Alliance

1.0

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1 File Index	1
1.1 File List	1
2 File Documentation	3
2.1 array.c File Reference	3
2.1.1 Detailed Description	4
2.1.2 Function Documentation	4
2.1.2.1 get_flat_c()	4
2.1.2.2 get_flat_r()	4
2.1.2.3 get_flatIndexComplex3D()	5
2.1.2.4 getIndChi()	5
2.1.2.5 getIndChiBufEL_c()	5
2.1.2.6 getIndChiBufEL_r()	7
2.1.2.7 getIndChiBufEM_c()	7
2.1.2.8 getIndChiBufEM_r()	8
2.1.2.9 multiply_ar_c()	8
2.1.2.10 multiply_ar_r()	8
2.2 diagnostics.c File Reference	8
2.2.1 Detailed Description	9
2.2.2 Function Documentation	9
2.2.2.1 diag_compute()	9
2.2.2.2 diag_computeFreeEnergy()	10
2.2.2.3 diag_computeFreeEnergyFields()	10
2.2.2.4 diag_computeKSpectrum()	10
2.2.2.5 diag_computeMSpectrum()	11
2.2.2.6 diag_computeSpectra()	11
2.2.2.7 diag_getShells()	11
2.2.2.8 diag_initSpec()	12
2.3 distrib.c File Reference	12
2.3.1 Detailed Description	12
2.3.2 Function Documentation	12
2.3.2.1 distrib_enforceReality()	12
2.3.2.2 distrib_getG()	13
2.3.2.3 distrib_getH()	13
2.3.2.4 distrib_getXGrad()	13
2.3.2.5 distrib_getYGrad()	14
2.3.2.6 distrib_getZGrad()	14
2.3.2.7 distrib_setZeroNHalf()	14
2.4 equation.c File Reference	15
2.4.1 Detailed Description	15
2.4.2 Function Documentation	16
2.4.2.1 equation_getLinearTerm()	16

2.4.2.2 equation_getNonlinearElectromagnetic()	16		
2.4.2.3 equation_getNonlinearElectrostatic()	16		
2.4.2.4 equation_getNonlinearProduct()	17		
2.4.2.5 equation_getNonlinearTerm()	17		
2.5 fields.c File Reference	18		
2.5.1 Detailed Description	19		
2.5.2 Function Documentation	19		
2.5.2.1 fields_getA()	19		
2.5.2.2 fields_getAFromH()	20		
2.5.2.3 fields_getB()	20		
2.5.2.4 fields_getBFromH()	20		
2.5.2.5 fields_getChi()	21		
2.5.2.6 fields_getChiA()	21		
2.5.2.7 fields_getChiB()	21		
2.5.2.8 fields_getChiPhi()	21		
2.5.2.9 fields_getFields()	21		
2.5.2.10 fields_getFieldsFromH()	22		
2.5.2.11 fields_getGradX()	22		
2.5.2.12 fields_getGradY()	22		
2.5.2.13 fields_getPhi()	23		
2.5.2.14 fields_getPhiFromH()	23		
2.5.2.15 fields_init()	23		
2.5.2.16 fields_sendG()	23		
2.6 init.c File Reference	24		
2.6.1 Detailed Description	25		
2.6.2 Macro Definition Documentation	25		
2.6.2.1 RANK_IO	25		
2.6.3 Function Documentation	25		
2.6.3.1 fill_rand()	25		
2.6.3.2 fill_randM0()	25		
2.6.3.3 fill_randSingleKM()	26		
2.6.3.4 init_conditions()	26		
2.6.3.5 init_energySpec()	26		
2.6.3.6 init_initEnums()	27		
2.6.3.7 init_printParameters()	27		
2.6.3.8 init_start()	27		
2.7 parameters_io.c File Reference	27		
2.7.1 Detailed Description			
2.7.2 Function Documentation			
2.7.2.1 init_global_size()	28		
2.7.2.2 read_parameters()	28		
2.7.2.3 read_parametersFromFile()	28		

2.8 solver.c File Reference	29
2.8.1 Detailed Description	29
2.8.2 Function Documentation	29
2.8.2.1 solver_init()	29
2.8.2.2 solver_makeStep()	29
2.9 space_config.c File Reference	30
2.9.1 Detailed Description	31
2.9.2 Function Documentation	31
2.9.2.1 free_wavespace()	31
2.9.2.2 space_generateMSpace()	31
2.9.2.3 space_generateWaveSpace()	31
Index	33

Chapter 1

File Index

1.1 File List

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

array.c	
Array manipulation module	3
diagnostics.c	
Diagnostics module	8
distrib.c	
Gyrokinetic distribution function module	2
equation.c	
Equation module	5
fields.c	
Field computation and manipulation module	8
init.c	
Initialization module for alliance	24
parameters_io.c	
Reads inpuit parameters from parameter file provided by user	27
solver.c	
Numerical solver	29
space_config.c	
Space configuration module	30

2 File Index

Chapter 2

File Documentation

2.1 array.c File Reference

array manipulation module

```
#include "array.h"
#include "utils_fftw.h"
#include "space_config.h"
```

