

Implementation of decay heat submodel in Lagrangian library

Manohar Kampili

Institute for Energy and Climate Research (IEK-6),
Forschungszentrum Jülich,
Germany

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Outline

1 Introduction

2 Theory

3 Insight into Lagrangian library

4 Creating a new submodel

5 Solver and test case

Introduction

- Aerosol: solid/liquid particles suspended in air/gas
- Nuclear aerosol: Generated in nuclear severe accidents
- Main carriers of the fission products
- Examples: CsI, CsOH, Ag etc ..

Physics of aerosols (1)

Apart from fluid dynamics,

- Particle transport
- Heat transfer between particles and fluid
- Chemistry among particles
- Agglomeration & De-agglomeration

Physics of aerosols (2)

Additional physics involved for nuclear aerosols,

- Decay heat associated with the particles
- Fractal nature
- Shape factors
- Change of chemical composition due to radioactive decay

Selection of solver

In general, flows in reactor containment are weakly compressible and turbulent.

The particle cloud is dynamic in nature along with chemistry.

Description of reactingParcelFoam:

Transient solver for compressible, turbulent flow with a reacting, multiphase particle cloud, and optional sources/constraints

Additional physics need to be added to this solver.

Adding decay heat model is the focus of this tutorial.

Outline

- 1 Introduction
- 2 **Theory**
- 3 Insight into Lagrangian library
- 4 Creating a new submodel
- 5 Linking it to reactingParcelFoam
- 6 Test case

Theory - Particle transport

Particles are tracked using Lagrangian method by calculating the forces.

- Sphere Drag
- Lift
- Gravitational force
- Brownian motion

Theory - Heat transfer between particle and fluid (1)

Heat transfer mechanisms are

- Conduction (Fourier's law)
- Convection (Ranz-Marshall correlation)
- Radiation (Blackbody radiation)

For nuclear aerosol particles, additionally

- Decay heat transport

Theory - Heat transfer between particle and fluid (2)

Energy balance within the particle

$$mC_p \frac{dT}{dt} = \dot{Q}$$

m is mass of the particle,

C_p is specific heat of the particle,

T is temperature of the particle,

t is time,

\dot{Q} is overall heat transfer rate from particle to fluid by all modes of heat transfer.

Theory - Heat transfer between particle and fluid (3)

By considering convection alone

$$mC_p \frac{dT}{dt} = hA(T - T_c)$$

h is convective heat transfer coefficient,

A is surface area of the particle,

T_c is carrier fluid temperature.

Analytical solution by variables separable method:

$$T = T_c + (T_i - T_c) * \exp\left(\frac{-6ht}{\rho C_p d}\right)$$

T_i is initial temperature of the particle,

ρ is density of the particle,

d is diameter of the particle.

Theory - Heat transfer between particle and fluid (4)

Convective heat transfer coefficient (h) calculation:

$$h = \frac{(Nu)k}{d}$$

k is thermal conductivity of the particle.

Ranz-Marshall correlation for Nusselt number:

$$Nu = 2 + 0.6Re^{\frac{1}{2}}Pr^{\frac{1}{3}}$$

Re is particle Reynolds number,

Pr is Prandtl number.

Note: If there is no heat transfer, Nusselt number and heat transfer coefficient would be 0.

Theory - Decay heat model

Decay heat of the radioactive fission product inside aerosol particle is transported by α, β, γ .

In the present case, it is assumed that all the decay heat is absorbed within the particle.

Governing equation:

$$mC_p \frac{dT}{dt} = \dot{Q} + \text{decayPower}$$

decayPower is amount of heat generated inside the particle in units *W/particle*

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Classification of parcel clouds

Cloud	submodels
KinematicCloud	Forces on particles
ThermoCloud	KinematicCloud + Heat transfer
KinematicCollidingCloud	KinematicCloud + Collision models
ReactingCloud	ThermoCloud + Phase change + Variable composition (single phase)
ReactingMultiphaseCloud	ReactingCloud + Multiphase composition + Devolatilisation + Surface reactions

Classification of Lagrangian solvers

Solver	Cloud
uncoupledKinematicParcelFoam	KinematicCloud
coalChemistryFoam	ThermoCloud
DPMFoam	KinematicCollidingCloud
reactingParcelFilmFoam	ReactingCloud
reactingParcelFoam	ReactingMultiphaseCloud

Understanding heat transfer model implementation (1)

Heat transfer from particle to fluid is calculated in `calcHeatTransfer` function defined in `ThermoParcel.C` file.

