

OpenMX tutorials

Magnons in Fe using OpenMX

Computational Materials Physics Bootcamp

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Fe using OpenMX: Ground-State, Band Structure, DOS

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1. Create a subdirectory inside your main directory using 'mkdir' command:

```
mkdir OpenMX-Fe
```

2. Copy all data from ../alikefay/OpenMX-Fe directory to your own subdirectory using cp command:

```
cp ../alikefay/OpenMX/* ./OpenMX-Fe
```

3. Go to your subdirectory using 'cd' command:

```
cd ./OpenMX-Fe
```

4. Open the Fe.in file and try to understand the keywords.

5. Now, if you feel ready for the run, it's time to submit the job. We need first to activate the module using:

```
module unload openmpi/4.1.4-intel-2023.0.0
```

```
module unload intel/2023.0.0
```

```
module load intel/2020.3
```

```
module load openmx
```

Fe using OpenMX: Ground-State, Band Structure, DOS

Fe.in

```
# File Name
System.CurrentDirectory      ./                # default=./
System.Name                  Fe
DATA.PATH                    /home/opt/pkg/openmx/3.9.9-intel-2020.3/DFT_DATA19
level.of.stdout              1                  # default=1 (1-3)
level.of.fileout              0                  # default=1 (0-2)

# Definition of Atomic Species
Species.Number               1
<Definition.of.Atomic.Species
| Fe Fe6.0S-s3p2d1 Fe_PBE19H
Definition.of.Atomic.Species>

# Atoms
Atoms.Number                  1
Atoms.SpeciesAndCoordinates.Unit Ang            # Ang|AU
<Atoms.SpeciesAndCoordinates
| 1 Fe 0.0000 0.0000 0.0000 9.0 7.0
Atoms.SpeciesAndCoordinates>

Atoms.UnitVectors.Unit       Ang                # Ang|AU
<Atoms.UnitVectors
| -1.43300 1.43300 1.43300
| 1.43300 -1.43300 1.43300
| 1.43300 1.43300 -1.43300
Atoms.UnitVectors>
```

Fe using OpenMX: Ground-State, Band Structure, DOS

Fe.in

```
# SCF or Electronic System
scf.XcType          GGA-PBE          # LDA|LSDA-CA|LSDA-PW|GGA-PBE
scf.SpinPolarization On              # On|Off|NC
scf.SpinOrbit.Coupling Off            # On|Off, default=off
HS.fileout          Off              # on|off, default=off
scf.ElectronicTemperature 300.0      # default=300 (K)
scf.energycutoff    600.0            # default=150 (Ry)
scf.maxIter         800              # default=40
scf.EigenvalueSolver band            # Recursion|Cluster|Band
scf.Kgrid            9 9 9           # means nk1xnk2xnk3
scf.Mixing.Type      rmm-diisk        # Simple|Rmm-Diis|Gr-Pulay
scf.Init.Mixing.Weight 0.300          # default=0.30
scf.Min.Mixing.Weight 0.001           # default=0.001
scf.Max.Mixing.Weight 0.400           # default=0.40
scf.Mixing.History    5               # default=5
scf.Mixing.StartPulay 6               # default=6
scf.criterion         1.0e-7          # default=1.0e-6 (Hartree)
Dos.fileout          on               # DOS calculation
Dos.Erange           -25.0 25.0       # DOS energy range (0=chemical potential)
Dos.Kgrid             27 27 27        # DOS k-point discretization

Band.dispersion      on               # on|off, default=off
Band.Nkpath          5
<Band.kpath
  40  0.0000 0.0000 0.0000 0.5000 -0.500 0.5000  G H
  50  0.5000 -0.500 0.5000 0.0000 0.0000 0.5000  H N
  30  0.0000 0.0000 0.5000 0.0000 0.0000 0.0000  N G
  30  0.0000 0.0000 0.0000 0.2500 0.2500 0.2500  G P
  40  0.2500 0.2500 0.2500 0.5000 -0.500 0.5000  P H
Band.kpath>

# MD or Geometry Optimization
MD.Type             NOMD              # NOMD|OptC1|OptC2|OptC3|OptC4|OptC5|OptC6|OptC7|RFC5|RFC6|RFC7
MD.maxIter          1                  # default=1
MD.Opt.criterion     3.0e-4            # default=0.0003 (Hartree/Bohr)
```