Macros

- #define CHI_EM 3
- #define CHI_EL 1
- #define FFT_OFFSET 2

Functions

```
    size_t get_flat_c (size_t is, size_t il, size_t im, size_t ix, size_t iy, size_t iz)
        returns flat index of the element of complex 6D array
    size_t getIndChiBufEM_c (size_t ix, size_t iy, size_t iz, size_t is, size_t ifield)
        returns flat index of an element of electromagnetic gyrokinetic potential in FOURIER SPACE
    size_t getIndChiBufEM_r (size_t ix, size_t iy, size_t iz, size_t is, size_t ifield)
        returns flat index of an element of electromagnetic gyrokinetic potential in POSITION SPACE
    size_t getIndChiBufEL_c (size_t ix, size_t iy, size_t iz, size_t is)
        returns returns flat index of an element of electrostatic gyrokinetic potential in FOURIER SPACE
    size_t getIndChiBufEL_r (size_t ix, size_t iy, size_t iz, size_t is)
        returns returns flat index of an element of electrostatic gyrokinetic potential in REAL SPACE
```

• size_t get_flat_r (size_t is, size_t il, size_t im, size_t ix, size_t iy, size_t iz)

returns flat index of the element of real 6D array

• size_t get_flatIndexComplex3D (size_t ix, size_t iy, size_t iz)

returns flat array of complex 3D array

- size_t getIndChi (size_t ix, size_t iy, size_t iz, size_t is)
- void multiply_ar_c (COMPLEX *ar1, COMPLEX *ar2, COMPLEX *ret)
- void multiply_ar_r (const double *ar1, const double *ar2, double *ret)

Variables

- struct array_size array_local_size
- struct array_size array_global_size
- · struct offset size array offset
- struct offset_size array_offset3D

2.1.1 Detailed Description

array manipulation module

contains functions which are supposed to make array manipulation simpler

2.1.2 Function Documentation

2.1.2.1 get_flat_c()

returns flat index of the element of complex 6D array

Parameters

is	species type	
il	Laguerre moment	
im	Hermite moment	
ix	kx index	
iy	ky index	
iz	kz index	

returns flattened index of a complex array from its 6D index. Flattened index then can be passed to distribution function 6D array to get a required element at position (is,il,im,ix,iy,iz).

2.1.2.2 get_flat_r()

```
size_t iy,
size_t iz )
```

returns flat index of the element of real 6D array

Parameters

is	species type	
il	Laguerre moment	
im	Hermite moment	
ix	x index	
iy	y index	
iz	z index	

returns flattened index of a real array from its 6D index. Flattened index then can be passed to distribution function 6D array to get a required element at position (is,il,im,ix,iy,iz).

2.1.2.3 get_flatIndexComplex3D()

returns flat array of complex 3D array

Parameters

ix	kx index
iy	ky index
iz	kz index

returns flattened index of a complex array from its 3D position index. Flattened index then can be passed to one of the fields ($\phi(\mathbf{k}), A_{||}(\mathbf{k}), B_{||}(\mathbf{k})$) 6D array to get a required element at position (ix,iy,iz).

2.1.2.4 getIndChi()

getIndChi(size_t ix,size_t iy, size_t iz, size_t is)

2.1.2.5 getIndChiBufEL_c()

```
size_t iy,
size_t iz,
size_t is)
```

returns returns flat index of an element of electrostatic gyrokinetic potential in FOURIER SPACE

Parameters

ix	kx index
iy	ky index
iz	kz index
is	particle species index

returns flattened index of a gyrokinetic potential $\chi^{\phi}(\mathbf{k})$ from its 4D index in FOURIER SPACE. flattened index is then can be used to access required value of the gyrokinetic potential at position (ix,iy,iz,is).

2.1.2.6 getIndChiBufEL_r()

returns returns flat index of an element of electrostatic gyrokinetic potential in REAL SPACE

Parameters

ix	x index
iy	y index
iz	z index
is	particle species index

returns flattened index of a gyrokinetic potential $\chi^{\phi}(\mathbf{r})$ from its 4D index in REAL SPACE. flattened index is then can be used to access required value of the gyrokinetic potential at position (ix,iy,iz,is).

2.1.2.7 getIndChiBufEM_c()

returns flat index of an element of electromagnetic gyrokinetic potential in FOURIER SPACE

Parameters

ix	kx index
iy	ky index
iz	kz index
is	particle species index
ifield	field type

returns flattened index of a gyrokinetic potential $\chi^{\phi,A,B}$ from its 4D index in FOURIER SPACE. flattened index is then can be used to access required value of the gyrokinetic potential at position (ix,iy,iz,is). Type of gyrokinetic potential is specified by ifield parameter. Use 0 is to access $\chi^{\phi}(\mathbf{k})$, 1 to access $\chi^{A}(\mathbf{k})$ and 2 to access $\chi^{B}(\mathbf{k})$.

2.1.2.8 getIndChiBufEM_r()

returns flat index of an element of electromagnetic gyrokinetic potential in POSITION SPACE

Parameters

ix	x index
iy	y index
iz	z index
is	particle species index
ifield	field type

returns flattened index of a gyrokinetic potential $\chi^{\phi,A,B}(\mathbf{k})$ from its 4D index in POSITION SPACE. flattened index is then can be used to access required value of the gyrokinetic potential at position (ix,iy,iz,is). Type of gyrokinetic potential is specified by ifield parameter. Use 0 is to access $\chi^{\phi}(\mathbf{r})$, 1 to access $\chi^{A}(\mathbf{r})$ and 2 to access $\chi^{B}(\mathbf{r})$.

