

# Hands-On: FukuiGrid - Fukui Functions $f^+(r)$ by Interpolation

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## 1 Preparing the Input Files

Before executing the code, ensure that you have the required input files in the working directory. You should have four CHGCAR files corresponding to the system at different charge states:

- CHGCAR for the neutral system ( $\delta N = 0.0$ ).
- CHGCAR files for the system with fraction charges ( $\delta N = +0.05, +0.10, +0.15$ ).

## 2 Example: ZrC (001) Surface

To run FukuiGrid, execute the following command in the terminal:

```
python FukuiGrid.py
```

Upon execution, the following main menu will be displayed:

```
***** Main Menu *****
1 — Fukui Function via Interpolation
2 — Fukui Potential via Electrodes
3 — Fukui Potential via SCPC
4 — Process Grid Data
5 — Perturbative Expansion
6 — Exit
```

Choose an option: 1

Select option 1 for Fukui Function via Interpolation:

```
Chose option 1: Fukui Function via Interpolation.
```

```
11 Electrophilic Fukui function  $f^-(r)$ .
```

```
12 Nucleophilic Fukui function  $f^+(r)$ .
```

```
Choose an option: 12
```

If you select option 12, you will be prompted to choose the  $\delta N$  values:

```
Do you want to use the default  $\delta N$  [0.0, +0.05, +.010 and +0.15]?
```

```
121 Yes, I would like to use the default array
```

```
122 No, I prefer to define my own values
```

Choose an option: 121

If you select option 121, you will be prompted to enter the CHGCAR files for the  $\delta N$  default values:

```
Name CHGCAR files with  $\delta N$ : 0.0, +0.05, +.010 and +0.15
```

```
Enter name of file 1: CHGCAR_00
```

```
Enter name of file 2: CHGCAR+005
```

```
Enter name of file 3: CHGCAR+010
```

```
Enter name of file 4: CHGCAR+015
```

Once the calculation is complete, the output file “CHGCAR\_FUKUI.vasp” will be generated. This file contains the computed Fukui function mapped onto the original charge density grid.

After execution, you will return to the main menu. To exit FukuiGrid, choose option 6:

Choose an option: 6

You selected option 6: Goodbye!

For the files available at <https://github.com/cacarden/FukuiGrid/tree/main/examples/Interpolation/fukui+>, which correspond to the ZrC (001) surface, the computed Fukui function  $f^+(r)$  is obtained and is shown in Fig. 1.

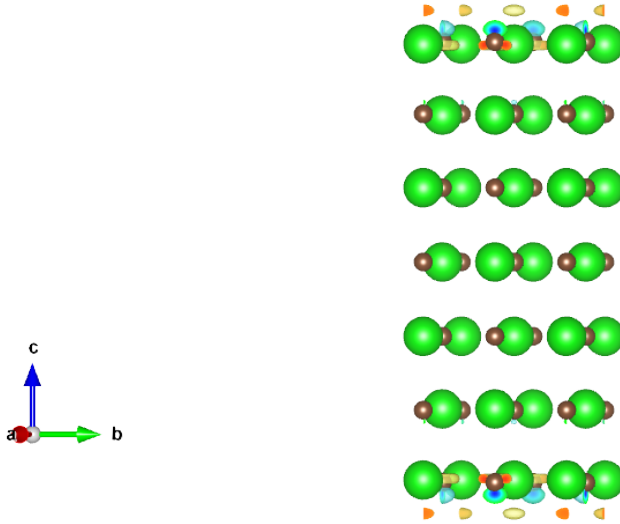


Figure 1: Isosurface representation of  $f^+(r)$  with isovalue  $0.0015 a_0^{-3}$ .