Fe using OpenMX: Ground-State, Band Structure, DOS

6. Now, you can submit the job using the following command,

```
nohup mpirun -np 5 openmx Fe.in > Fe.out -f openmp &
```

7. The Fe.out is the job log and you can see the progress of the scf loops by opening it.
8. When the job is finished open the Fe.out and find the magnetic moment of Fe atom and compare it with the experimental value of $2.22 \mu_B$. What is the ground state energy? How many scf loops are performed?

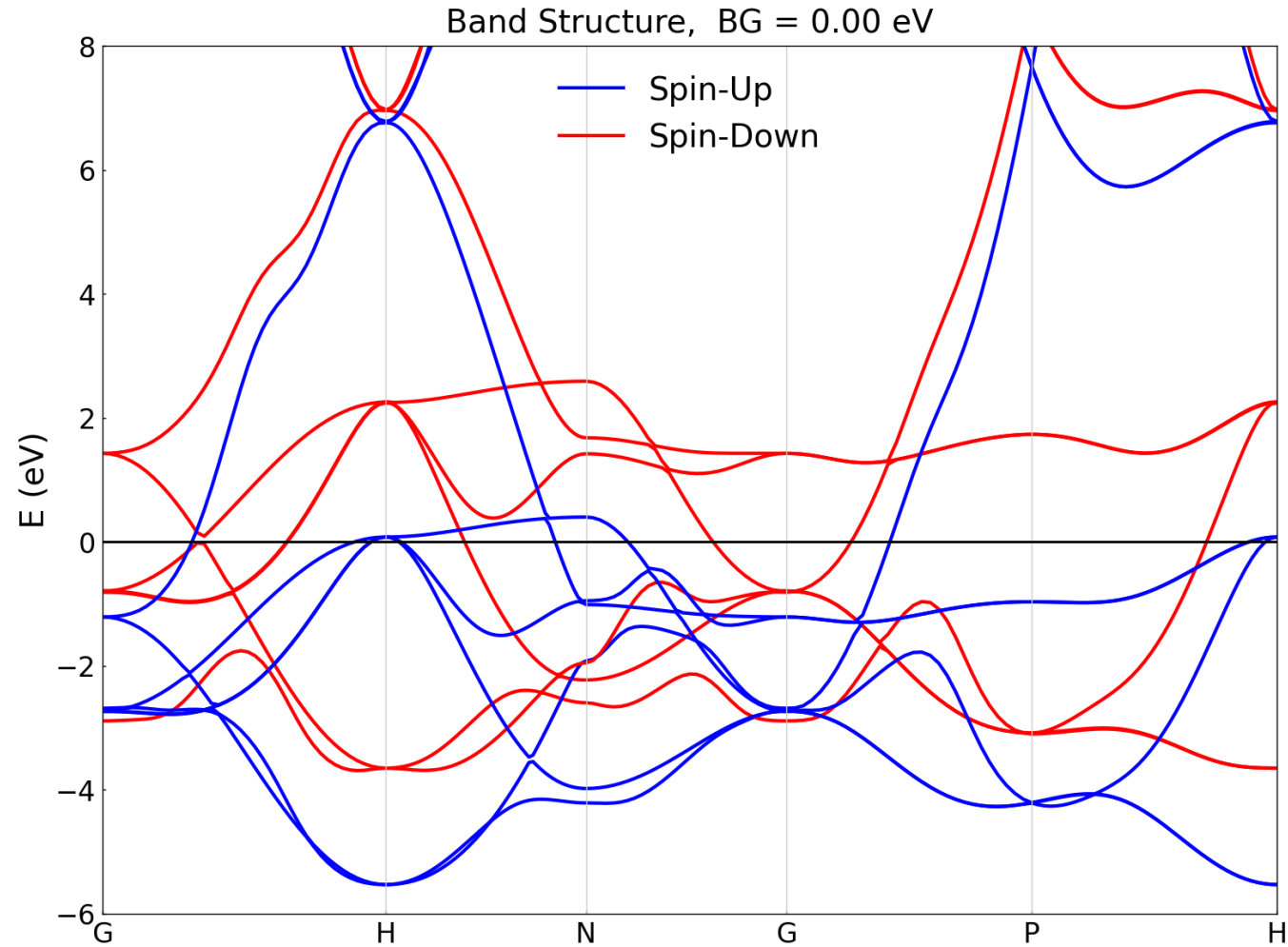
9. Use the following command to prepare the band structure data for plotting,

```
bandgnu13 Fe.Band
```