2.1.2.9 multiply_ar_c()

multiply_ar_c(COMPLEX *ar1, COMPLEX *ar2, COMPLEX *ret)

2.1.2.10 multiply_ar_r()

multiply_ar_r(const double *ar1, const double *ar2, double *ret)

2.2 diagnostics.c File Reference

diagnostics module

```
#include "diagnostics.h"
#include "parameters_io.h"
```

Macros

- #define TO_ROOT 0
- #define BUFFER_SIZE 1

Functions

- void diag_computeSpectra (const COMPLEX *g, const COMPLEX *h, int timestep) general function to compute k or m spectra
- · void diag_initSpec ()

initialize spectra computation

void diag_computeFreeEnergy (COMPLEX *g, COMPLEX *h)

compute free energy

- void diag_computeKSpectrum (const COMPLEX *g, const COMPLEX *h, double *spec)
- void diag_computeMSpectrum (const COMPLEX *g, const COMPLEX *h, double *spec)

computes free energy spectra in m space

void diag_getShells ()

computes shells from parameters

• double diag_computeFreeEnergyFields (COMPLEX *g, COMPLEX *fields)

to be done later

• void diag_compute (COMPLEX *g, COMPLEX *h, int timestep)

computes all diagnostics

Variables

```
double * diag_kSpec = 0
```

used to store free energy k spectra

• double * diag_mSpec = 0

used to store free energy m spectra

• double * diag_shells = 0

used to store positions of k shells required to compute k spectra

double diag_freeEnergy

free energy

2.2.1 Detailed Description

diagnostics module

different diagnostic tools are gathered in this module

2.2.2 Function Documentation

2.2.2.1 diag_compute()

computes all diagnostics

Parameters

g	modified distribution function
h	distribution function
iter	current time step

2.2.2.2 diag_computeFreeEnergy()

compute free energy

Parameters

g	modified gyrokinetic distribution function
h	gyrokintic distribution function

computes free energy as $W=2.\Re(\sum_{k_x,k_y,k_z>0,m,l,s}g*\bar{h})$, taking into account reality condition.

2.2.2.3 diag_computeFreeEnergyFields()

to be done later

Parameters

g	
fields	computes free energy from the fields and distribution function.

2.2.2.4 diag_computeKSpectrum()

Parameters

g modified gyrokinetic distribution fund	
h	gyrokinetic distribution function
spec	spectra array

computes free energy k_{\perp} spectra $W(k_i^{shell})=\frac{1}{N}\sum_{k_{i-1}^{shell}<|\mathbf{k}_{\perp}|< k_i^{shell}}\sum_{k_z,l,m,s}g\bar{h}$ where N is a number of wave vectors between shells k_{i-1}^{shell} and k_i^{shell}

2.2.2.5 diag_computeMSpectrum()

computes free energy spectra in m space

Parameters

g	modified gyrokinetic distribution function
h	gyrokinetic distribution function
spec	spectra array

computes free energy m spectra as $W(m) = \sum_{k_x, k_y, k_z, l, s} g \bar{h}$

2.2.2.6 diag_computeSpectra()

general function to compute k or m spectra

Parameters

g	gyrokinetic distribution function
h	distribution function
timestep	current time step

function computes spectra at timestep as given in parameter file. k_{\perp} spectra is computed using diag_computeKSpectrum, and m spectra is computed using diag_computeMSpectrum

2.2.2.7 diag_getShells()

```
diag_getShells ( )
```

computes shells from parameters

computes positions of k_shells in between last_shell and first_shell as provided by user in parameter file. Position of i^{th} shell is computed as $k_i^{shell} = (last_shell - first_shell)/(k_shells) \cdot i$

2.2.2.8 diag_initSpec()

```
void diag_initSpec ( )
```

initialize spectra computation

Prepares free energy spectra computation. For spectra in k: allocates diag_kSpec array used to store k spectra. Allocates diag_shells array and fills it with shell positions k^{shells} , used for binning of wave vectors when computing k_{\perp} spectra. For spectra in m: allocates diag_mSpec array used to store m spectra. Called in init_start function

2.3 distrib.c File Reference

```
gyrokinetic distribution function module
```

```
#include "distrib.h"
```

Functions

- void distrib_getH (COMPLEX *h, const COMPLEX *g)
 computes h from g
- $\bullet \ \ \text{void} \ \ \frac{\text{distrib_getG}}{\text{getG}} \ \ (\text{COMPLEX} \ *g, \ \text{const} \ \ \text{COMPLEX} \ *h) \\$

computes g from h

• void distrib_getXGrad (const COMPLEX *in, COMPLEX *out)

Computes gradient in kx direction.

void distrib_getYGrad (const COMPLEX *in, COMPLEX *out)

Computes gradient in ky direction.

• void distrib_getZGrad (const COMPLEX *in, COMPLEX *out)

Computes gradient in kz direction.

- void distrib_enforceReality (COMPLEX *f)
 - enforces reality condition on distribution function array
- void distrib_setZeroNHalf (COMPLEX *f)

sets all Nk/2 modes to zero

2.3.1 Detailed Description

gyrokinetic distribution function module

everything required to perform different manipulations to distribution functions

2.3.2 Function Documentation

2.3.2.1 distrib_enforceReality()

```
void distrib_enforceReality ( {\tt COMPLEX} * f )
```

enforces reality condition on distribution function array

2.3 distrib.c File Reference

Parameters

f complex array for which reality condition will be forced.

