```
| OpenFOAM: The Open Source CFD Toolbox
 || / F ield
 | / O peration /
  \parallel / A nd
               / www.openfoam.com
  W Manipulation /
  Copyright (C) 2017-2020 OpenCFD Ltd.
License
  This file is part of OpenFOAM.
  OpenFOAM is free software: you can redistribute it and/or modify it
  under the terms of the GNU General Public License as published by
  the Free Software Foundation, either version 3 of the License, or
  (at your option) any later version.
  OpenFOAM is distributed in the hope that it will be useful, but WITHOUT
  ANY WARRANTY; without even the implied warranty of MERCHANTABILITY or
  FITNESS FOR A PARTICULAR PURPOSE. See the GNU General Public License
  for more details.
  You should have received a copy of the GNU General Public License
  along with OpenFOAM. If not, see <a href="http://www.gnu.org/licenses/">http://www.gnu.org/licenses/>.</a>.
Class
  Foam::meltingEvaporationModels::LeeCNT
Description
  Mass transfer LeeCNT model. Simple model driven by field value difference as:
     |dot\{m\}| = C | rho | alpha (T - T_{activate}) / T_{activate} |
  \f7
  where C is a model constant.
  if C > 0:
     |dot{m} = C |rho |alpha (T - T_{activate})/T_{activate}
    for |f| T > T_{activate} |f|
  and
     |f[mDot = 0.0 | f] \text{ for } |f[T < T_{activate}| |f]
  if C < 0:
     \| \dot{m} = -C \| \rho \| \alpha \( T_{\activate} \) - T)/T_{\activate}
    for \f[ T < T_{activate} \f]
  and
```

```
|f| |dot\{m\}| = 0.0 |f| |f| |T| > T_{activate} |f|
  Based on the reference:
  -# W. H. LeeCNT. "A Pressure Iteration Scheme for Two-Phase Modeling".
  Technical Report LA-UR 79-975. Los Alamos Scientific Laboratory,
  Los Alamos, New Mexico. 1979.
Usage
  Example usage:
  \verbatim
    massTransferModel
      (solid to liquid)
        type
                 LeeCNT;
        C
                 40;
        Tactivate 302.78;
  \endverbatim
  Where:
  ltable
    Property | Description | Required | Default value
    Tactivate | Activation temperature | yes
    C | Model constant | yes
    includeVolChange | Volumen change | no
    species | Specie name on the other phase | no | none
  lendtable
SourceFiles
  LeeCNT.C
|*_____*/
#ifndef meltingEvaporationModels_LeeCNT_H
#define meltingEvaporationModels_LeeCNT_H
#include "InterfaceCompositionModel.H"
namespace Foam
namespace meltingEvaporationModels
{
             Class LeeCNT Declaration
template<class Thermo, class OtherThermo>
class LeeCNT
```

```
public InterfaceCompositionModel<Thermo, OtherThermo>
{
  // Private Data
    //- Condensation coefficient [1/s]
    dimensionedScalar C_;
    volScalarField interfaceVolume_;
    //- Phase transition temperature
    const dimensionedScalar Tactivate_;
    //- Phase minimum value for activation
    scalar alphaMin_;
    //- Planck constant [J.s]
    const dimensionedScalar planck_;
    //- Boltzmann constant [J/K]
    const dimensionedScalar boltzmann_;
    //- Activation energy of water molecules passing through water-ice interface [J]
    const dimensionedScalar deltag_;
    //- Number of water molecule in a water volume [m3]
    const dimensionedScalar nL_;
    //- Superficial free energy of the water-ice interface [J/m2]
    const dimensionedScalar gammaYW_;
    //- Latent heat per volume [J/m3]
    const dimensionedScalar hLV_;
    //- Shape coefficient of nucleation
    const dimensionedScalar alphaEY_;
public:
  //- Runtime type information
  TypeName("LeeCNT");
  // Constructors
    //- Construct from components
    LeeCNT
       const dictionary& dict,
       const phasePair& pair
    );
  //- Destructor
  virtual ~LeeCNT() = default;
```

```
// Member Functions
    //- Explicit total mass transfer coefficient
    virtual tmp<volScalarField> Kexp
       const volScalarField& field
    );
    //- Implicit mass transfer coefficient
    virtual tmp<volScalarField> KSp
       label modelVariable,
       const volScalarField& field
    );
    //- Explicit mass transfer coefficient
    virtual tmp<volScalarField> KSu
       label modelVariable,
       const volScalarField& field
    );
    //- Return T transition between phases
    virtual const dimensionedScalar& Tactivate() const;
    //- Add/subtract alpha*div(U) as a source term
    //- for alpha, substituting div(U) = mDot(1/rho1 - 1/rho2)
    virtual bool includeDivU();
    virtual const dimensionedScalar& planck() const;
    virtual const dimensionedScalar& boltzmann() const;
    virtual const dimensionedScalar& deltag() const;
    virtual const dimensionedScalar& nL() const;
    virtual const dimensionedScalar& gammaYW() const;
    virtual const dimensionedScalar& hLV() const;
    virtual const dimensionedScalar& alphaEY() const;
};
} // End namespace meltingEvaporationModels
} // End namespace Foam
```

#ifdef NoRepository # include "LeeCNT.C" #endif
//***********************
#endif
// ************************************