

# SPHERLSgen Reference Manual

## 1.0

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# Chapter 1

## SPHERLSgen Class Index

### 1.1 SPHERLSgen Class List

Here are the classes, structs, unions and interfaces with brief descriptions:

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## Chapter 2

# SPHERLSgen File Index

### 2.1 SPHERLSgen File List

Here is a list of all documented files with brief descriptions:

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## Chapter 3

# SPHERLSgen Class Documentation

### 3.1 eos Class Reference

```
#include <eos.h>
```

#### Public Member Functions

- [eos](#) ()
- [eos](#) (int [nNumT](#), int [nNumRho](#))
- [eos](#) (const [eos](#) &ref)
- [~eos](#) ()
- [eos & operator=](#) (const [eos](#) &eosRightSide)
- void [readAscii](#) (std::string sFileName)
- void [readBobsAscii](#) (std::string sFileName)
- void [writeAscii](#) (std::string sFileName)
- void [readBin](#) (std::string sFileName) throw (exception2)
- void [writeBin](#) (std::string sFileName)
- double [dGetPressure](#) (double dT, double dRho)
- double [dGetEnergy](#) (double dT, double dRho)
- double [dGetOpacity](#) (double dT, double dRho)
- double [dDRhoDP](#) (double dT, double dRho)
- double [dSoundSpeed](#) (double dT, double dRho)
- void [getEKappa](#) (double dT, double dRho, double &dE, double &dKappa)
- void [getPEKappa](#) (double dT, double dRho, double &dP, double &dE, double &dKappa)
- void [getPEKappaGamma](#) (double dT, double dRho, double &dP, double &dE, double &dKappa, double &dGamma)
- void [getPKappaGamma](#) (double dT, double dRho, double &dP, double &dKappa, double &dGamma)
- void [gamma1DelAdC\\_v](#) (double dT, double dRho, double &dGamma1, double &dDelAd, double &dC\_v)
- void [getPAndDRhoDP](#) (double dT, double dRho, double &dP, double &dDRhoDP)
- void [getEAndDTDE](#) (double dT, double dRho, double &dE, double &dTDE)

## Public Attributes

- int `nNumRho`
- int `nNumT`
- double `dXMassFrac`
- double `dYMassFrac`
- double `dLogRhoMin`
- double `dLogRhoDelta`
- double `dLogTMin`
- double `dLogTDelta`
- double \*\* `dLogP`
- double \*\* `dLogE`
- double \*\* `dLogKappa`

### 3.1.1 Detailed Description

This class holds an equation of state as well as many functions useful for manipulating it

### 3.1.2 Constructor & Destructor Documentation

#### 3.1.2.1 `eos::eos ()`

Constructor, doesn't really do anything

#### 3.1.2.2 `eos::eos (int nNumT, int nNumRho)`

Constructor, allocates memory for the 2D arrays

#### Parameters:

- ← *nNumT* number of temperatures in the equaiton of state table
- ← *nNumRho* number of densities in the equaiton of state table

#### 3.1.2.3 `eos::eos (const eos & ref)`

Copy constructor, simply constructs a new eos object from another eos object

#### 3.1.2.4 `eos::~~eos ()`

Destructor, deletes dynamic arrays

### 3.1.3 Member Function Documentation

**3.1.3.1 double eos::dDRhoDP (double  $dT$ , double  $dRho$ )**

This function calculates the partial derivative of density w.r.t. pressure

**Parameters:**

- ←  $dT$  temperature at which the derivative is to be computed
- ←  $dRho$  density at which the derivative is to be computed

**Returns:**

the partial derivative of density w.r.t. pressure.

**3.1.3.2 double eos::dGetEnergy (double  $dT$ , double  $dRho$ )**

This function linearly interpolates the energy to a given temperature and density. Note that both  $dT$  and  $dRho$  are not in log space.

**Parameters:**

- ←  $dT$  temperature to interpolate to.
- ←  $dRho$  density to interpolate to.

**Returns:**

the interpolated energy.

**3.1.3.3 double eos::dGetOpacity (double  $dT$ , double  $dRho$ )**

This function linearly interpolates the opacity to a given temperature and density. Note that both  $dT$  and  $dRho$  are not in log space.