Enforces reality condition f(k) = conj(f(-k)) in plane kz = 0. For a given kx, it first checks where modes -kx are located using the mpi_w -where mpi_w -where

```
where_neg = mpi_whereIsX[kxNeg * 2];
```

If -kx is stored on a different processor, MPI_VECTOR with a 4D data slice f(kx,kz = 0) is sent to this processor, to the buffer array:

```
mpi_sendVector(&f[ind6D], buffer, where_pos, where_neg);
```

if the data stored on the same processor, no vector is being sent. Reality condition is fulfilled in a loop over all other coordinates:

```
f[ind6D_neg] = conj(buffer[ind6D_pos]);
```

2.3.2.2 distrib_getG()

```
distrib_getG (  \label{eq:complex} {\tt COMPLEX} \, * \, g, \\ {\tt const} \, {\tt COMPLEX} \, * \, h \, )
```

computes g from h

Parameters

g	complex array to store g
h	complex array with h

computes modified gyrokinetic distribution function g from gyrokinetic distribution function h. Please note that before calling this function gyrokinetic potentials must be computed

2.3.2.3 distrib_getH()

computes h from g

Parameters

h	complex array to store h
g	complex array with g

computes gyrokinetic distribution function h from modified gyrokinetic distribution function g. Please note that before calling this function gyrokinetic potentials must be computed

2.3.2.4 distrib getXGrad()

```
void distrib_getXGrad (
```

```
const COMPLEX * in,
COMPLEX * out )
```

Computes gradient in kx direction.

Parameters

in	complex array. Distribution function of which gradient will be taken
out	complex array, where gradient is stored

Computes gradient in kx direction as following: grad(f) = i * kx * f

2.3.2.5 distrib_getYGrad()

Computes gradient in ky direction.

Parameters

in	complex array. Distribution function of which gradient will be taken
out	complex array, where gradient is stored

Computes gradient in ky direction as following: grad(f) = i * ky * f

2.3.2.6 distrib_getZGrad()

Computes gradient in kz direction.

Parameters

in	complex array. Distribution function of which gradient will be taken
out	complex array, where gradient is stored

Computes gradient in kz direction as following: grad(f) = i * kz * f

2.3.2.7 distrib_setZeroNHalf()

```
void distrib_setZeroNHalf ( {\tt COMPLEX} \ * \ f \ )
```

sets all Nk/2 modes to zero

Parameters

f complex array

sets Nkx/2, Nky/2 and Nz/2 modes of distribution function to zero. Due to reality condition, for kz yhe last mode should be set to zero.

2.4 equation.c File Reference

equation module

```
#include "equation.h"
#include "array.h"
#include "fields.h"
#include "distrib.h"
```

Macros

- #define CHI EM 3
- #define CHI EL 1
- #define CHI_PHI 0
- #define CHI A 1
- #define CHI_B 2

Functions

void equation_getLinearTerm (const COMPLEX *in, const COMPLEX *plus_boundary, const COMPLEX *minus_boundary, COMPLEX *out)

computes linear term

• void equation_getNonlinearElectromagnetic (double *in, double *chiAr, double *out, double sign)

returns nonlinear electromagnetic term

void equation_getNonlinearElectrostatic (double *in, double *chiAr, double *out, double sign)

returns nonlinear electrostatic term

• void equation_getNonlinearProduct (double *in, double *chiAr, double *out, double sign)

chooses between computing electrostatic or electromagnetic term

 $\bullet \ \ void\ equation_getNonlinearTerm\ (const\ COMPLEX\ *h,\ COMPLEX\ *out)$

computes nonlinear term

void equation_getRHS (const COMPLEX *in_g, COMPLEX *in_h, COMPLEX *out)

2.4.1 Detailed Description

equation module

module required to compute RHS of the equation.

2.4.2 Function Documentation

2.4.2.1 equation_getLinearTerm()

computes linear term

Parameters

in	complex array
out	complex array
plus_boundary	complex array
minus_boundary	complex array

computes linear term ${\tt out}$ from distribution function ${\tt in}$.

2.4.2.2 equation_getNonlinearElectromagnetic()

returns nonlinear electromagnetic term

Parameters

in	input double array
chiAr	input double array
out	output double array
sign	should be 1 or -1

performs multiplication between input 6D complex array in and gyrokinetic potential array chiAr, in such way that the structure of the product is the same as nonlinear term of drift-kinetic equations. Used by equation_getNonlinearProduct. sign is used to determine the sign of the resulting product. See equation_getNonlinearTerm for explanation.

2.4.2.3 equation_getNonlinearElectrostatic()

```
double * chiAr,
double * out,
double sign )
```

returns nonlinear electrostatic term

Parameters

in	input double array
chiAr	input double array
out	output double array
sign	should be 1 or -1

 $see\ equation_getNonlinear Electromagnetic\ for\ explanation$

2.4.2.4 equation_getNonlinearProduct()

chooses between computing electrostatic or electromagnetic term

Parameters

in	input double array
chiAr	input double array
out	output double array
sign	should be 1 or -1

 $\label{thm:thm:depending} \begin{subarray}{l} depending on flag \verb| systemType| provided by user in parameter file, chooses between equation_getNonlinearElectrostatic and equation_getNonlinearElectromagnetic \\ \end{subarray}$

2.4.2.5 equation_getNonlinearTerm()

computes nonlinear term

Parameters

h	input complex array
out	output complex array

function returns nonlinear term. First it takes y gradient of distribution function h, and x gradient of gyrokinetic

```
potentials chi, and transforms them to real space:
```

```
distrib_getYGrad(h, fftw_hBuf); fields_getGradX(fftw_chiBuf); fftw_c2r(); fftw_c2r_chi(); after that, it computes \frac{\partial h}{\partial y} \frac{\partial x}{\partial x} part of the poisson brackets: equation_getNonlinearProduct((double *)fftw_hBuf, (double *)fftw_chiBuf, buffer, 1.); with the result stored in buffer after that, it computes x gradient of h and y gradient of gyrokinetic potential chi, and transforms results to real space: distrib_getXGrad(h, fftw_hBuf); fields_getGradY(fftw_chiBuf); fftw_c2r();
```

and computes second part of the poisson brackets $-\frac{\partial h}{\partial x}\frac{\partial \chi}{\partial y}$ and adds the result to buffer. buffer is then transformed back to Fourier space, and dealiasing is performed.