This file located at
`$FOAM_SRC/lagrangian/intermediate/parcels/Templates/
ThermoParcel/`

Since this is a very big function, it is explained in parts here.

This function returns the particle new temperature after the heat transfer from particle to fluid in a specific timestep.

Understanding heat transfer model implementation (2)

```
260     if (!td.cloud().heatTransfer().active())
261     {
262         return T_;
263     }
```

It means, if the heat transfer model is not active (specified in `reactingCloud1Properties` in constant directory of case), then it returns the original temperatures of the particles as new temperatures.

Understanding heat transfer model implementation (3)

```
269 scalar htc = td.cloud().heatTransfer().htc(d, Re, Pr, kappa, NCpW);
270
271 if (mag(htc) < ROOTVSMALL && !td.cloud().radiation())
272 {
273     return
274         max
275         (
276             T_ + dt*Sh/(this->volume(d)*rho*Cp_),
277             td.cloud().constProps().TMin()
278         );
279 }
```

htc is heat transfer coefficient calculated from heatTransfer model. And the remaining code in if loop, calculates new temperatures of the particles if htc value is negligible and radiation model is not active. These new temperatures are due to enthalpy transfer by phase change Sh.

Understanding heat transfer model implementation (4)

```
281    htc = max(htc, ROOTVSMALL);  
282    const scalar As = this->areaS(d);  
283  
284    scalar ap = Tc_ + Sh/(As*htc);  
285    scalar bp = 6.0*(Sh/As + htc*(Tc_ - T_));
```

scalar As is surface area of the particle. ap, bp are intermediate values to calculate new temperatures using analytical solution.

For better understanding, let's assume there is no phase change, which means value of Sh is 0.

```
ap = Tc_  
bp = 6.0*htc*(Tc_ - T_)
```

Where Tc_ is carrier fluid temperature. T_ is particle initial temperature.

Understanding heat transfer model implementation (5)

```

if (td.cloud().radiation())
{
    tetIndices tetIs = this->currentTetIndices();
    const scalar Gc = td.GInterp().interpolate(this->position(), tetIs);
    const scalar sigma = physicoChemical::sigma.value();
    const scalar epsilon = td.cloud().constProps().epsilon0();

    // Assume constant source
    scalar s = epsilon*(Gc/4.0 - sigma*pow4(T_));

    ap += s/htc;
    bp += 6.0*s;
}
bp /= rho*d*Cp_*(ap - T_) + ROOTVSMALL;

```

This piece of code changes the values of ap and bp , if the radiation model is active. If it is not active, then $ap = T_{c_}$, $bp = 6*htc/(rho*d*Cp)$.

Understanding heat transfer model implementation (6)

```

302     IntegrationScheme<scalar>::integrationResult Tres =
303     td.cloud().TIntegrator().integrate(T_, dt, ap*bp, bp);

```

These two lines calculates the new temperatures of the particles using the analytical solution, which is equivalent to integration during a time step.

This integrate function is defined in

`$FOAM_SRC/lagrangian/intermediate/IntegrationScheme/Analytical/Analytical.C`

This is exactly equivalent to

$$T = T_c + (T_i - T_c) * \exp\left(\frac{-6ht}{\rho C_p d}\right)$$

Understanding heat transfer model implementation (7)

```
316     Sph = dt*htc*As;  
317  
318     dhsTrans += Sph*(Tres.average() - Tc_);
```

```
dhsTrans = htc * As * dt * (Tres.average() - Tc_)
```

`Tres.average()` is average particle temperature during timestep `dt`. `dhsTrans` is heat transferred from particle to fluid. Value of `htc` is obtained from the heat transfer model specified in `reactingCloud1Properties` file in `constant` directory of case.

This is exactly equivalent to

$$Q = hA(T - T_c)dt$$

Understanding heat transfer model implementation (8)

Heat transfer models are located at

`$FOAM_SRC/lagrangian/intermediate/submodels/Thermodynamic/HeatTransferModel`

In `NoHeatTransfer/NoHeatTransfer.C`

```
template<class CloudType>
Foam::scalar Foam::NoHeatTransfer<CloudType>::Nu
(
    const scalar,
    const scalar
) const
{
    return 0.0;
}
```

Nusselt number value returned is always zero, which is justified since there is no heat transfer.