10. Next use the python script in your directory to plot the band structure using the command,

```
Python3 Plot_band.py
```

Fe using OpenMX: Ground-State, Band Structure, DOS



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11. Now, we are ready to calculate the DOS. The keyword setting in the code are:

```
Dos.fileout      on          # DOS calculation
Dos.Erange      -25.0 25.0   # DOS energy range (0=chemical potential)
Dos.Kgrid       27 27 27    # DOS k-point discretization
```

Since the keyword is already set in the input files, you just need to do the post processing. Use the following command to calculate the total DOS (PDOS)

```
DosMain Fe.Dos.val Fe.Dos.vec
```

Then, you are interactively asked from the program as follow:

```
DosMain Fe.Dos.val Fe.Dos.vec
Max of Spe_Total_CNO = 14
1 1 1 101 102 103 101 102 103 201 202 203 204 205
<Fe.Dos.val>
<Fe>
Which method do you use?, Tetrahedron(1), Gaussian Broadening(2)
1
Do you want Dos(1) or PDos(2)?
2

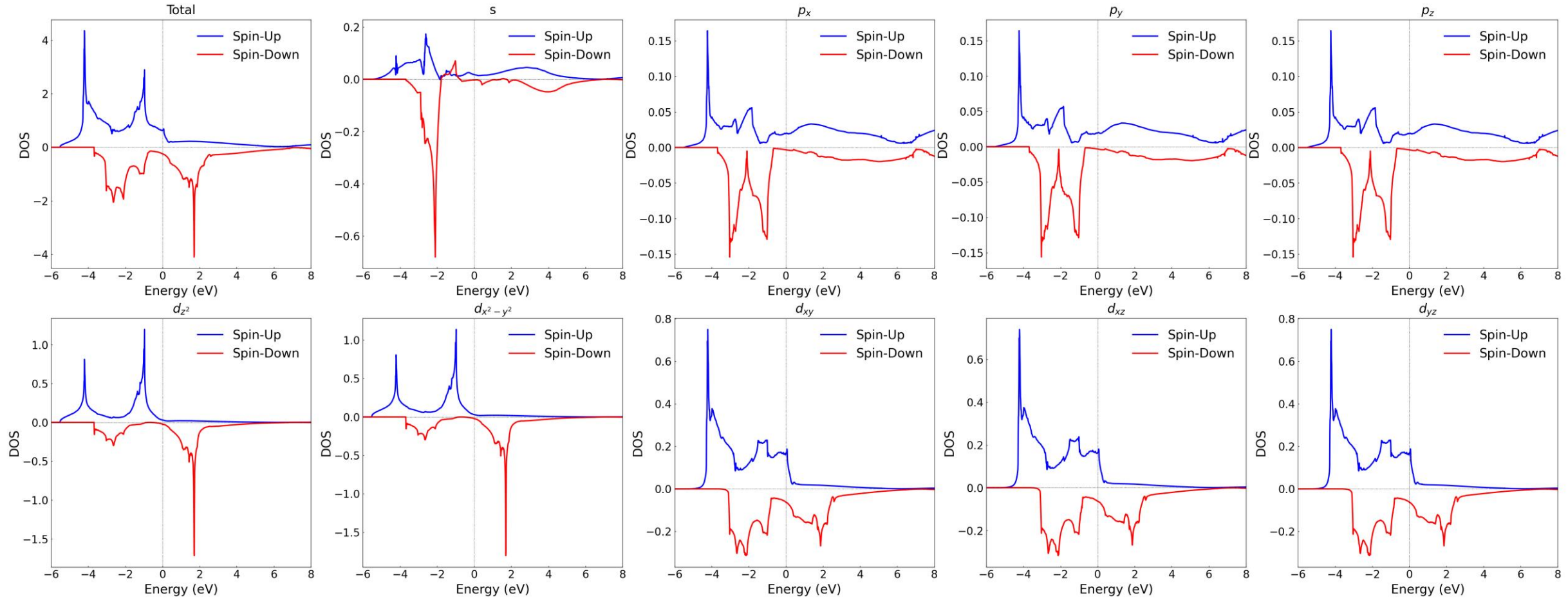
Number of atoms=1
pdos_n=1
1
<Spectra_Tetrahedron> start
```