2.5 fields.c File Reference

field computation and manipulation module

```
#include "fields.h"
```

fftw_c2r_chi();

Macros

- #define CHI_PHI 0
- #define CHI A 1
- #define CHI_B 2

Functions

```
· void fields_init ()
     intializes fields

    void fields_getA (const COMPLEX *g)

      compute A field

    void fields_getB (const COMPLEX *g0, const COMPLEX *g1)

      computes B potential

    void fields_getPhi (const COMPLEX *g0, const COMPLEX *g1)

      computes phi potential

    void fields getFields (COMPLEX *g00, COMPLEX *g10, COMPLEX *g01)

      wrapper to get all the fields simultaneously

    void fields_getChi ()

     computes gyrokinetic potentials chi
void fields_getChiPhi ()
     computes chiPhi gyrokinetic potential from phi potential

    void fields getChiB ()

     computes chiB gyrokinetic potential from B potential

    void fields getChiA ()
```

computes chiA gyrokinetic potential from A potential

2.5 fields.c File Reference 19

```
    void fields_sendG (COMPLEX *g)
```

sends moments of gyrokinetic distribution function which are required to compute fields

• void fields_getFieldsFromH (COMPLEX *h00, COMPLEX *h10, COMPLEX *h01)

wrapper to get all the fields simultaneously, computed from gyrokinetic distribution function

void fields_getAFromH (const COMPLEX *h)

compute A field

• void fields_getBFromH (const COMPLEX *h0, const COMPLEX *h1)

computes B potential

void fields_getPhiFromH (const COMPLEX *h)

computes phi potential

void fields getGradX (COMPLEX *out)

computes chi gradient in x direction

void fields_getGradY (COMPLEX *out)

computes chi gradient in y direction

Variables

- · struct fields_fields fields_fields
- struct fields_chi fields_chi
- double * A denom
- double * qnvTsJ
- double * I_B
- double * I_phi
- double * a_pot
- double * b_pot
- double * c_pot
- double * phiB_denom
- int * global_nm_index
- COMPLEX * g00
- COMPLEX * g10
- COMPLEX * g01

2.5.1 Detailed Description

field computation and manipulation module

Rerquired to compute $A_{||}(\mathbf{k}), B_{||}(\mathbf{k}), \phi(\mathbf{k})$ potentials, as well as gyrokinetic potentials $\chi_s^A(\mathbf{k}), \chi_s^B(\mathbf{k}), \chi_s^\phi(\mathbf{k})$

2.5.2 Function Documentation

2.5.2.1 fields_getA()

```
void fields_getA ( {\tt const~COMPLEX~*~g~)}
```

compute A field

Parameters

g

4D complex array (kx,ky,kz,s). Must be first Hermite and zeroth Laguerre moment of modified gyrokinetic distribution function g.

computes $A_{||}(\mathbf{k})$ potential from g_{s0}^1 (g parameter)

2.5.2.2 fields_getAFromH()

```
void fields_getAFromH ( {\tt const~COMPLEX~*~h~)}
```

compute A field

Parameters

h 4D complex array (kx,ky,kz,s). Must be first Hermite and zeroth Laguerre moment of gyrokinetic distribution function h.

computes $A_{||}(\mathbf{k})$ potential from h^1_{s0} (h parameter)

2.5.2.3 fields_getB()

```
void fields_getB ( {\tt const~COMPLEX~*~g0,} {\tt const~COMPLEX~*~g1~)}
```

computes B potential

Parameters

g0	4D complex array (kx,ky,kz,s). Zeroth Hermite and Laguerre moment of modified gyrokinetic distribution
	function.
g1	4D complex array (kx,ky,kz,s). Zeroth Hermite and first Laguerre moment of the modified distribution
	function.

Computes $B_{\perp}(\mathbf{k})$ from g_{s0}^0 (g0 parameter) and g_{s0}^1 (g1 parameter).

2.5.2.4 fields_getBFromH()

computes B potential

Parameters

h0	4D complex array (kx,ky,kz,s). Zeroth Hermite and Laguerre moment of gyrokinetic distribution function.
h1	4D complex array (kx,ky,kz,s). Zeroth Hermite and first Laguerre moment of distribution function.

2.5 fields.c File Reference

Computes $B_{\perp}({\bf k})$ from h_{s0}^0 (h0 parameter) and h_{s0}^1 (h1 parameter).