Understanding heat transfer model implementation (9)

Ranz-Marshall heat transfer model is located at

\$FOAM_SRC/lagrangian/intermediate/submodels/Thermodynamic/HeatTransferModel

In RanzMarshall.C

```
template<class CloudType>
Foam::scalar Foam::RanzMarshall<CloudType>::Nu
(
    const scalar Re,
    const scalar Pr
) const
{
    return 2.0 + 0.6*sqrt(Re)*cbrrt(Pr);
}
```

Nusselt number value returned is a function of Reynolds number and Prandtl number, which is exactly equivalent to the equation showed in Theory (slide 12).

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Decay heat model implementation (1)

Decay heat of the particle is a source term to the energy balance equation.

$$mC_p \frac{dT}{dt} = \dot{Q} + decayPower$$

In order to implement the decay heat model explained in Theory, calculation ap , bp should be understood clearly for different modes of heat transfer

For convection alone:

$$ap = Tc$$

$$bp = \frac{6 * htc}{\rho * d * Cp}$$

Decay heat model implementation (3)

Following the same trend for decay heat contribution:

$$ap = Tc + \frac{Sh}{(As)(htc)} + \frac{\epsilon * (\frac{Gc}{4} - \sigma * T^4)}{htc} + \frac{decayPower}{(As)(htc)}$$

$$bp = \frac{6 * htc}{\rho * d * Cp} + 6(\frac{Sh}{As}) + 6 * \epsilon * (\frac{Gc}{4} - \sigma * T^4) + \frac{6 * decayPower}{As}$$

decayPower is in the units of W / particle.

Steps for developing new submodel

- User copy of Lagrangian library
- Identifying the files that need to be changed
- Modifying one by one file
- Compiling
- Linking it solver

User copy of Lagrangian library

Create an user copy of lagrangian library that is needed along with Make directory. Modify the files file in Make directory as following.

OF1706+

```
cp -r $FOAM_SRC/lagrangian/intermediate $WM_PROJECT_USER_DIR/src
```

```
cd $WM_PROJECT_USER_DIR/src/lagrangian/intermediate/Make
```

```
sed -i s/FOAM_LIB_BIN/FOAM_USER_LIB_BIN/g files
```

```
sed -i s/liblagrangianIntermediate/libmyLagrangianIntermediate/g files
```

Identifying files for modifying

In this tutorial, directory structure of heat transfer model is replicated for developing decay heat models.

Let's find out heatTransfer keyword using grep command.

```
cd $WM_PROJECT_USER_DIR/src/lagrangian/intermediate  
grep -r -n heatTransfer .
```

This will result different instances of heatTransfer in following files:

- ThermoCloud.H
- ThermoCloudI.H
- ThermoCloud.C
- ThermoParcel.C
- ReactingMultiphaseParcel.C
- ReactingParcel.C
- HeatTransferModelNew.C
- HeatTransferModel.H

Modifications in ThermoCloud.H (1)

```
vi $WM_PROJECT_USER_DIR/src/lagrangian/intermediate/clouds/  
Templates/ThermoCloud/ThermoCloud.H
```

- First modification is forward declaration of DecayHeatModel class (just below the HeatTransferModel) as following:

```
template<class CloudType>  
class HeatTransferModel;
```

```
template<class CloudType>  
class DecayHeatModel;
```

Modifications in ThermoCloud.H (2)

- Second modification is adding `decayHeatModel_` in protected member data

```
//- Heat transfer model
autoPtr<HeatTransferModel<ThermoCloud<CloudType>>>
heatTransferModel_;
//- decay heat model
autoPtr<DecayHeatModel<ThermoCloud<CloudType>>>
decayHeatModel_;
```

- Third modification is adding `decayHeat()` member function

```
// Sub-models
//- Return reference to heat transfer model
inline const HeatTransferModel<ThermoCloud<CloudType>>&
heatTransfer() const;
//- Return reference to decay heat model
inline const DecayHeatModel<ThermoCloud<CloudType>>&
decayHeat() const;
```

