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# NHDSPy Documentation

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nhds.py is a wrapper for the NHDS dispersion relation solver. For the original NHDS code see <https://github.com/danielver02/NHDS>.

If you use this software to produce data for publication, please cite the NHDS paper: <http://iopscience.iop.org/article/10.3847/2515-5172/aabfe3>



# CHAPTER 1

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## Installing

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To install nhdsby, run

```
pip install nhdsby
```





## CHAPTER 2

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### Running

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NHDS is written in fortran, so requires a working fortran compiler. When nhds.py is first run it attempts to compile NHDS; if it is not successfully it gracefully exits and prints the fortran compile error.



## CHAPTER 3

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### Example

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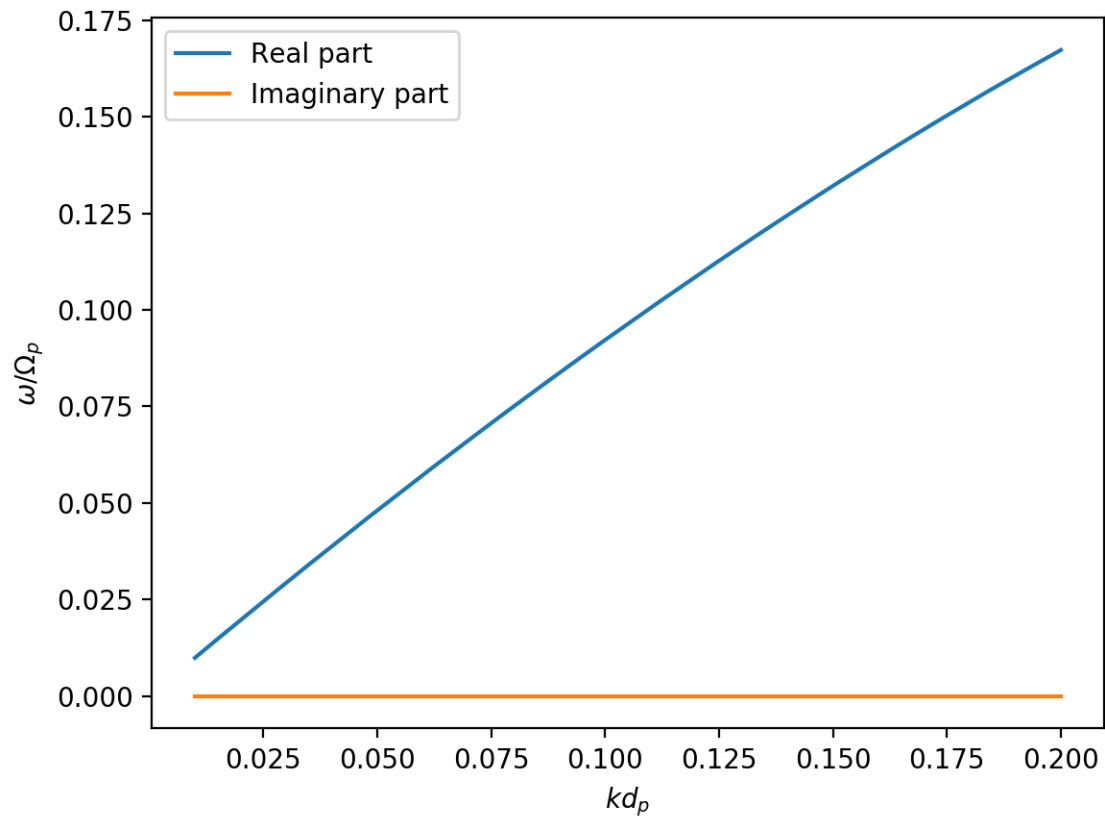
```
import nhdspy
import matplotlib.pyplot as plt

electrons = nhdspy.Species(-1, 1 / 1836, 1, 0, 1, 1)
protons = nhdspy.Species(1, 1, 1, 0, 1, 1)

va = 1e-4
theta_kB = 0.001
omega_0 = 0.009 - 0.0001 * 1j
kzmin = 0.01
kzmax = 0.2
input = nhdspy.InputParams([electrons, protons], omega_0, theta_kB,
                           va, kzmin, kzmax)
output = nhdspy.run(input)

fig, ax = plt.subplots()
ax.plot(output.kz, output.omega_real, label='Real part')
ax.plot(output.kz, output.omega_imag, label='Imaginary part')
ax.set_xlabel('$kd_{p}$')
ax.set_ylabel(r'$\omega / \Omega_{p}$')
ax.legend()

plt.show()
```



## 3.1 Code reference

### 3.1.1 nhds py Package

#### Functions

<code>format_input_file(input)</code>	Function to create input file.
<code>run(input)</code>	Run the dispersion solver for a given input.

#### format\_input\_file

`nhds py.format_input_file(input)`  
Function to create input file.

##### Parameters

**input** [InputParams]

## run

`nhdsipy.run(input)`

Run the dispersion solver for a given input.

