WIEN2k 24.1 Install with OneAPI (ifx)

August 31, 2025

Linux Operating System Version (lsb_release -a): Ubuntu 24.04.3 LTS Fortran Compiler Version (ifx -v): ifx version 2025.2.1 WIEN2k Version (cat \$WIENROOT/WIEN2k VERSION): 24.1 (Release 1/8/2024)

The Ubuntu 24.04.3 LTS was already installed, but if that is needed it should be available at:

https://ubuntu.com/download/desktop

Operating system packages needed were installed before proceeding with step 1 given below. If that hasn't been done, then install them:

username@computername:~\$ sudo apt update

. . .

username@computername:~ \$ sudo apt upgrade

. . .

username@computername:~ \$ sudo apt install tesh make autoconf libtool git

. .

username@computername:~ \$ sudo apt install gnuplot gnuplot-x11 libgd-dev gedit

. . .

username@computername:~\$ sudo apt install libquadmath0 libglu1-mesa-dev

username@computername:~\$ sudo apt install libtogl-dev tcl-dev tk-dev libfftw3-dev libxmu-dev

. .

Note: If prompted with "Do you want to continue [Y/n]", enter: Y.

The XCrySDen 1.6.2 was installed using:

username@computername:~ \$ sudo apt install xcrysden

username@computername:~ \$ gedit ~/.bashrc

XCRYSDEN_TOPDIR=/usr/bin XCRYSDEN_LIB_BINDIR=/usr/lib export XCRYSDEN_TOPDIR XCRYSDEN_LIB_BINDER

username@computername:~ \$ source ~/.bashrc

Of note, for XCrySDen to work with Wayland, toglOpt(accum) needs set to false in the /usr/share/xcrysden/Tcl/custom-definitions file:

https://www.mail-archive.com/wien@zeus.theochem.tuwien.ac.at/msg23440.html

Restart the computer.