Modification in ThermoCloudI.H (1)

- Defining the inline function `decayHeat()`:

```
template<class CloudType>
inline const Foam::HeatTransferModel<Foam::ThermoCloud<CloudType>>&
Foam::ThermoCloud<CloudType>::heatTransfer() const
{
    return heatTransferModel_;
}

template<class CloudType>
inline const Foam::DecayHeatModel<Foam::ThermoCloud<CloudType>>&
Foam::ThermoCloud<CloudType>::decayHeat() const
{
    return decayHeatModel_;
}
```

Modifications in ThermoCloud.C (1)

- First modification is including DecayHeatModel.H in the header files

```
#include "HeatTransferModel.H"  
#include "DecayHeatModel.H"
```

- Second modification is adding decayHeatModel_ in cloudReset member function

```
heatTransferModel_.reset(c.heatTransferModel_.ptr());  
  
decayHeatModel_.reset(c.decayHeatModel_.ptr());
```

Modifications in ThermoCloud.C (2)

- Third modification is in the first member function, as following:

```
heatTransferModel_.reset
(
    HeatTransferModel<ThermoCloud<CloudType>>::New
    (
        this->subModelProperties(),
        *this
    ).ptr()
);

decayHeatModel_.reset
(
    DecayHeatModel<ThermoCloud<CloudType>>::New
    (
        this->subModelProperties(),
        *this
    ).ptr()
);
```

Modifications in ThermoCloud.C (3)

- Modification in first constructor

```
heatTransferModel_(nullptr),  
decayHeatModel_(nullptr),
```

- Modification in second constructor

```
heatTransferModel_(c.heatTransferModel_->clone()),  
decayHeatModel_(c.decayHeatModel_->clone()),
```

- Modification in third constructor

```
heatTransferModel_(nullptr),  
decayHeatModel_(nullptr),
```

Modification in ThermoParcel.C

In calcHeatTransfer function, modify the few lines in the middle as following:

```
..
```

```
..
```

```
htc = max(htc, ROOTVSMALL);
```

```
const scalar As = this->areaS(d);
```

```
const scalar Vs = this->volume(d);
```

```
scalar decay= td.cloud().decayHeat().decayPower();
```

```
scalar ap = Tc_ + (Sh/(As*htc)) + (decay/(As*htc));
```

```
scalar bp = 6.0*(Sh/As + htc*(Tc_ - T_)+(decay/As));
```

```
if (td.cloud().radiation())
```

```
{
    tetIndices tetIs = this->currentTetIndices();
    const scalar Gc = td.GInterp().interpolate(this->position(), tetIs)
```

```
..
```

```
..
```

Creating DecayHeatModel directory

HeatTransferModel directory is taken as basis for creating DecayHeatModel directory. In these files, replace all the instances of heatTransfer with decayHeat, HeatTransfer with DecayHeat, htm with dhm, RanzMarshall with constantDecayHeat.

```
cd $WM_PROJECT_USER_DIR/src/intermediate/submodels/Thermodynamic
```

```
cp -r HeatTransferModel DecayHeatModel
```

```
cd DecayHeatModel
```

```
sed -i s/heatTransfer/decayHeat/g *
```

```
sed -i s/HeatTransfer/DecayHeat/g *
```

```
sed -i s/htm/dhm/g *
```

```
sed -i s/RanzMarshall/constantDecayHeat/g *
```


DecayHeatModel directory

In this DecayHeatModel directory, we have 3 sub-directories with following files:

- DecayHeatModel
 - DecayHeatModel.H
 - DecayHeatModel.C
 - DecayHeatModelNew.C
- NoDecayHeat
 - NoDecayHeat.H
 - NoDecayHeat.C
- constantDecayHeat
 - constantDecayHeat.H
 - constantDecayHeat.C

Modifications in DecayHeatModel.H and DecayHeatModel.C

- Declare a member function `decayPower()` and delete old member functions if any in `DecayHeatModel.H`.

```
// Member Functions
// Evaluation
//- Retrun decay Power
virtual scalar decayPower() const = 0;
```

- Define the member function `decayPower()` as following in `DecayHeatModel.C`:

```
// * * * * * Member Functions * * * * * //

template<class CloudType>
Foam::scalar Foam::DecayHeatModel<CloudType>::decayPower() const
{
    const scalar decayPower_ = this -> decayPower();
    return decayPower_;
}
```