**Parameters:**

- ←  $dT$  temperature to interpolate to.
- ←  $dRho$  density to interpolate to.

**Returns:**

the interpolated opacity.

**3.1.3.4 double eos::dGetPressure (double  $dT$ , double  $dRho$ )**

This function linearly interpolates the pressure to a given temperature and density. Note that both  $dT$  and  $dRho$  are not in log space.

**Parameters:**

- ←  $dT$  temperature to interpolate to.
- ←  $dRho$  density to interpolate to.

**Returns:**

the interpolated pressure.

### 3.1.3.5 double eos::dSoundSpeed (double $dT$ , double $dRho$ )

This function calculates the adiabatic sound speed

**Parameters:**

- ←  $dT$  temperature at which the derivative is to be computed
- ←  $dRho$  density at which the derivative is to be computed

**Returns:**

the sound speed.

### 3.1.3.6 void eos::gamma1DelAdC\_v (double $dT$ , double $dRho$ , double & $dGamma1$ , double & $dDelAd$ , double & $dC_v$ )

This function calculates gamma1 and the adiabatic gradient

**Parameters:**

- ←  $dT$  temperature at which the derivative is to be computed
- ←  $dRho$  density at which the derivative is to be computed
- $dGamma1$  gamma1
- $dDelAd$  adiabatic gradient
- $dC_v$  specific heat at constant volume

### 3.1.3.7 void eos::getEAndDTDE (double $dT$ , double $dRho$ , double & $dE$ , double & $dDTDE$ )

This function calculates the partial derivative of temperature w.r.t. energy and the energy

**Parameters:**

- ←  $dT$  temperature at which the derivative is to be computed
- ←  $dRho$  density at which the derivative is to be computed
- $dE$  energy at  $dT$  and  $dRho$
- $dDTDE$  derivative of temperature w.r.t. energy at constant density

### 3.1.3.8 void eos::getEKappa (double $dT$ , double $dRho$ , double & $dE$ , double & $dKappa$ )

This function linearly interpolates the three dependent quantities (Pressure, Energy , Opacity) to a given temperature and density. Note that both  $dT$  and  $dRho$  are not in log space.

**Parameters:**

- ←  $dT$  temperature to interpolate to.
- ←  $dRho$  density to interpolate to.
- $dE$  energy at  $dT$  and  $dRho$ .
- $dKappa$  opacity at  $dT$  and  $dRho$ .

### 3.1.3.9 void eos::getPAndDRhoDP (double *dT*, double *dRho*, double & *dP*, double & *ddRhoDP*)

This function calculates the partial derivative of density w.r.t. pressure and the pressure

#### Parameters:

- ← *dT* temperature at which the derivative is to be computed
- ← *dRho* density at which the derivative is to be computed
- *dP* pressure at *dT* and *dRho*
- *ddRhoDP* derivative of density w.r.t. pressure at constant temperature

### 3.1.3.10 void eos::getPEKappa (double *dT*, double *dRho*, double & *dP*, double & *dE*, double & *dKappa*)

This function linearly interpolates the three dependent quantities (Pressure, Energy , Opacity) to a given temperature and density. Note that both *dT* and *dRho* are not in log space.

#### Parameters:

- ← *dT* temperature to interpolate to.
- ← *dRho* density to interpolate to.
- *dP* pressure at *dT* and *dRho*.
- *dE* energy at *dT* and *dRho*.
- *dKappa* opacity at *dT* and *dRho*.

### 3.1.3.11 void eos::getPEKappaGamma (double *dT*, double *dRho*, double & *dP*, double & *dE*, double & *dKappa*, double & *dGamma*)

This function linearly interpolates the energy and opacity to a given temperature and density. Note that both *dT* and *dRho* are not in log space.

#### Parameters:

- ← *dT* temperature to interpolate to.
- ← *dRho* density to interpolate to.
- *dP* pressure at *dT* and *dRho*.
- *dE* energy at *dT* and *dRho*.
- *dKappa* opacity at *dT* and *dRho*.
- *dGamma* adiabatic index at *dT* and *dRho*.

### 3.1.3.12 void eos::getPKappaGamma (double *dT*, double *dRho*, double & *dP*, double & *dKappa*, double & *dGamma*)

This function linearly interpolates the energy and opacity to a given temperature and density. Note that both *dT* and *dRho* are not in log space.