Installed OneAPI [https://www.intel.com/content/www/us/en/docs/oneapi/installation-guide-linux/2025-2/hpc-apt.html#HPC-APT] using:

```
username@computername:~$ wget -O- https://apt.repos.intel.com/intel-gpg-keys/GPG-PUB-
KEY-INTEL-SW-PRODUCTS.PUB \
| gpg --dearmor | sudo tee /usr/share/keyrings/oneapi-archive-keyring.gpg > /dev/null
username@computername:~$ echo "deb [signed-by=/usr/share/keyrings/oneapi-archive-
keyring.gpg] https://apt.repos.intel.com/oneapi all main" | sudo tee
/etc/apt/sources.list.d/oneAPI.list
username@computername:~$ sudo apt update
username@computername:~\$ sudo apt install intel-hpckit
Do you want to continue? [Y/n] Y
done.
Configuring autoload of sep5.service service ...
Created symlink /etc/systemd/system/multi-user.target.wants/sep5.service → /usr/
lib/systemd/system/sep5.service.
done.
Setting up intel-oneapi-openmp-common-2025.2 (2025.2.1-7) ...
Setting up intel-oneapi-mpi-devel-2021.16 (2021.16.1-803) ...
Setting up intel-oneapi-common-oneapi-vars (2025.2.0-377) ...
/var/lib/dpkg/info/intel-oneapi-common-oneapi-vars.postinst: 27: cannot create /
tmp//oneapi/-2025.2-link-tlt.log: Directory nonexistent
Setting up intel-oneapi-tcm-1.4 (1.4.0-345) ...
Setting up intel-oneapi-dpcpp-debugger-2025.2 (2025.2.0-388) ...
Setting up intel-pti-0.13 (0.13.1-10) ...
Setting up intel-oneapi-compiler-shared-common-2025.2 (2025.2.1-7) ...
Setting up intel-oneapi-mpi-devel (2021.16.1-803) ...
Setting up intel-oneapi-dpcpp-ct-2025.2 (2025.2.0-517) ...
Setting up intel-oneapi-dal-common-devel-2025.8 (2025.8.0-11) ...
Setting up intel-oneapi-compiler-dpcpp-eclipse-cfg-2025.2 (2025.2.1-7) ...
Setting up intel-oneapi-compiler-fortran-common-2025.2 (2025.2.1-7) ...
Setting up intel-oneapi-mkl-classic-include-2025.2 (2025.2.0-628) ...
Setting up intel-oneapi-dal-common-2025.8 (2025.8.0-11) ...
Setting up intel-oneapi-ishmem-1.4 (1.4.0-244) ...
Setting up intel-oneapi-dev-utilities-eclipse-cfg-2025.2 (2025.2.1-30) ...
Setting up intel-oneapi-ippcp-2025.2 (2025.2.0-447) ...
Setting up intel-oneapi-ccl-devel-2021.16 (2021.16.1-9) ...
Setting up intel-oneapi-compiler-dpcpp-cpp-common-2025.2 (2025.2.1-7) ...
```

```
Setting up intel-oneapi-ishmem-devel-1.4 (1.4.0-244) ...
Setting up intel-oneapi-umf-0.11 (0.11.0-394) ...
Setting up intel-pti-dev-0.13 (0.13.1-10) ...
Setting up intel-oneapi-openmp-2025.2 (2025.2.1-7) ...
Setting up intel-oneapi-tbb-2022.2 (2022.2.0-507) ...
Setting up intel-oneapi-tbb-devel-2022.2 (2022.2.0-507) ...
Setting up intel-oneapi-ccl-devel (2021.16.1-9) ...
/var/lib/dpkg/info/intel-oneapi-ccl-devel.postinst: 27: cannot create /tmp//onea
pi/-2021.16-link-tlt.log: Directory nonexistent
Setting up intel-oneapi-mkl-sycl-include-2025.2 (2025.2.0-628) ...
Setting up intel-oneapi-ipp-2022.2 (2022.2.0-582) ...
Setting up intel-oneapi-dpcpp-ct (2025.2.0-517) ...
Setting up intel-oneapi-compiler-shared-runtime-2025.2 (2025.2.1-7) ...
Setting up intel-oneapi-compiler-shared-2025.2 (2025.2.1-7) ...
Setting up intel-oneapi-mkl-core-2025.2 (2025.2.0-628) ...
Setting up intel-oneapi-ipp-devel-2022.2 (2022.2.0-582) ...
Setting up intel-oneapi-tbb-devel (2022.2.0-507) ...
Setting up intel-oneapi-mkl-cluster-2025.2 (2025.2.0-628) ...
Setting up intel-oneapi-ippcp-devel-2025.2 (2025.2.0-447) ...
Setting up intel-oneapi-dev-utilities-2025.2 (2025.2.1-30) ...
Setting up intel-pti-dev (0.13.1-10) ...
/var/lib/dpkg/info/intel-pti-dev.postinst: 27: cannot create /tmp//oneapi/-0.13-
link-tlt.log: Directory nonexistent
Setting up intel-oneapi-ishmem-devel (1.4.0-244) ...
/var/lib/dpkg/info/intel-oneapi-ishmem-devel.postinst: 27: cannot create /tmp//o
neapi/-1.4-link-tlt.log: Directory nonexistent
Setting up intel-oneapi-dev-utilities (2025.2.1-30) ...
Setting up intel-oneapi-mkl-core-devel-2025.2 (2025.2.0-628) ...