2.5.2.5 fields_getChi()

```
void fields_getChi ( )
```

computes gyrokinetic potentials chi

Wrapper for functions fields_getChiPhi, fields_getChiA, fields_getChiB

2.5.2.6 fields_getChiA()

```
void fields_getChiA ( )
```

computes chiA gyrokinetic potential from A potential

computes $chi_s^A(\mathbf{k})$

2.5.2.7 fields_getChiB()

```
void fields_getChiB ( )
```

computes chiB gyrokinetic potential from B potential

computes $\chi_s^B(\mathbf{k})$

2.5.2.8 fields_getChiPhi()

```
fields_getChiPhi ( )
```

computes chiPhi gyrokinetic potential from phi potential

computes $chi_s^{\phi}(\mathbf{k})$

2.5.2.9 fields_getFields()

wrapper to get all the fields simultaneously

Parameters

g00	4D complex array (kx,ky,kz,s). Zeroth Hermite and Laguerre moment of modified gyrokinetic distribution function.
g10	4D complex array (kx,ky,kz,s). Must be first Hermite and zeroth Laguerre moment of modified gyrokinetic distribution function g.
Generate	4D complex array (kx,ky,kz,s). Zeroth Hermite and first Laguerre moment of the modified distribution function.

Wrapper for functions fields_getPhi, fields_getB, fields_getA.

2.5.2.10 fields getFieldsFromH()

wrapper to get all the fields simultaneously, computed from gyrokinetic distribution function

Parameters

h00	4D complex array (kx,ky,kz,s). Zeroth Hermite and Laguerre moment of gyrokinetic distribution function.
h10	4D complex array (kx,ky,kz,s). Must be first Hermite and zeroth Laguerre moment of gyrokinetic distribution
	function h.
h01	4D complex array (kx,ky,kz,s). Zeroth Hermite and first Laguerre moment of the distribution function.

Wrapper for functions fields_getPhiFromH, #fields_get_BFromH, fields_getAFromH.

2.5.2.11 fields getGradX()

computes chi gradient in x direction

Parameters

```
out output complex array of size (kx,ky,kz,s,Nfields).
```

computes gradient in x direction for chi potentials. Nfields<\tt> can be 1 or 3, and chosen automatically at start of the simulation depending on the simulation type (electrostatic or electromagnetic)

2.5.2.12 fields_getGradY()

computes chi gradient in y direction

Parameters

```
out output complex array of size (kx,ky,kz,s,Nfields).
```

computes gradient in y direction for chi potentials. Nfields < tt> can be 1 or 3, and chosen automatically at start of the simulation depending on the simulation type

2.5 fields.c File Reference 23

```
(electrostatic or electromagnetic)
```

2.5.2.13 fields getPhi()

computes phi potential

Parameters

g0	4D complex array (kx,ky,kz,s). Zeroth Hermite and Laguerre moment of modified gyrokinetic distribution
	function.
g1	4D complex array (kx,ky,kz,s). Zeroth Hermite and first Laguerre moment of the modified distribution
	function.

Computes $\phi(\mathbf{k})$ from g_{s0}^0 (g0 parameter) and g_{s0}^1 (g1 parameter).

2.5.2.14 fields_getPhiFromH()

```
void fields_getPhiFromH ( {\tt const~COMPLEX~*~h~)}
```

computes phi potential

Parameters

h | 4D complex array (kx,ky,kz,s). Zeroth Hermite and Laguerre moment of gyrokinetic distribution function.

Computes $\phi(\mathbf{k})$ from h_{s0}^0 (h parameter)

2.5.2.15 fields_init()

```
void fields_init ( )
```

intializes fields

pre-computes some comstants required to compute fields. Called in init_start function

2.5.2.16 fields_sendG()

```
fields_sendG ( {\tt COMPLEX} \ * \ g \ )
```

sends moments of gyrokinetic distribution function which are required to compute fields

Parameters

g complex array. Modified or non-modified gyrokinetic distribution function

```
sends g_{s0}^{1}(\mathbf{k})f, g_{s1}^{0}(\mathbf{k})f, g_{s0}^{0}(\mathbf{k})f
```

to all processes to compute potentials locally.

2.6 init.c File Reference

initialization module for alliance.

```
#include "init.h"
#include "distrib.h"
```

Macros

• #define RANK_IO 0

Functions

```
• void init_start (char *filename)
```

initialization of ALLIANCE

void init_printParameters ()

parameter output

• void init_initEnums ()

enumerator initialization

void fill_rand (COMPLEX *ar1)

fills the inital conditions randomly

• void fill_randM0 (COMPLEX *ar1)

fill zeroth Hermite moment with random values

void fill_randSingleKM (COMPLEX *ar1)

fill single chosen wavevector and Hermite moment

void init_conditions (COMPLEX *data)

distribution function initialization

• double init_energySpec (double k, double m, double amp, double disp)

returns energy spectrum

Variables

- · enum adiabatic kinetic
- enum electromagnetic systemType
- · enum initial initialConditions

2.6 init.c File Reference 25

2.6.1 Detailed Description

initialization module for alliance.

all the inititalization routines are here.

2.6.2 Macro Definition Documentation

2.6.2.1 RANK_IO

```
#define RANK_IO 0
```

defines rank of the processor used to output information to console

2.6.3 Function Documentation

2.6.3.1 fill_rand()

```
void fill_rand ( {\tt COMPLEX} \ * \ data \ )
```

fills the inital conditions randomly

Parameters

data

complex 6d array to fill initializes distribution with spectrum defined in init_energySpec This function is supposed to be used in-module only and should not be used elsewhere outside init.c file.

2.6.3.2 fill_randM0()

```
void fill_randM0 ( {\tt COMPLEX * data })
```

fill zeroth Hermite moment with random values

Parameters

data | complex 6D array to fill

fills 0-th Hermite moment of a distribution function ar1 with random values This function is supposed to be used in-module only and should not be used elsewhere outside init.c file.

