

Namelist "Radiative_Heat_Transfer_DOM"

Not for the release SUNFLUIDH_EDU .

This data set is used to define the radiative problem. Otherwise, it can be omitted.

This module considers the Radiative Transfer Equation (RTE) for an emitting-absorbing non-scattering medium enclosed by diffuse boundaries. To take into account the gas behavior, it considers both gray-gas assumption as well as real gas behavior through the Spectral-Line-Weighted-Sum-of-Gray-Gases (SLW) model.

The final RTE-SLW problem is then discretized with the Discrete Ordinates Method (DOM).

The DOM discretizes the 4π steradians integration in a set of M discrete directions represented by their direct cosines and corresponding weights $\vec{q}_m = (\vec{s}_m, \omega_m) = (\mu_m, \eta_m, \xi_m, \omega_m)$ for all $m \in [1, M]$.

The SLW model will change the spectral integration in a weighted sum of N_g gray-gases represented by their absorption coefficient and corresponding weights (κ_j, a_j) for all $j \in [1, N_g]$.

Thus, the resulting **RTE-SLW-DOM** problem for emitting-absorbing non-scattering medium stands as below :

$$\begin{aligned} \vec{s}_m \cdot \nabla I_j^m(x_i, \vec{s}_m) = \kappa_j \left[a_j I_b(T(x_i)) - I_j^m(x_i, \vec{s}_m) \right]; \quad \forall (m, j) \in [M, N_g] \end{aligned}$$

where I_j^m is the radiative intensity for the virtual gray-gas j in direction m and I_b is the blackbody radiative intensity.

The **dimensional** radiative source term S_r and boundary net radiative heat flux q_r^{net} are defined as :

$$\begin{aligned} S_r(x_i, T) &= - \sum_{j=0}^{N_g} \kappa_j \left[\sum_{m=1}^M \omega_m I_j^m(x_i, \vec{s}_m) - 4 a_j \sigma_B(T(x_i))^4 \right] \\ &= \varepsilon_{wall} \left[\sigma_B(T(x_i^{wall}))^4 - \sum_{j=0}^{N_g} \sum_{m: \vec{s}_m \cdot \vec{n} > 0} \omega_m \vec{s}_m \cdot \vec{n} I_j^m(x_i^{wall}, \vec{s}_m) \right] \end{aligned}$$

where σ_B is the Stefan-Boltzmann constant, ε_{wall} is the boundary emissivity and \vec{n} is the normal to the wall pointing out of the domain.



This radiative solver implementation considers only **cartesian** problems and **does not support immersed bodies**.

Full data set of the namelist

```
&Radiative_Heat_Transfer_DOM activateRadiation = .false., RadiativePeriod
= 1, FirstIterations = 20,
                    RadiativeLocalIterations = 1, RadiativeConvergenceTolerance =
1.E-15,
```

```
    WallRadcoeff = 1.0 , VolRadCoeff = 1.0, RadiativeScheme =
"STEP",
    ActivateGas = .false., NbGas = 1, ka_max = 0.0, ka_min = 0.0,
    Pref = 101325.0, Tref = 300., Href = 1, speca = "H2O", xaref =
0.07, xaUniform = 0.07,
    Squad = 8, WallEmissivity = 0.0 0.0 0.0 0.0 0.0 0.0 /
```

Definition of the data set for the DOM-RTE problem

activateRadiation

- Type : Boolean value
- This option activates the radiative module.
 - .false. : no radiation considered
 - .true. : radiation problem is considered
- Default value = .false.

