

## Example: orbital decomposition of electron density and envelope wavefunctions

This file is intended to detail the calculation presented in Sections 4.3.3 of our *BinPo* manuscript. Importantly, we remind you that any *BinPo* component looks for the input parameters at the command-line at first, whereas the omitted and more advanced parameters are set by default from the configuration file. Keeping this in mind, we will show how the configuration files are set for each case and what to type from command-line. For a description of all the parameters in the configuration files, see `~/config_files/help_config.md`. Those parameters updatable by command-line in any *BinPo* components (*BP-component.py*) can be checked by means of help command as:

```
$ python BP-component.py -h
```

### Orbital decomposition of electron density

Firstly, we will show how the configuration file *orb\_density.yaml* is set:

```
3 ---
4 DENSITY_ANALYSIS :
5     identifier : "sto_01"
6     sqrt_kgrid_numbers : 25
7     k_shift : [0.0, 0.0]
8     PLOT_ADJUST :
9         plotstyle : 'ggplot'
10        linewidth : 1.7
11        markersize : 8
12        axis_adjust : [0.2, 0.15, 0.9, 0.9] #[left, bottom, right, top]
13        fig_size : [8,6]
14        color_seq : k,r,lime,b
15        fill_curves : yes
16        alpha : 0.25
17        title : "Orbital decomposition of electron density"
18        title_size : 14
19    LABELS :
20        xlabel : 'planes'
21        xfontsize : 16
22        ylabel : 'electron density [elecs./plane]'
23        yfontsize : 16
24        ticksize : 12
25        legends : yes
26        legend_size : 15
27    SAVING :
28        save_data : yes
29        save_plot : yes
30        format : ".png"
31        dpi : 300
32 ...
```

We will compute an electron density profile with the contributions from different orbital characters:

```
$ python BP-orb_density.py -id run2 -aa 0.6
```

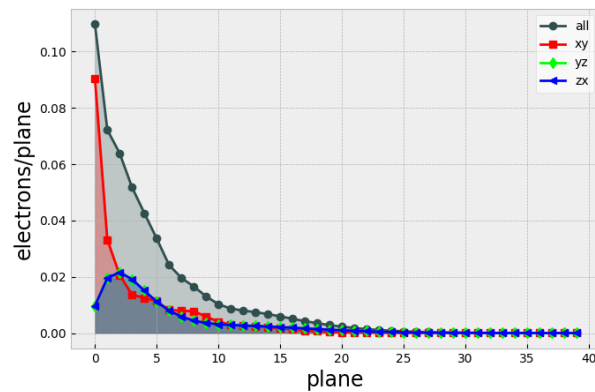
where again we recalled the SC-solution under the “run2” identifier. We also set the curves opacity (*aa*) to 0.6. The output in command-line will be is:

```

DETAILS:
  Identifier: run2
  Surface : STO(100)
  Number of planes: 40
  K-grid: 25 x 25
  Temperature: 10 K
  Fermi level: 11.47650 eV

```

whereas the interactive output plot will be:



### Envelope wavefunctions at the $\Gamma$ -point

The *envelope\_wfs.yaml* configuration file looks:

```

3  ---
4  ENVELOPE_WAVEFUNCTIONS:
5      identifier : "sto_01"
6      intensity_factor : 0.35
7      number_of_wavefunctions : 8
8      PLOT_ADJUST :
9
10         plotstyle : 'ggplot'
11         linewidth_wfs : 2.5
12         linewidth_V : 1.8
13         markersize : 8
14         axis_adjust : [0.2, 0.15, 0.9, 0.9] #[left, bottom, right, top]
15         fig_size : [8,6]
16         color_V : "k"
17         color_wfs : "teal"
18         alpha : 0.3
19         title : ""
20         title_size : 12
21
22  LABELS :
23
24     xlabel : 'planes'
25     xfontsize : 16
26     ylabel : 'Energy [eV]'
27     yfontsize : 16
28     ticksize : 12
29     legends : yes
30     legend_size : 12
31
32  SAVING :
33
34     save_data : yes
35     save_plot : yes
36     format : ".png"
37     dpi : 300
38  ...

```

Now, we will compute the envelope wavefunctions by typing:

```
$ python BP-envelope_wfs.py -id run2 -sf 3.0 -nw 8 -xy -2 22.5 -0.35 0.06
```

Here, besides the identifier, we specify the scale factor (*sf*) which affects the amplitude of the envelope wavefunctions, the number of these wavefunctions (*nw*) to compute and the x and y limits of the plot by means of the *xy* parameter. The output in command-line will be:

```
DETAILS:
Identifier: run2
Surface : STO(100)
Number of planes: 40
Temperature: 10 K
Fermi level: 11.47650 eV
Number of wavefunctions: 8
```

whereas the output interactive plot will be:

