appendCurrent()	Appends currentList to pointInTimeList
appendList()	Append input list to pointInTimeList
convertTime()	Converts time for better clarity
printCurrent()	Print currentList as table to terminal, optional message
print()	Print pointInTimeList as table to terminal, optional msg
write(), read()	Write to file, read from file
set/getCurrent()	Set/get value of desired facet in currentList
setLast(), getLast()	Set/get value of desired facet from pointInTimeList

ProblemDef	
resultpath	Path of result folder
outFile	Path of file that contains terminal output
loadbufferPath	Path of loadbuffer file
hitbufferPath	Path of hitbuffer file
simulationTime, unit ⇒simulationTimeMS	Computation time of each iteration in milliseconds
maxTime, maxUnit ⇒maxTimeS	Maximal total simulated time in seconds
iterationNumber	Number of iterations
particleDia	Diameter of particles
E_de, H_vap, W_tr	Parameters to calculate binding energy, see equation A.7
sticking	Sticking factor for all facets
targetParticles/-Error	Target values for each iteration
hitRatioLimit	threshold of hitratio at which hits are ignored
coveringMinThresh	Minimum covering, multiplication to this if covering low
Tmin, maxStepSize	Minimum/ Maximum step size
maxSimPerIt	Maximun simulation steps per iteration
histSize	Size of history lists (most recent values in memory)
vipFacets	alterning: vip facet and target error, e.g. 1 0.001 3 0.002
readInputfile()	Initialization from input file
printInputfile()	Print to terminal

C.2. New Functions

C.2.1. molflowlinux_main.cpp

Preprocessing	
parametercheck()	Checks validity of input parameters from input file Defines values for ProblemDef object `p`
importBuff()	Import load- and hitbuffer to main process
MPI_Bcast()	Send loadbuffer to sub processes
LoadSimulation()	Load geometry from loadbuffer
initCoveringThresh()	Initialize covering threshold
`simHistory`	Initialize SimulationHistory object

Simulation Loop	
initbufftozero()	Reset all hitbuffer counters except covering
MPI_Bcast()	Send `simHistory→coveringList` and `simHistory→currentStep` to sub processes
setCoveringThreshold()	Sets covering threshold for each facet
UpdateSojourn()	Sets sojourn variables for each facet
UpdateDesorptionRate()	Sets desorption for each facet, ends simulation if 0
simulateSub()	Simulation on sub processes
MPI_Send(), MPI_Recv()	Send sub hitbuffer to main process
UpdateMCMainHits()	Add simulation results from sub hitbuffer to main hitbuffer
UndoSmallCovering()	Divide covering if it was multiplied before
UpdateErrorMain()	Calculate and save error of iteration to simHistory
UpdateCovering()	Calculate and save new covering to simHistory
UpdateCoveringphys()	Saves current covering to hitbuffer
`simHistory→coveringList` `simHistory→errorList`	Adapt size to `p→histSize` if necessary
End simulation if maximum simulation time is reached	

Postprocessing	
exportBuff()	Export final hitbuffer
simHistory→write()	Export simulation history

C.2.2. SimulationLinux.cpp

simulateSub()	
targetParticles, targetError	Calculate target values from overall target and number sub processes
`simHistory->updateHistory()`	Reset and Update SimulationHistory object from hitbuffer
smallCovering, smallCoveringFactor	If smallCovering: Covering is multiplied by smallCoveringFactor to improve statistics
SimulationRun()	Simulate for desired simulation time
UpdateError()	Calculate current error of sub process
CheckErrorSub()	Checks if normal facets reached targetError and if vip facets reached own target
UpdateMCSubHits()	Save simulation results to hitbuffer

Small covering	
CheckSmallCovering()	If smallCovering, find smallCoveringFactor to reach `p→coveringMinThresh`
UndoSmallCovering()	If smallCovering, divide by previously determined smallCoveringFactor

C.2.3. Iteration.cpp

Set Covering Threshold to avoid negative covering	
initCoveringThresh()	Initializes size of covering threshold vector
setCoveringThreshold()	Sets covering threshold for each facet

C.2.4. Buffer.cpp

Buffer functions	
Databuff struct()	signed int size BYTE *buff
checkReadable()	Checks if file can be opened for reading
checkWriteable()	Checks if file can be openend or created for writing
importBuff()	Imports buffer file to Databuff struct
exportBuff()	Exports Databuff struct to buffer file

C.2.5. Calculations in SimulationCalc.cpp etc.

SimulationCalc.cpp	
getCovering()	Get covering from hitbuffer
getHits()	Get number of hits from hitbuffer
getnbDesorbed()	Get number of total desorbed molecules from hitbuffer
calcNmono()	see equation A.2
calcdNsurf()	see equation A.3
calcCoverage()	see equation A.5
calcStep()	see equation A.6
calcEnergy()	see equation A.7
calcStickingnew()	sets sticking coefficient to p→sticking
calcDesorption(), calcDesorptionRate()	see equation A.10
GetMoleculesPerTP()	see equation A.12
calctotalDesorption	calculates desorption for `startFromSource()`
calcPressure(), calcParticleDensity()	TODO has to be verified

worker.cpp	
CalcTotalOutgassingWorker()	see equation A.11, calculates outgassing for `startFromSource()`

SimulationLinux.cpp	
convertunit()	Converts simutime*unit to milliseconds

C.2.6. UpdateSubProcess.cpp

Update sHandle paramters from hitbuffer	
UpdateSticking()	Updates sticking
UpdateDesorptionRate()	Updates desorption rate
UpdateSojourn()	Updates sojourn freqency and energy

Error calculations	
UpdateErrorSub()	Updates error for current iteration, see equation A.14
UpdateError()	Sums up error of normal facets & weights by facet area
CheckErrorSub()	Checks if normal & vip facets reached respective target

Update hitbuffer	
initbufftozero()	Sets hitbuffer except covering to zero
UpdateMCSubHits()	Saves simulation results from sHandle into hitbuffer

C.2.7. UpdateMainProcess.cpp

Update main hitbuffer from sub hitbuffer	
UpdateMCMainHits()	Add simulation results from sub hitbuffer to main hitbuffer

Update real covering in hitbuffer	
getStepSize()	Calculates step size for current step, see equation A.13
manageStepSize()	Adapts step size if desRate · step size > than covering
UpdateCovering()	Uses step size and Krealvirt to calculate new covering Saved to `simHistory→coveringList`
UpdateCoveringphys()	Saves current real covering to hitbuffer
UpdateErrorMain()	Calculates total error, see equation A.14 Saved to `simHistory→errorList`
CalcPerIteration()	Calculates total error and covering over all facets per iteration