By answering these questions, you can calculate DOS or PDOS for a specific atom and orbitals.

Fe using OpenMX: Ground-State, Band Structure, DOS

Now, use the following command to plot the DOS

```
python3 Plot-DOS.py
```



Fe using OpenMX: Heisenberg Exchange Parameters



Fe: Adiabatic Magnon Dispersion

In this exercise, you will learn how to calculate the Heisenberg exchange parameters of ferromagnetic Fe using OpenMX. In the next step, we will calculate the magnon dispersion of Fe within spin wave theory.

1. Create a subdirectory in your directory and call it 'Fe-Jx'

```
mkdir ./Fe-Jx
```

2. First copy the following directory to you own directory using cp command:

```
cp ../alikefay/Fe-Jx/* ./Fe-Jx
```

3. 'cd' to the subdirectory and open the Fe.in file. Try to understand all the keywords. As you can see, we are using many k-points for this calculation. This is because we want to get good convergence and accuracy for the Heisenberg exchange parameter.
4. In addition, the basis set of Fe is chosen as **Fe4.0H-s2p2d2**. This is because we want very localized magnetic moments for adopting the Heisenberg model.

Fe: Adiabatic Magnon Dispersion

Fe.in

```
# File Name
System.CurrentDirectory      ./                # default=./
System.Name                  Fe
DATA.PATH                    /home/opt/pkg/openmx/3.9.9-intel-2020.3/DFT_DATA19
level.of.stdout              1                  # default=1 (1-3)
level.of.fileout             0                  # default=1 (0-2)

# Definition of Atomic Species
Species.Number               1
<Definition.of.Atomic.Species
| Fe Fe4.0H-s2p2d2 Fe_PBE19H
Definition.of.Atomic.Species>

# Atoms
Atoms.Number                  1
Atoms.SpeciesAndCoordinates.Unit Ang           # Ang|AU
<Atoms.SpeciesAndCoordinates
| 1 Fe 0.0000 0.0000 0.0000 9.0 7.0
Atoms.SpeciesAndCoordinates>

Atoms.UnitVectors.Unit       Ang               # Ang|AU
<Atoms.UnitVectors
| -1.43300 1.43300 1.43300
| 1.43300 -1.43300 1.43300
| 1.43300 1.43300 -1.43300
Atoms.UnitVectors>
```

Fe: Adiabatic Magnon Dispersion

```
# SCF or Electronic System
scf.XcType          GGA-PBE
scf.SpinPolarization On
scf.SpinOrbit.Coupling Off
HS.fileout          On
scf.ElectronicTemperature 300.0
scf.energycutoff    600.0
scf.maxIter         800
scf.EigenvalueSolver band
scf.Kgrid            27 27 27
scf.Mixing.Type      rmm-diisk
scf.Init.Mixing.Weight 0.300
scf.Min.Mixing.Weight 0.001
scf.Max.Mixing.Weight 0.400
scf.Mixing.History    5
scf.Mixing.StartPulay 6
scf.criterion        1.0e-7

Band.dispersion      on
Band.Nkpath          5
<Band.kpath
  40  0.0000 0.0000 0.0000 0.5000 -0.500 0.5000  G H
  50  0.5000 -0.500 0.5000 0.0000 0.0000 0.5000  H N
  30  0.0000 0.0000 0.5000 0.0000 0.0000 0.0000  N G
  30  0.0000 0.0000 0.0000 0.2500 0.2500 0.2500  G P
  40  0.2500 0.2500 0.2500 0.5000 -0.500 0.5000  P H
Band.kpath>

# MD or Geometry Optimization
MD.Type             NOMD
MD.maxIter          1
MD.Opt.criterion    3.0e-4

# LDA|LSDA-CA|LSDA-PW|GGA-PBE
# On|Off|NC
# On|Off, default=off
# on|off, default=off
# default=300 (K)
# default=150 (Ry)
# default=40
# Recursion|Cluster|Band
# means nk1xnk2xnk3
# Simple|Rmm-Diis|Gr-Pulay
# default=0.30
# default=0.001
# default=0.40
# default=5
# default=6
# default=1.0e-6 (Hartree)

# on|off, default=off

# NOMD|OptC1|OptC2|OptC3|OptC4|OptC5|OptC6|OptC7|RFC5|RFC6|RFC7
# default=1
# default=0.0003 (Hartree/Bohr)
```