Modification in DecayHeatModelNew.C

Change the error message as following:

FatalErrorInFunction

```
<< "Unknown decay heat model type"
<< modelType << nl << nl
<< "Valid decay heat model types are:" << nl
<< dictionaryConstructorTablePtr_->sortedToc()
<< exit(FatalError);
```

Modifications in NoDecayHeat.H and NoDecayHeat.C

- Declare a member function `decayPower()` and delete old member functions in `NoDecayHeat.H`.

```
// Member Functions
```

```
//- Flag to indicate whether model activates heat transfer model  
virtual bool active() const;
```

```
//- decay Power  
virtual scalar decayPower() const;
```

- Define the member function `decayPower()` as following in `NoDecayHeat.C`: This function should return zero always.

```
// * * * * * Member Functions * * * * * //
```

```
template<class CloudType>  
Foam::scalar Foam::NoDecayHeat<CloudType>::decayPower() const  
{  
    return 0.0;  
}
```

Modifications in constantDecayHeat.H

- Declare a new member function `decayPower()` and delete old member functions in `constantDecayHeat.H`.

```
// Member Functions
```

```
//- Flag to indicate whether model activates heat transfer model
virtual bool active() const;
```

```
//- decay Power
virtual scalar decayPower() const;
```

- Declare a new member data `decayPower_`

```
template<class CloudType>
class constantDecayHeat
:
public DecayHeatModel<CloudType>
{
//private data
```

```
scalar decayPower_;
```

Modification in constantDecayHeat.C (1)

- Initialize the decayPower_ with input value in the costructor as following:

```
// * * * * * Constructors * * * * * //

template<class CloudType>
Foam::exponentialDecayHeat<CloudType>::constantDecayHeat
(
    const dictionary& dict,
    CloudType& cloud
)
:
    DecayHeatModel<CloudType>(dict, cloud, typeName),
    decayPower_(readScalar(this->coeffDict().lookup("decayPower")))
{}

```

Modification in constantDecayHeat.C (2)

- Define the member function `decayPower()` as following in `constantDecayHeat.C` to return private member data `decayPower_`

```
// * * * * * Member Functions * * * * *

template<class CloudType>
bool Foam::constantDecayHeat<CloudType>::active() const
{
    return false;
}

template<class CloudType>
Foam::scalar Foam::constantDecayHeat<CloudType>::decayPower() const
{
    return decayPower_;
}
```

Compiling decay heat models

Heat transfer models are compiled using
`makeParcelHeatTransferModels.H` located at

`$WM_PROJECT_USER_DIR/src/intermediate/parcels/include/`

Following the same trend, create a file `makeParcelDecayHeatModels.H` in the same directory.

Code in makeParcelDecayHeatModels.H

```

#ifndef makeParcelDecayHeatModels_H
#define makeParcelDecayHeatModels_H
// * * * * *
#include "NoDecayHeat.H"
#include "constantDecayHeat.H"
// * * * * *

#define makeParcelDecayHeatModels(CloudType) \
\
    makeDecayHeatModel(CloudType); \
\
    makeDecayHeatModelType(NoDecayHeat, CloudType); \
    makeDecayHeatModelType(constantDecayHeat, CloudType);

// * * * * *
#endif

```

Linking ParcelDecayHeatModels to ThermoParcelSubmodels

In order to link the decay heat models to ThermoParcelSubmodels, there are two modifications required in **makeBasicThermoParcelSubmodels.C**, located in the directory at

```
$WM_PROJECT_USER_DIR/src/lagrangian/intermediate/parcels/derived/
basicThermoParcel/
```

- Including header file makeParcelDecayHeatModels.H

```
// Thermodynamic
#include "makeParcelHeatTransferModels.H"
#include "makeParcelDecayHeatModels.H"
```

- Adding thermo variant of parcelDecayHeatModels as following:

```
// Thermo sub-models
makeParcelHeatTransferModels(basicThermoCloud);

// decay heat models
makeParcelDecayHeatModels(basicThermoCloud);
```