RadiativePeriod

- Type : Integer value
- This option set the periodicity of resolution of the Radiative problem in time iteration.
- Default value = 1

FirstIterations

- Type : Integer value
- In the case that no restart fields are available (start radiation from scratch), the solver will iterates “FirstIterations” times before entering the time loop.
- Default value = 20

RadiativeLocalIterations

- Type : Integer value
- Number of sub-iteration for the RTE solving at each radiative iteration.
- Default value = 1

RadiativeConvergenceTolerance

- Type : Real value
- Convergence criteria on the wall Fluxes and radiative source term for the sub-iteration.
- Default value = 1.E-15

WallRadcoeff

- Type : Real value
- Prescaler on the net radiative heat flux $\$q_r^{\{net\}}\$$ at walls.
- **For debugging only.**
- Default value = 1.0

VolRadcoeff

- Type : Real value
- Prescaler on the radiative source term $\$S_r\$$.
- **For debugging only.**
- Default value = 1.0

RadiativeScheme

- Type : Character string with a maximum size of 20
- Name of the interpolation scheme used in the Discrete Ordinates Method.
- Available values :
 - “STEP” : first order interpolation scheme (robust)
 - “DIAMOND” : second order centered interpolation scheme (could lead to negative intensity)
 - “LATHROP” : second order interpolation scheme with limiter (time-consuming)
- Default value = “STEP”

SQuad

- Type : Integer value.
- Order N of the level symmetric angular quadrature ($\$S_N\$$)
- This quadrature leads to $M = (N+2) \times N$ directions in volume and half at boundaries
- Available values are 2, 4, 6, 8, 10, 12, 14
- Default value = 8

Tref

- Type : Real value.
- Reference temperature $\$T_{\{ref\}}\$$ in [$\$K\$$].
- Default value = Fluid_Properties%Reference_Temperature

Href

- Type : Real value.
- Reference Length $\$H_{\{ref\}}\$$ in [$\$m\$$].
- Default value = Nondimensionalization%Reference_Length

WallEmissivity

- Type : Real array of size 6.
- Boundaries emissivities $\backslash varepsilon$ sorted as (x-,x+,y-,y+,z-,z+).
- Default values = 0.0 0.0 0.0 0.0 0.0 0.0

Definition of the data set for the SLW model

activateGas

- Type : Boolean value
- This option activates the SLW module.
 - .false. : **transparent medium** under gray-gas assumption is considered (i.e. $\kappa = 0$)
 - .true. : Gas absorption and emission is considered
- Default value = .false.



if activateGas == .false., the settings below are unnecessary.

NbGas

- Type : Integer value
- This option sets the number of weighted gray-gases N_g used in the SLW model.
 - NbGas $= 1$: gray-gas assumption with $\kappa = ka_{min}$
 - NbGas ≥ 2 : SLW model is employed
- Default value = 1



Setting **NbGas $= 1$ and $ka_{min} = 0$** is equivalent to **wall-to-wall radiation** due to the presence of transparent medium

ka_min , ka_max

- Type : Real values
- These options set the lower and upper bounds of dimensional absorbing coefficient [m^{-1}] for the SLW model.
 - if NbGas $= 1$: $\kappa = ka_{min}$, **ka_max is useless**
 - if NbGas ≥ 2 : $ka_{min} < \kappa_j < ka_{max}$ for all $j \in [1, N_g]$
- Default value = [ka_{min} , ka_{max}] = [0.0 , 0.0]

SPECA

- Type : Character string with a maximum size of 3
- Name of the absorbing species when NbGas ≥ 2 (SLW model).
 - **if “NbGas” $\neq 1$: useless**
- Available values :
 - “H2O” : \$air-H_2O\$ mixture
 - “CO2” : \$air-CO_2\$ mixture
- Default value = “H2O”

xaref

- Type : Real value
- This option set the reference molar fraction x_{ref} of the absorbing species for the SLW model.
 - **if “NbGas” $\neq 1$: useless**
- Default value = 0.07

xaUniform

- Type : Real value
- As long as the SLW model is not coupled with species equations, this option set a uniform molar fraction x_a of the absorbing species in the overall domain.
 - **if NbGas $\neq 1$: useless**
- Default value = 0.07

Pref

- Type : Real value.
- Reference pressure P_{ref} in [Pa].
 - **if “NbGas” $\neq 1$: useless**
- Default value = obtained from Fluid_Properties quantities

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<https://sunfluidh.lisn.upsaclay.fr/> - Documentation du code de simulation numérique SUNFLUIDH

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