Group 4

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3.1 File List

Here is a list of all files with brief descriptions:

Chapter 4

Namespace Documentation

4.1 constants Module Reference

Module contains definitions of useful constants.

Variables

- integer, parameter dp =selected_real_kind(15, 300)

  Double-precision real kind.
- real(kind=dp), parameter pi = 4.0_dp∗atan(1.0_dp)

  The circle constant, pi.
- real(kind=dp), parameter tau = 2.0_dp∗pi

  2∗pi

4.1.1 Detailed Description

Module contains definitions of useful constants.

This program creates a CASTEP cell file consisting of a random arrangement of Ca and Mg atoms in a crystal structure with oxygen, according to a user specification.

4.1.2 Variable Documentation

4.1.2.1 integer, parameter constants::dp =selected_real_kind(15, 300)

Double-precision real kind.

Definition at line 12 of file constants.f90.

4.1.2.2 real(kind=dp), parameter constants::pi = 4.0_dp∗atan(1.0_dp)

The circle constant, pi.

Definition at line 15 of file constants.f90.

4.1.2.3 real(kind=dp), parameter constants::tau = 2.0_dp∗pi

2∗pi

Definition at line 16 of file constants.f90.

4.2 convex_hull Namespace Reference

Variables

- string file_name = 'madeup.dat'
- data = np.genfromtxt(file_name, usecols = (0, 1))
- n = len(data)
- convex_hull = open('convex_hull.dat', 'w')
- int hv_new = 0
- float theta_max = -2.0
- int theta = -1
- max_loc = i

4.2.1 Variable Documentation

4.2.1.1 convex_hull.convex_hull = open('convex_hull.dat', 'w')

Definition at line 18 of file convex_hull.py.

4.2.1.2 convex_hull.data = np.genfromtxt(file_name, usecols = (0, 1))

Definition at line 12 of file convex_hull.py.

4.2.1.3 string convex_hull.file_name = 'madeup.dat'

Definition at line 11 of file convex_hull.py.

4.2.1.4 convex_hull.hv_new = 0

Definition at line 22 of file convex_hull.py.

4.2.1.5 convex_hull.max_loc = i

Definition at line 35 of file convex_hull.py.

4.2.1.6 convex_hull.n = len(data)

Definition at line 15 of file convex_hull.py.

4.2.1.7 int convex_hull.theta = -1

Definition at line 31 of file convex_hull.py.

4.2.1.8 convex_hull.theta_max = -2.0

Definition at line 27 of file convex_hull.py.

Chapter 5

File Documentation

5.1 constants.f90 File Reference

Modules

- module constants

  Module contains definitions of useful constants.

Variables

- integer, parameter constants::dp =selected_real_kind(15, 300)

  Double-precision real kind.
- real(kind=dp), parameter constants::pi = 4.0_dp∗atan(1.0_dp)

  The circle constant, pi.
- real(kind=dp), parameter constants::tau = 2.0_dp∗pi

  2∗pi

5.2 convex_hull.py File Reference

Namespaces

- convex_hull

Variables

- string convex_hull.file_name = 'madeup.dat'
- convex_hull.data = np.genfromtxt(file_name, usecols = (0, 1))
- convex_hull.n = len(data)
- convex_hull.convex_hull = open('convex_hull.dat', 'w')
- int convex_hull.hv_new = 0
- float convex_hull.theta_max = -2.0
- int convex_hull.theta = -1
- convex_hull.max_loc = i

5.3 lattice.f90 File Reference

Functions/Subroutines

- program lattice

  Fortran 2003 program to generate an initial lattice.
- subroutine sheet_init (L, V, sheet)

  Initialises a thin sheet of randomly-arranged atoms.
- subroutine cube_init (L, cube, prop)

  Initialises a cube of randomly-arranged atoms.
- subroutine write_cube (cube, L, prop, fileno)

  Writes a generated cube to a CASTEP cell file.
- subroutine write_sheet (sheet, L, V, fileno)

  Writes a generated sheet to a CASTEP cell file.

5.3.1 Function/Subroutine Documentation

5.3.1.1 subroutine lattice::cube_init (integer, intent(in) *L,* integer, dimension(:,:,:), intent(inout), allocatable *cube,* real(kind=dp), intent(inout) *prop*)

Initialises a cube of randomly-arranged atoms.

Parameters

L	(input, integer) length of a side of the cube
cube	(output, integer array) the generated cube
prop	(input, real) proportion of metal ions that are calcium

Definition at line 92 of file lattice.f90.

5.3.1.2 program lattice ( )

Fortran 2003 program to generate an initial lattice.

Todo Produce a pair of sheets with a vacuum between them

This program creates a CASTEP cell file consisting of a random arrangement of Ca and Mg atoms in a crystal structure with oxygen, according to a user specification.

Definition at line 8 of file lattice.f90.

5.3.1.3 subroutine lattice::sheet_init (integer, intent(in) *L,* integer, intent(in) *V,* integer, dimension(:,:,:), intent(inout), allocatable *sheet*)

Initialises a thin sheet of randomly-arranged atoms.

Parameters

L	(input, integer) length of a side of the sheet
V	(input, integer) total height of the structure
sheet	(output, integer array) the generated sheet

Definition at line 54 of file lattice.f90.

5.3.1.4 subroutine lattice::write_cube (integer, dimension(:,:,:), intent(in), allocatable *cube,* integer, intent(in) *L,* real(kind=dp), intent(in) *prop,* integer, intent(in) *fileno*)

Writes a generated cube to a CASTEP cell file.

Parameters

cube	(input, integer array) the cube of atoms
L	(input, real) length of a side of the cube
prop	(input, real) proportion of metal ions that are calcium
fileno	(input, integer) the memory unit corresponding to the file to which to write

Definition at line 161 of file lattice.f90.

5.3.1.5 subroutine lattice::write_sheet (integer, dimension(:,:,:), intent(in), allocatable *sheet,* integer, intent(in) *L,* integer, intent(in) *V,* integer, intent(in) *fileno*)

Writes a generated sheet to a CASTEP cell file.

Parameters

sheet	(input, integer array) the sheet of atoms
L	(input, real) length of a side of the sheet
V	(input, integer) total height of the structure
fileno	(input, integer) the memory unit corresponding to the file to which to write

Definition at line 228 of file lattice.f90.