Setting up intel-oneapi-ipp-devel (2022.2.0-582) ...
/var/lib/dpkg/info/intel-oneapi-ipp-devel.postinst: 27: cannot create /tmp//onea
pi/-2022.2-link-tlt.log: Directory nonexistent
Setting up intel-oneapi-ippcp-devel (2025.2.0-447) ...
/var/lib/dpkg/info/intel-oneapi-ippcp-devel.postinst: 27: cannot create /tmp//on
eapi/-2025.2-link-tlt.log: Directory nonexistent
Setting up intel-oneapi-compiler-dpcpp-cpp-runtime-2025.2 (2025.2.1-7) ...
Setting up intel-oneapi-mkl-sycl-stats-2025.2 (2025.2.0-628) ...
Setting up intel-oneapi-mkl-sycl-rng-2025.2 (2025.2.0-628) ...
Setting up intel-oneapi-dal-2025.8 (2025.8.0-11) ...
Setting up intel-oneapi-compiler-fortran-runtime-2025.2 (2025.2.1-7) ...
Setting up intel-oneapi-mkl-sycl-vm-2025.2 (2025.2.0-628) ...
Setting up intel-oneapi-mkl-sycl-data-fitting-2025.2 (2025.2.0-628) ...
Setting up intel-oneapi-dnnl-2025.2 (2025.2.0-561) ...
Setting up intel-oneapi-compiler-fortran-2025.2 (2025.2.1-7) ...
Setting up intel-oneapi-mkl-sycl-dft-2025.2 (2025.2.0-628) ...
Setting up intel-oneapi-mkl-sycl-blas-2025.2 (2025.2.0-628) ...
```

```
Setting up intel-oneapi-mkl-cluster-devel-2025.2 (2025.2.0-628) ...
Setting up intel-oneapi-dpcpp-cpp-2025.2 (2025.2.1-7) ...
Setting up intel-oneapi-dnnl-devel-2025.2 (2025.2.0-561) ...
Setting up intel-oneapi-mkl-sycl-lapack-2025.2 (2025.2.0-628) ...
Setting up intel-oneapi-mkl-sycl-sparse-2025.2 (2025.2.0-628) ...
Setting up intel-oneapi-mkl-classic-devel-2025.2 (2025.2.0-628) ...
Setting up intel-oneapi-dnnl-devel (2025.2.0-561) ...
Setting up intel-oneapi-mkl-sycl-2025.2 (2025.2.0-628) ...
Setting up intel-oneapi-compiler-dpcpp-cpp-2025.2 (2025.2.1-7) ...
Setting up intel-oneapi-mkl-sycl-devel-2025.2 (2025.2.0-628) ...
Setting up intel-oneapi-compiler-dpcpp-cpp (2025.2.1-7) ...
Setting up intel-oneapi-mkl-devel-2025.2 (2025.2.0-628) ...
Setting up intel-oneapi-mkl-devel (2025.2.0-628) ...
Setting up intel-oneapi-dal-devel-2025.8 (2025.8.0-11) ...
Setting up intel-oneapi-dal-devel (2025.8.0-11) ...
Setting up intel-oneapi-hpc-toolkit (2025.2.1-40) ...
Setting up intel-hpckit (2025.2.1-40) ...
username@computername:~\$ echo "source /opt/intel/oneapi/setvars.sh intel64" >> ~/.bashrc
username@computername:~\$ grep oneapi ~/.bashrc
source /opt/intel/oneapi/setvars.sh intel64
username@computername:~$ source ~/.bashrc
:: initializing one API environment ...
 bash: BASH VERSION = 5.2.21(1)-release
 args: Using "$@" for setvars.sh arguments: intel64
:: advisor -- latest
:: ccl -- latest
:: compiler -- latest
:: dal -- latest
:: debugger -- latest
:: dev-utilities -- latest
:: dnnl -- latest
:: dpcpp-ct -- latest
:: dpl -- latest
:: ipp -- latest
:: ippcp -- latest
:: ishmem -- latest
:: mkl -- latest
:: mpi -- latest
:: pti -- latest
:: tbb -- latest
:: umf -- latest
:: vtune -- latest
:: oneAPI environment initialized ::
```

```
username@computername:~$ ifx -v
   ifx version 2025.2.1
   username@computername:~\$ icx -v
   Intel(R) oneAPI DPC++/C++ Compiler 2025.2.1 (2025.2.0.20250806)
   username@computername:~$ mpiifx -v
   ifx version 2025.2.1
1. The following was entered into the terminal to install libxc:
   username@computername:~$ cd ~
   username@computername:~$ wget
   https://gitlab.com/libxc/libxc/-/archive/7.0.0/libxc-7.0.0.tar.bz2
   username@computername:~\$ tar xvf libxc-7.0.0.tar.bz2
   username@computername:~$ cd libxc-7.0.0/
   username@computername:~/libxc-7.0.0$ autoreconf -i --force
   username@computername:~/libxc-7.0.0$ ./configure FC=ifx CC=icx
   --prefix=$HOME/libxc-7.0.0
   username@computername:~/libxc-7.0.0$ make
   username@computername:~/libxc-7.0.0$ make check
   PASS: xc-run testsuite
   Testsuite summary for libxc 7.0.0
   #TOTAL: 1
   # PASS: 1
   # SKIP: 0
   # XFAIL: 0
   #FAIL: 0
   # XPASS: 0
   # ERROR: 0
```