Fe: Adiabatic Magnon Dispersion

4. Another important keyword is '**HS.fileout On**'. In fact, jx is a post processing calculation and we need first to run a ground state calculation. Please note that current version of OpenMX only support colinear calculation for jx.

5. Activate the module using:

```
module unload openmpi/4.1.4-intel-2023.0.0
module unload intel/2023.0.0
module load intel/2020.3
module load openmx
```

5. Now you can run the job. While you are waiting for the run to be finished, open the Fe.config file and try to understand the keywords.

```
nohup mpirun -np 5 openmx Fe.in > Fe.out -f openmp &
```

6. When the calculation is finished, run the jx using the following command,

```
nohup mpirun -np 5 jx Fe.scfout Fe.config > Jx.log -f openmp &
```

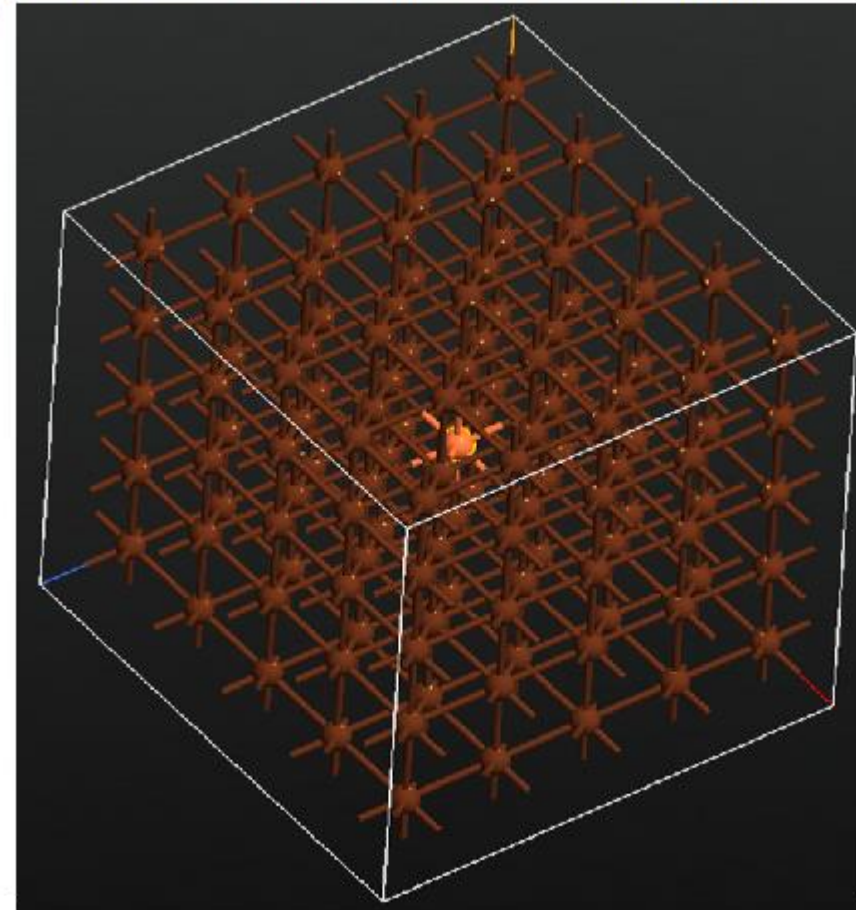
Fe: Adiabatic Magnon Dispersion

Fe.config

```
Flag.PeriodicSum      Off      # default - off (NOTE: No role in cluster calculations.)
Num.Poles             60      # number of poles NP for the nite pole approximation of Fermi function
Num.Kgrid            27 27 27 # These values should be same or a bit larger than those in OpenMX calculation.
Num.ij.pairs         124      # NOTE: Number of ij pairs.
Bunch.ij.pairs       124      # default - 1
```

<ijpairs.cellid

```
1  1 -2 -2 -2
1  1 -2 -2 -1
1  1 -2 -2  0
1  1 -2 -2  1
1  1 -2 -2  2
1  1 -2 -1 -2
1  1 -2 -1 -1
1  1 -2 -1  0
1  1 -2 -1  1
1  1 -2 -1  2
1  1 -2  0 -2
1  1 -2  0 -1
1  1 -2  0  0
1  1 -2  0  1
1  1 -2  0  2
1  1 -2  1 -2
1  1 -2  1 -1
1  1 -2  1  0
1  1 -2  1  1
1  1 -2  1  2
```



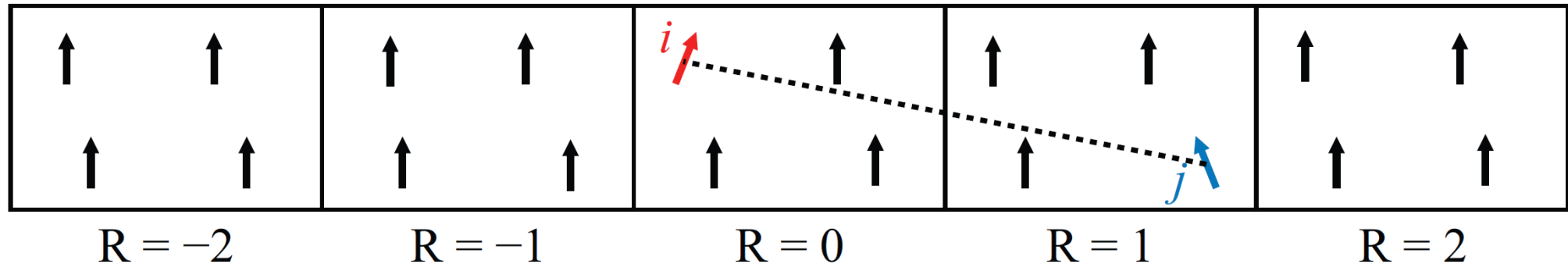
Fe: Adiabatic Magnon Dispersion

The code calculates the Heisenberg constants as,

$$J_{i0,j\mathbf{R}} = \frac{1}{2} \sum_{p=1}^{N_P} \tilde{R}_p \sum_{\mu,\nu \in i} \sum_{\mu',\nu' \in j} \text{Re} \left\{ [\hat{P}_i]_{\nu\mu} G_{i\mu,j\nu'}^+(\downarrow, \tilde{z}_p, \mathbf{R}) [\hat{P}_j]_{\nu'\mu'} G_{j\mu',i\nu}^+(\uparrow, \tilde{z}_p, -\mathbf{R}) \right\}$$

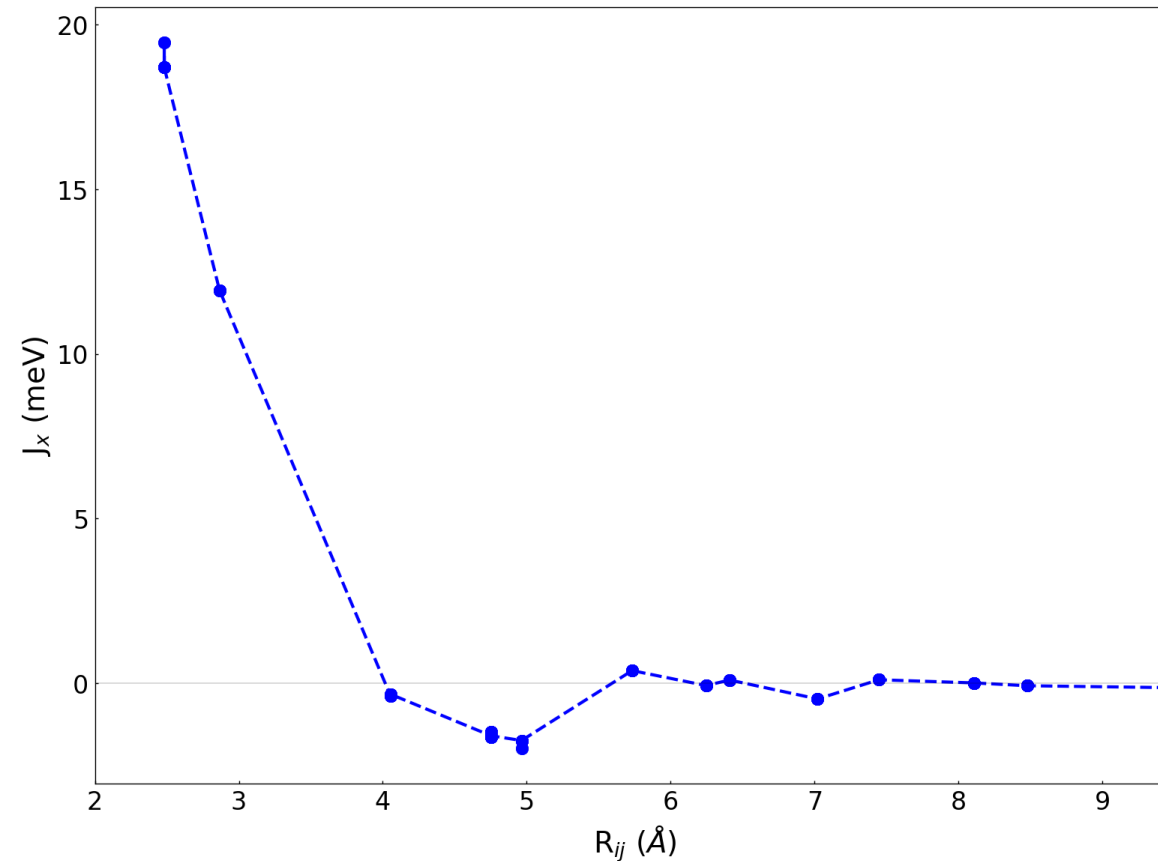
$$G_{j\mu',i\nu}^+(\uparrow, \varepsilon, -\mathbf{R}) \equiv \int d^3 \left(\frac{ka}{2\pi} \right) e^{i\mathbf{k} \cdot \mathbf{R}} \sum_n \frac{C_{j\mu',n\uparrow}(\mathbf{k}) C_{i\mu,n\uparrow}(\mathbf{k})}{\varepsilon + i\eta - \varepsilon_{n\uparrow}(\mathbf{k})}$$

$$G_{i\mu,j\nu'}^+(\downarrow, \varepsilon, \mathbf{R}) \equiv \int d^3 \left(\frac{ka}{2\pi} \right) e^{-i\mathbf{k} \cdot \mathbf{R}} \sum_{n'} \frac{C_{i\mu,n'\downarrow}(\mathbf{k}) C_{j\nu',n'\downarrow}(\mathbf{k})}{\varepsilon + i\eta - \varepsilon_{n'\downarrow}(\mathbf{k})}.$$



Fe: Adiabatic Magnon Dispersion

Here, we calculated the Heisenberg exchange constants for all neighbors in the demonstrated supercell of the Fe atom in the middle of the supercell (The supercell contains 124 atoms with one centered and 123 neighboring Fe atoms). When the calculation is done run: `python3 Plot-HeisP.py`



Fe: Adiabatic Magnon Dispersion

In the next step, run the following script: `python3 Plot-Magnon.py`

It plots the calculated exchange constants as well as the adiabatic magnon dispersion along high symmetry lines in the Brillouin zone using the spin wave theory:

$$\hbar\omega_{\mathbf{k}} = 2S \sum_j J_{ij} \left[1 - e^{i\mathbf{k} \cdot (\mathbf{R}_i - \mathbf{R}_j)} \right]$$