2.6.3.3 fill_randSingleKM()

fill single chosen wavevector and Hermite moment

Parameters

```
data complex 6D array
```

initializes single wavevector and Hermite moment of a distribution function with random variable. This function is only for in-module use and should not be used elsewhere outside init.c file.

2.6.3.4 init_conditions()

```
void init_conditions ( {\tt COMPLEX} \ * \ {\tt data} \ )
```

distribution function initialization

Parameters

data	complex 6D array
------	------------------

initializes distribution function with chosen method (see fill_rand, fill_randM0, fill_randSingleKM)

2.6.3.5 init_energySpec()

returns energy spectrum

Parameters

k	a wavenumber at which spectrum is computed
m	Hermite moment at which amplitude is computed
amp	amplitude of the spectrum
disp	dispersion of the spectrum

computes spectrum of form $A \cdot k^2 exp(-2k^2/\sigma^2)$, where $\sigma = disp$, and A = amp This function is supposed to be

used in-module only and should not be used elsewhere outside init.c file.

2.6.3.6 init_initEnums()

```
void init_initEnums ( )
```

enumerator initialization

initializes enumerators, which are then used to define if system is adiabatic or kinetic, electromagnetic or electrostatic, and type of initial conditions

2.6.3.7 init_printParameters()

```
void init_printParameters ( )
```

parameter output

prints parameters of the simulation

2.6.3.8 init_start()

initialization of ALLIANCE

Parameters

lename specifies parameter filena	те
-----------------------------------	----

initializes all the modules required for ALLIANCE to work.

2.7 parameters_io.c File Reference

reads inpuit parameters from parameter file provided by user

```
#include "parameters_io.h"
#include "utils_fftw.h"
```

Macros

• #define VERBOSE 0

Functions

```
• void init_global_size ()
```

initializes global size of the 6D array

• void read_parameters (char *filename)

reads parameters from user parameter file.

• void read_parametersFromFile (char *filename)

Variables

• struct system_param parameters

2.7.1 Detailed Description

reads inpuit parameters from parameter file provided by user

2.7.2 Function Documentation

2.7.2.1 init_global_size()

```
void init_global_size ( )
```

initializes global size of the 6D array

 $initializes \verb| array_local_size| | \textit{structure with global simulation size}.$

2.7.2.2 read_parameters()

reads parameters from user parameter file.

Reads parameters from user parameter file. All the parameters are stored in the parameters structure

2.7.2.3 read_parametersFromFile()

read_parametersFromFile(char *filename):

2.8 solver.c File Reference

2.8 solver.c File Reference

```
numerical solver
```

```
#include "solver.h"
```

Macros

- #define **SOLVERTYPE** RK4
- #define IORANK 0

Functions

```
    void solver_init ()
        initializes solver
    void solver_makeStep (COMPLEX **g, COMPLEX *h)
        iterate solver forward
```

Variables

- enum solverType solverType
- struct solver solver
- struct rk4 rk4

2.8.1 Detailed Description

numerical solver

2.8.2 Function Documentation

2.8.2.1 solver_init()

```
void solver_init ( )
initializes solver
initializes solver with the solverType.
```

2.8.2.2 solver_makeStep()

```
void solver_makeStep ( {\tt COMPLEX~**~g,} {\tt COMPLEX~*~h~)}
```

iterate solver forward

Parameters

g	address of the 6D complex array. Modified gyrokinetic distribution function	
h	6D complex array. Gyrokinetic distribution function	

solves one simulation time step

2.9 space_config.c File Reference

space configuration module

```
#include "space_config.h"
#include <complex.h>
```

Macros

• #define MINUS_I -1.j

Functions

```
    void space_init ()
        initializes wave space. Called in init_start() function.
    void space_generateWaveSpace ()
        generates wave space
```

• void space_generateMSpace () generates Hermite space

• void free_wavespace ()

deallocates all the arrays

Variables

```
• double space_Lx = 100
```

- double space_Ly = 100
- double space_Lz = 100
- double * space_kx
- double * space_ky
- double * space_kz
- double * space_kPerp
- double * space_kPerp2
- double * space_kSq
- double * space_sqrtM
- size t * space_globalMIndex
- COMPLEX * space_iKx
- COMPLEX * space_iKy
- COMPLEX * space_iKz

2.9.1 Detailed Description

space configuration module

creates k and m spaces

2.9.2 Function Documentation

2.9.2.1 free_wavespace()

```
free_wavespace ( )
```

deallocates all the arrays

to be added

2.9.2.2 space_generateMSpace()

```
space_generateMSpace ( )
```

generates Hermite space

to be added

2.9.2.3 space generateWaveSpace()

```
void space_generateWaveSpace ( )
```

generates wave space

generates wave number arrays space_kx, space_ky, space_kz of lengths nkx,nky,nkz for a numerical box of size [lx, ly, lz] in kx,ky,kz directions as following:

```
[0, pi / lx, 2 pi / lx, ..., (n / 2 + 1) pi / lx, - (n / 2) pi / lx, ..., - pi / lx] generates arrays space_iKx, space_iKy, space_iKz, of lengths nkx,nky,nkz. These arrays are later used to compute gradients by fields_getGradX, fields_getGradY, distrib_getXGrad, distrib_getYGrad, distrib_getZGrad.
```