Page **5** of **16**

. . .

```
username@computername:~/libxc-7.0.0$ make install
   username@computername:~/libxc-7.0.0$ ls ~/libxc-7.0.0/lib
   libxc.a libxcf03.a libxcf03.la libxc.la pkgconfig
2. The following was entered into the terminal to install fftw:
   username@computername:~/libxc-7.0.0$ cd ~
   username@computername:~\$ wget https://www.fftw.org/fftw-3.3.10.tar.gz
   username@computername:~\$ tar xvf fftw-3.3.10.tar.gz
   username@computername:~$ cd fftw-3.3.10/
   username@computername:~/fftw-3.3.10\$ ./configure FCC=ifx CC=icx MPICC=mpiicx
   CFLAGS="-gcc-sys" --enable-mpi --prefix=$HOME/fftw-3.3.10
   username@computername:~/fftw-3.3.10$ make
   username@computername:~/fftw-3.3.10$ make install
   username@computername:~/fftw-3.3.10\$ ls ~/fftw-3.3.10/include/ ~/fftw-3.3.10/lib
   /home/username/fftw-3.3.10/include/:
                     fftw31-mpi.f03 fftw3-mpi.h
   fftw3.f fftw3.h
   fftw3.f03 fftw31.f03 fftw3-mpi.f03 fftw3q.f03
   /home/username/fftw-3.3.10/lib:
   cmake libfftw3.a libfftw3.la libfftw3 mpi.a libfftw3 mpi.la pkgconfig
3. The following was entered into the terminal to install ELPA:
   username@computername:~/fftw-3.3.10$ cd ~
   username@computername:~$ wget
   https://gitlab.mpcdf.mpg.de/elpa/elpa/-/archive/new release 2025.06.001/elpa-
   new release 2025.06.001.tar.gz
   username@computername:~\$ tar xvf elpa-new release 2025.06.001.tar.gz
   username@computername:~$ cd elpa-new release 2025.06.001/
   username@computername:~/elpa-new release 2025.06.001$./autogen.sh
   username@computername:~/elpa-new release 2025.06.001$ ./configure FC=mpiifx CC=mpiicx
```

SCALAPACK LDFLAGS="-L\$MKLROOT/lib/intel64 -lmkl scalapack lp64 -lmkl intel lp64

```
-lmkl sequential -lmkl core -lmkl blacs intelmpi lp64 -lpthread -lm -ldl -liomp5 -lm -Wl,-
rpath,$MKLROOT/lib/intel64" SCALAPACK FC FLAGS="-L$MKLROOT/lib/intel64"
-lmkl scalapack lp64 -lmkl intelmpi lp64 -lmkl sequential -lmkl core
-lmkl blacs intelmpi lp64 -lpthread -lm -I$MKLROOT/include/intel64/lp64" --
prefix=$HOME/elpa-new release 2025.06.001 FCFLAGS=-O3 CFLAGS="-O3 -mfma
-funsafe-math-optimizations -ftree-vectorize" LIBS="-lmpi -lmpifort" --disable-avx512
The following ELPA2 kernels will be build:
real generic
real generic simple
real generic simple block4
real generic simple block6
real sse block2
real sse block4
real sse block6
real sse assembly
real avx block2
real avx block4
real avx block6
real avx2 block2 (default)
real avx2 block4
real avx2 block6
complex generic
complex generic simple
complex sse block1
complex sse block2
complex sse assembly
complex avx block1
complex avx block2
complex avx2 block1 (default)
complex avx2 block2
This version of ELPA support the minimal API version: 20170403
The current API version is: 20250131
This version of ELPA support the minimal autotuning version: 20171201
The current autotune version is: 20250131
username@computername:~/elpa-new release 2025.06.001$ make
username@computername:~/elpa-new release 2025.06.001$ make install
username@computername:~/elpa-new release 2025.06.001$ ls ~/elpa-
new release 2025.06.001/lib ~/elpa-new release 2025.06.001/include/elpa-2025.06.001/elpa/
/home/username/elpa-new release 2025.06.001/include/elpa-2025.06.001/elpa/:
```

```
elpa_configured_options.h elpa_explicit_name.h elpa_generated.h elpa.h elpa_version.h elpa_constants.h elpa_generated_c_api.h elpa_generic.h elpa_simd_constants.h /home/username/elpa-new_release_2025.06.001/lib: libelpa.a libelpa.la libelpa.so libelpa.so libelpa.so.19 libelpa.so.19.4.1 pkgconfig
```