Linking ParcelDecayHeatModels to ReactingMultiphaseParcelSubmodels

To link the decay heat models to ReactingMultiphaseParcelSubmodels, there are two modifications required in **makeBasicReactingMultiphaseParcelSubmodels.C**, located in the directory at

```
$WM_PROJECT_USER_DIR/src/lagrangian/intermediate/parcels/derived/  
basicReactingMultiphaseParcel/
```

- Including header file makeParcelDecayHeatModels.H

```
// Thermodynamic  
#include "makeParcelHeatTransferModels.H"  
#include "makeParcelDecayHeatModels.H"
```

- Adding thermo variant of parcelDecayHeatModels as following:

```
// Thermo sub-models  
makeParcelHeatTransferModels(basicReactingMultiphaseCloud);  
  
// decay heat models  
makeParcelDecayHeatModels(basicReactingMultiphaseCloud);
```

Modifications in Make/options file for compilation

Modifications required in options file in order to include all the files related to decayHeatModels.

```
EXE_INC = \  
-I. \  
-Ilagrangian/intermediate/lnInclude \  
..  
..
```

and

```
LIB_LIBS = \  
-L$(FOAM_USER_LIBBIN) \  
..  
..
```

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Solver and test case

Creating myReactingParcelFoam:

```
cp -r $FOAM_SOLVERS/lagrangian/reactingParcelFoam $WM_PROJECT_USER_DIR/
applications
```

```
cd $WM_PROJECT_USER_DIR/applications
```

```
rm -rf reactingParcelFoam/simpleReactingParcelFoam
```

```
sed -i s/reactingParcelFoam/myReactingParcelFoam/g *
```

```
sed -i s/FOAM_APPBIN/FOAM_USER_APPBIN/g *
```

Now user version of reactingParcelFoam is ready to compile with name myReactingParcelFoam.

Compiling solver with Lagrangian library

- Copy the new Lagrangian library into solver:

```
cp -r $WM_PROJECT_USER_DIR/src/lagrangian $WM_PROJECT_USER_DIR/
applications/myReactingParcelFoam
```

```
cd $WM_PROJECT_USER_DIR/applications/myReactingParcelFoam
```

- Modify options file in Make directory of myReactingParcelFoam

```
EXE_INC = \
-I. \
-Ilagrangian/intermediate/lnInclude \
```

and

```
EXE_LIBS = \
-L$(FOAM_USER_APPBIN) \
-lmyLagrangianIntermediate \
```

- Compile

```
wmake all
```

verticalChannel case without decayHeat (1)

```
cd $WM_PROJECT_USER_DIR/run
```

```
cp -r $FOAM_TUTORIALS/lagrangian/reactingParcelFoam/verticalChannel  
verticalChannelWithoutDecay
```

```
cd verticalChannelWithoutDecay
```

- Modifications in constant/reactingCloud1Properties file:

```
constantProperties
{
    rho0          4510;
    T0            473;
    Cp0           200;

    constantVolume true;
}
```


verticalChannel case without decayHeat (2)

- Modifications in injectionModels dictionary of constant/reactingCloud1Properties file:

```

injectionModels
{
    model1
    {
        type                patchInjection;
        massTotal            2.36143e-5;
        parcelBasisType     fixed;
        SOI                  0.0;
        duration            0.1;
        nParticle            1;
        parcelsPerSecond    100;
        patch               inletCentral;
        U0                   (0 1.08 0);
        flowRateProfile      constant 1;
        sizeDistribution
        {
            type            uniform;
            uniformDistribution
            {
                minValue    1e-3;
                maxValue    1e-3;
            }
        }
    }
}

```

verticalChannel case without decayHeat (3)

- Modifications in constant/reactingCloud1Properties file:

```
decayHeatModel none;

singleMixtureFractionCoeffs
{
    phases
    (
        gas
        {
        }
        liquid
        {
        }
        solid
        {
            ash 1;
        }
    );
    YGasTot0      0;
    YLiquidTot0    0;
    YSolidTot0     1;
}
```

verticalChannel case with decayHeat

```
cd $WM_PROJECT_USER_DIR/run

cp -r verticalChannelWithoutDecay verticalChannelWithDecay

vi verticalChannelWithDecay/constant/reactingCloud1Properties
```

- Modifications in constant/reactingCloud1Properties file:

```
decayHeatModel constantDecayHeat;
```

```
constantDecayHeatCoeffs
{
    decayPower 100000;
}
```