### Parameters

**input** [InputParams]

## Classes

<code>InputParams(species, omega_guess, ...[, ...])</code>	Input parameters for calculating a dispersion relation.
<code>Result(input, run_output)</code>	Result of running the dispersion solver.
<code>Species(q, m, n, v_d, t_anis, beta_par)</code>	A single bi-Maxwellian species.

## InputParams

```
class nhdsipy.InputParams (species, omega_guess, propagation_angle, va, kzmin, kzmax,
                             kzsteps=200, numiter=1000, det_D_threshold=1e-16, n_bessel=1000,
                             bessel_zero=1e-50, vxsteps=100, vysteps=100, vzsteps=100)
```

Bases: object

Input parameters for calculating a dispersion relation.

### Parameters

**species** [list of *Species*] List of the particle species for which to calculate the dispersion.  
**Note** that the order of the species matters, as some of the parameters below are defined relative to the first species in the *species* list.

**omega\_guess** [complex] Initial guess for the frequency, normalised to the gyro frequency of the first species.

**propagation\_angle** [float] Propagation angle to calculate dispersion relation for. Defined with respect to the magnetic field.

**va** [float] Ratio of the Alfvén speed to the speed of light. The number density in the Alfvén speed is taken from the first species.

**kzmin** [float] Start of kz range.

**kzmax** [float] End of kz range.

### Other Parameters

**kzsteps** [int, optional] Number of steps in kz to calculate dispersion relation at. Points are linearly spaced between *kzmin* and *kzmax*.

**numiter** [int, optional] Maximum number of iterations in the Newton method

**det\_D\_threshold** [float, optional] Threshold for the determinant of the dispersion tensor. If  $\det D \leq \det\_D\_threshold$ , the Newton iteration will be stopped.

**n\_bessel** [int, optional] Maximum of sum in Bessel functions (both regular and modified). Can be very low (e.g., 3) for quasi-parallel propagation.

**bessel\_zero** [float, optional] If  $I_n$  is less than this value, higher  $n$  are neglected.

**vxsteps, vysteps, vzsteps** [int, optional] Steps in the x,y,z directions.

### Attributes

*nspecies* Number of species.

*total\_charge* Total charge density.

*total\_current* Total current density. Result is normalised to the first species' charge, number density, and the Alfvén speed.

### Attributes Summary

<i>nspecies</i>	Number of species.
<i>total_charge</i>	Total charge density.
<i>total_current</i>	Total current density. Result is normalised to the first species' charge, number density, and the Alfvén speed.

### Attributes Documentation

#### **nspecies**

Number of species.

#### **total\_charge**

Total charge density. Result is normalised to the first species' charge and number density.

#### **total\_current**

Total current density. Result is normalised to the first species' charge, number density, and the Alfvén speed.

## Result

**class** `nhdspy.Result` (*input, run\_output*)

Bases: `object`

Result of running the dispersion solver.

#### Attributes

***EyEx*** The electric field component ratio  $E_y/E_x$ .

***EzEx*** The electric field component ratio  $E_z/E_x$ .

***kz*** Wave vector normalised to the proton inertial length ( $kd_p$ ).

***omega\_imag*** Imaginary part of frequency normalised to the proton gyro-frequency ( $\gamma/\Omega_p$ ).

***omega\_real*** Real part of frequency normalised to the proton gyro-frequency ( $\omega/\Omega_p$ ).

### Attributes Summary

<i>EyEx</i>	The electric field component ratio $E_y/E_x$ .
<i>EzEx</i>	The electric field component ratio $E_z/E_x$ .
<i>kz</i>	Wave vector normalised to the proton inertial length ( $kd_p$ ).

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<i>omega_imag</i>	Imaginary part of frequency normalised to the proton gyro-frequency ( $\gamma/\Omega_p$ ).
<i>omega_real</i>	Real part of frequency normalised to the proton gyro-frequency ( $\omega/\Omega_p$ ).

## Attributes Documentation

### **EyEx**

The electric field component ratio  $E_y/E_x$ . Returned as complex numbers.

### **EzEx**

The electric field component ratio  $E_z/E_x$ . Returned as complex numbers.

### **kz**

Wave vector normalised to the proton inertial length ( $kd_p$ ).

### **omega\_imag**

Imaginary part of frequency normalised to the proton gyro-frequency ( $\gamma/\Omega_p$ ).

### **omega\_real**

Real part of frequency normalised to the proton gyro-frequency ( $\omega/\Omega_p$ ).

## Species

**class** `nhds.py.Species` (*q, m, n, v\_d, t\_anis, beta\_par*)

Bases: `object`

A single bi-Maxwellian species.

### Parameters

**q** [float] Particle charge, in units of proton charge. e.g. for a proton enter 1, for an electron enter -1.

**m** [float] Particle mass, in units of proton mass. e.g. for a proton enter 1.

**n** [float] Number density, relative to the number density used to define the Alfvén speed.

**v\_d** [float] Drift speed as a fraction of the Alfvén speed.

**t\_anis** [float] Temperature anisotropy (perpendicular temperature divided by parallel temperature).

**beta\_par** [float] Parallel beta (thermal pressure divided by magnetic pressure).





## CHAPTER 4

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### Indices and tables

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