Index

```
array.c, 3
                                                              distrib.c, 13
                                                         distrib getXGrad
     get_flat_c, 4
     get flat r, 4
                                                              distrib.c, 13
     get flatIndexComplex3D, 5
                                                         distrib getYGrad
     getIndChi, 5
                                                              distrib.c, 14
     getIndChiBufEL_c, 5
                                                         distrib_getZGrad
     getIndChiBufEL r, 7
                                                              distrib.c, 14
     getIndChiBufEM_c, 7
                                                         distrib setZeroNHalf
                                                              distrib.c, 14
     getIndChiBufEM_r, 8
     multiply ar c, 8
                                                         equation.c, 15
     multiply ar r, 8
                                                              equation getLinearTerm, 16
diag compute
                                                              equation getNonlinearElectromagnetic, 16
     diagnostics.c. 9
                                                              equation getNonlinearElectrostatic, 16
diag computeFreeEnergy
                                                              equation getNonlinearProduct, 17
     diagnostics.c, 10
                                                              equation getNonlinearTerm, 17
diag computeFreeEnergyFields
                                                         equation getLinearTerm
     diagnostics.c. 10
                                                              equation.c, 16
diag_computeKSpectrum
                                                         equation_getNonlinearElectromagnetic
     diagnostics.c, 10
                                                              equation.c, 16
diag_computeMSpectrum
                                                         equation_getNonlinearElectrostatic
     diagnostics.c, 11
                                                              equation.c, 16
                                                         equation_getNonlinearProduct
diag_computeSpectra
     diagnostics.c, 11
                                                              equation.c, 17
diag getShells
                                                         equation getNonlinearTerm
     diagnostics.c, 11
                                                              equation.c, 17
diag_initSpec
                                                         fields.c, 18
     diagnostics.c, 11
                                                              fields getA, 19
diagnostics.c, 8
                                                              fields_getAFromH, 20
    diag compute, 9
                                                              fields_getB, 20
     diag_computeFreeEnergy, 10
                                                              fields getBFromH, 20
     diag_computeFreeEnergyFields, 10
                                                              fields_getChi, 21
     diag computeKSpectrum, 10
                                                              fields_getChiA, 21
     diag computeMSpectrum, 11
                                                              fields getChiB, 21
     diag computeSpectra, 11
                                                              fields getChiPhi, 21
     diag getShells, 11
                                                              fields getFields, 21
     diag initSpec, 11
                                                              fields getFieldsFromH, 22
distrib.c, 12
                                                              fields getGradX, 22
    distrib_enforceReality, 12
                                                              fields getGradY, 22
     distrib_getG, 13
                                                              fields_getPhi, 23
     distrib getH, 13
     distrib_getXGrad, 13
                                                              fields_getPhiFromH, 23
                                                              fields init, 23
     distrib_getYGrad, 14
                                                              fields sendG, 23
     distrib getZGrad, 14
                                                         fields_getA
     distrib setZeroNHalf, 14
                                                              fields.c, 19
distrib_enforceReality
                                                         fields getAFromH
     distrib.c, 12
                                                              fields.c. 20
distrib getG
                                                         fields getB
     distrib.c, 13
                                                              fields.c, 20
distrib_getH
```

34 INDEX

fields_getBFromH	init_printParameters, 27
fields.c, 20	init_start, 27
fields_getChi	RANK_IO, 25
fields.c, 21	init_conditions
fields_getChiA	init.c, 26
fields.c, 21	init_energySpec
fields_getChiB	init.c, 26
fields.c, 21	init_global_size
fields getChiPhi	parameters io.c, 28
fields.c, 21	init initEnums
fields getFields	init.c, 27
fields.c, 21	init printParameters
fields_getFieldsFromH	init.c, 27
	, and the second
fields.c, 22	init_start
fields_getGradX	init.c, 27
fields.c, 22	multiply or o
fields_getGradY	multiply_ar_c
fields.c, 22	array.c, 8
fields_getPhi	multiply_ar_r
fields.c, 23	array.c, 8
fields_getPhiFromH	07
fields.c, 23	parameters_io.c, 27
fields_init	init_global_size, 28
fields.c, 23	read_parameters, 28
fields_sendG	read_parametersFromFile, 28
fields.c, 23	BANK 18
fill rand	RANK_IO
init.c, 25	init.c, 25
fill randM0	read_parameters
init.c, 25	parameters_io.c, 28
fill_randSingleKM	read_parametersFromFile
	parameters_io.c, 28
init.c, 26	
free_wavespace	solver.c, 29
space_config.c, 31	solver_init, 29
ant flat a	solver_makeStep, 29
get_flat_c	solver_init
array.c, 4	solver.c, 29
get_flat_r	solver_makeStep
array.c, 4	solver.c, 29
get_flatIndexComplex3D	space_config.c, 30
array.c, 5	free_wavespace, 31
getIndChi	space_generateMSpace, 31
array.c, 5	space generateWaveSpace, 31
getIndChiBufEL_c	space_generateMSpace
array.c, 5	space_generatewopace space config.c, 31
getIndChiBufEL_r	. – •
array.c, 7	space_generateWaveSpace
getIndChiBufEM_c	space_config.c, 31
array.c, 7	
getIndChiBufEM_r	
array.c, 8	
-y -) -	
init.c, 24	
fill_rand, 25	
fill_randM0, 25	
fill_randSingleKM, 26	
init_conditions, 26	
init_energySpec, 26	
init initEnums, 27	
IIII,	