Here decayPower 100000 W/particle is taken as a random value to see the effect of decayHeat in 0.1 s. Run both the cases with myReactingParcelFoam solver

```
cd verticalChannelWithoutDecay

myReactingParcelFoam >& log &

cd ..

cd verticalChannelWithDecay

myReactingParcelFoam >& log &
```

PostProcessing with foamLog and gnuplot (1)

Tracking the highest and lowest temperatures of the particles

- copy paste the foamLog.db file to case

```
cp $WM_PROJECT_DIR/bin/tools/foamLog.db verticalChannelWithoutDecay
```

```
vi verticalChannelWithoutDecay/foamLog.db
```

- Add the following lines in the end to existing foamLog.db file:

```
#particle minimum Temp
particleMinTemp/Temperature min/max      =
```

```
#particle maximum Temp
particleMaxTemp/Temperature min/,
```

```
cp verticalChannelWithoutDecay/foamLog.db verticalChannelWithDecay
```

- Running foamLog in both cases

```
foamLog -case verticalChannelWithoutDecay
```

```
foamLog -case verticalChannelWithDecay
```

PostProcessing with foamLog and gnuplot (2)

Tracking the highest and lowest temperatures of the particles

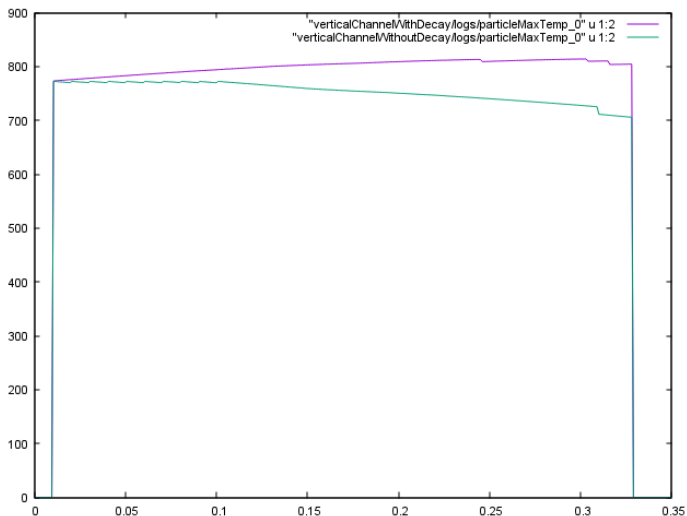
- Plotting with gnuplot

gnuplot

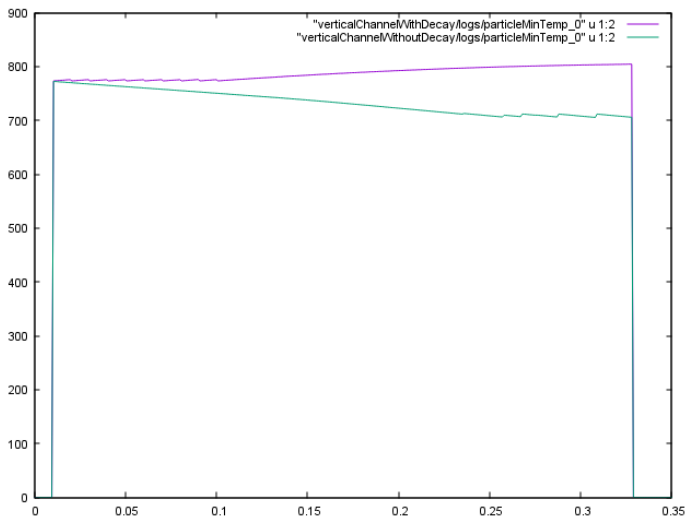
```
plot "verticalChannelWithoutDecay/postProcessing/logs/particleMinTemp.xy" u 1:2 w l,
"verticalChannelWithDecay/postProcessing/logs/particleMinTemp.xy" u 1:2 w l
```

```
plot "verticalChannelWithoutDecay/postProcessing/logs/particleMaxTemp.xy" u 1:2 w l,
"verticalChannelWithDecay/postProcessing/logs/particleMaxTemp.xy" u 1:2 w l
```

Maximum temperatures comparison



Minimum temperatures comparison



Thank you