#### Parameters:

- ← *dT* temperature to interpolate to.
- ← *dRho* density to interpolate to.
- *dP* pressure at *dT* and *dRho*.
- *dKappa* opacity at *dT* and *dRho*.
- *dGamma* adiabatic index at *dT* and *dRho*.

### 3.1.3.13 eos & eos::operator= (const eos & *eosRightSide*)

Assignment operator, assigns one eos object to another.

### 3.1.3.14 void eos::readAscii (std::string *sFileName*)

This fuction reads in an ascii file and stores it in the current object.

#### Parameters:

- ← *sFileName* name of the equation of state file to read from.

### 3.1.3.15 void eos::readBin (std::string *sFileName*) throw (exception2)

This fuction reads in a binary file and stores it in the current object.

#### Parameters:

- ← *sFileName* name of the equation of state file to read from.

### 3.1.3.16 void eos::readBobsAscii (std::string *sFileName*)

This fuction reads in an ascii file and stores it in the current object. The ascii file is in Bob's format.

#### Parameters:

- ← *sFileName* name of the equation of state file to read from.

**3.1.3.17 void eos::writeAscii (std::string *sFileName*)**

This function writes the equation of state stored in the current object to an ascii file.

**Parameters:**

← *sFileName* name of the file to write the equation of state to.

**3.1.3.18 void eos::writeBin (std::string *sFileName*)**

This function writes the equation of state stored in the current object to a binary file.

**Parameters:**

← *sFileName* name of the file to write the equation of state to.

**3.1.4 Member Data Documentation****3.1.4.1 double\*\* eos::dLogE**

2D array of log10 energies. `dLogE[i][j]` gives the log10 energy at log10 density of `eos::dLogRhoDelta*i+eos::dLogRhoMin`, and at log10 temperature of `eos::dLogTDelta*j+eos::dLogTMin`.

**3.1.4.2 double\*\* eos::dLogKappa**

2D array of log10 opacities. `dLogKappa[i][j]` gives the log10 opacity at log10 density of `eos::dLogRhoDelta*i+eos::dLogRhoMin`, and at log10 temperature of `eos::dLogTDelta*j+eos::dLogTMin`.

**3.1.4.3 double\*\* eos::dLogP**

2D array of log10 pressures. `dLogP[i][j]` gives the log10 pressure at log10 density of `eos::dLogRhoDelta*i+eos::dLogRhoMin`, and at log10 temperature of `eos::dLogTDelta*j+eos::dLogTMin`.

**3.1.4.4 double eos::dLogRhoDelta**

Increment of the density between table entries in log10.

**3.1.4.5 double eos::dLogRhoMin**

Minimum density of the table in log10.

**3.1.4.6 double eos::dLogTDelta**

Increment of the temperature between table entries in log10.

**3.1.4.7 double [eos::dLogTMin](#)**

Minimum temperature of the table in log10.

**3.1.4.8 double [eos::dXMassFrac](#)**

Hydrogen mass fraction of the composition used to generate the equation of state table.

**3.1.4.9 double [eos::dYMassFrac](#)**

Helium mass fraction of the composition used to generate the equation of state table.

**3.1.4.10 int [eos::nNumRho](#)**

Number of densities in the equation of state table

**3.1.4.11 int [eos::nNumT](#)**

Number of temperatures in the equation of state table

The documentation for this class was generated from the following files:

- [/home/cgeroux/SPHERLS\\_new/src/eos.h](#)
- [/home/cgeroux/SPHERLS\\_new/src/eos.cpp](#)



## Chapter 4

# SPHERLSgen File Documentation

### 4.1 /home/cgeroux/SPHERLS\_new/src/eos.cpp File Reference

```
#include <string>
#include <fstream>
#include <sstream>
#include <iostream>
#include <cmath>
#include "eos.h"
#include "exception2.h"
```

#### 4.1.1 Detailed Description

Implements the eos (equation of state) class defined in [eos.h](#)

## 4.2 /home/cgeroux/SPHERLS\_new/src/eos.h File Reference

```
#include <string>
#include "exception2.h"
```

### Classes

- class [eos](#)

### 4.2.1 Detailed Description

Header file for [eos.cpp](#)

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