4. Go to:

http://www.wien2k.at/reg_user/index.html

- 5. Click "Code download (after registration)"
- 6. Enter your username and password that you were given when you purchased WIEN2k.
- 7. Click the link "here" in "You can download the complete source code (about 170 MB) from here" to download "WIEN2k_24.1.tar". You should now have the source code package for WIEN2k 24.1.
- 8. If you have the folder "WIEN2k" in your Linux operating system home directory (for example at /home/username/WIEN2k, where username is replaced by your user name). You can rename it something else, such as "WIEN2k23.2".
- 9. The following was entered into the terminal to install WIEN2k 24.1:

```
username@computername:~/elpa-new release 2025.06.001$ cd ~
username@computername:~$ mkdir WIEN2k
username@computername:~$ cd WIEN2k
username@computername:~/WIEN2k$ ls
WIEN2k 24.1.tar
username@computername:~/WIEN2k$ tar xvf WIEN2k_24.1.tar
username@computername:~/WIEN2k$ gunzip *.gz
username@computername:~/WIEN2k$ chmod +x ./expand lapw
username@computername:~/WIEN2k$ export LC NUMERIC=en US.UTF-8
username@computername:~/WIEN2k$ ./expand lapw
continue (y/n)
y
...
Applied patches at: https://github.com/gsabo/WIEN2k-Patches/tree/master/24.1
username@computername:~/WIEN2k$ ./siteconfig
continue or stop (c/s) c
```

```
Press RETURN to continue
 Selection: LI
  Press RETURN to continue
   Your compiler: ifx
   Your compiler: icx
  Press RETURN to continue
Hit Enter to continue
Would you like to use LIBXC (needed ONLY for self-consistent gKS mGGA calculations, for
the stress tensor and experts who want to play with different DFT options. It must have been
installed before)? (y,N):
y
Do you want to automatically search for LIBXC installations? (Y,n):
Please specify a comma separated list of directories to search! (If no list is entered, /usr/lib64,
/usr/local and /opt will be searched as default):
/home/username/libxc-7.0.0
More than one libxc was found in the specified directory(ies).
Pick one of the following (enter the number of the line of your choice or 0 to manually specify a
path)!:
/home/username/libxc-7.0.0/lib/libxc.a
/home/username/libxc-7.0.0/src/.libs/libxc.a
1
  Press RETURN to continue
Do you want to automatically search for FFTW installations? (Y,n):
Please specify a comma separated list of directories to search! (If no list is entered,
/usr/lib64, /usr/local and /opt will be searched as default):
/home/username/fftw-3.3.10
Finding the required fftw3 library-files in /home/username/fftw-3.3.10 ....
/home/username/fftw-3.3.10/.libs/libfftw3.a
/home/username/fftw-3.3.10/lib/libfftw3.a
```

```
Do you want to use a FFTW version from the list above? (Y,n):
Please enter the line number of the chosen version or enter 0 to manually specify your choice!:
2
Please specify the target achitecture of your FFTW library (e.g. lib64) or accept present choice
(enter): lib
Please specify the name of your FFTW library or accept present choice (enter): fftw3
Is this correct? (Y,n): Y
Current settings:
 M OpenMP switch:
                           -gopenmp
 O Compiler options:
                          -O -FR -mp1 -w -prec div -pc80 -pad -ip -DINTEL VML
-traceback -assume buffered io -I$(MKLROOT)/include
                        \label{lem:condition} $(FOPT) - L$(MKLROOT)/lib/$(MKL_TARGET\_ARCH) - lpthread
 L Linker Flags:
-lm -ldl -liomp5
 P Preprocessor flags
                         '-DParallel'
 R R LIBS (LAPACK+BLAS): -lmkl intel lp64 -lmkl_intel_thread -lmkl_core
 F FFTW options:
                         -DFFTW3 -I/home/username/fftw-3.3.10/include
   FFTW-LIBS:
                         -L/home/username/fftw-3.3.10/lib -lfftw3
 X LIBX options:
                         -DLIBXC -I/home/username/libxc-7.0.0/include
   LIBXC-LIBS:
                         -L/home/username/libxc-7.0.0/lib -lxcf03 -lxc
 S Save and Quit
Selection: S
  Press RETURN to continue
Shared Memory Architecture? (y/N):y
 Do you know/need a command to bind your jobs to specific nodes?
 (like taskset -c). Enter N / your specific command: N
 Do you have MPI, ScaLAPACK, ELPA, or MPI-parallel FFTW installed and intend
 to run finegrained parallel?
 (y/N) y
 Your compiler: mpiifx
Do you want to use a present ScaLAPACK installation? (Y,n): Y
```

```
Do you want to use the MKL version of ScaLAPACK? (Y,n):Y
Do you use Intel MPI? (Y,n):Y
Is this correct? (Y,n): Y
  Press RETURN to continue
Do you want to use ELPA? (y,N):
Do you want to automatically search for ELPA installations? (Y,n):
Please specify a comma-separated list of directories to search! (If no list is entered,
/usr/lib64 /usr/local and /opt will be searched as default):
/home/username/elpa-new release 2025.06.001
Finding the required ELPA files in /home/username/elpa-new release 2025.06.001 ....
More than one version of ELPA found:
/home/username/elpa-new release 2025.06.001/.libs
/home/username/elpa-new release 2025.06.001/lib
Pick one (enter line number) or enter 0 to manually specify a path!:
Present root directory of ELPA is: /home/username/elpa-release 2025.01.001/
Please specify the ROOT-path of your ELPA installation (like /usr/local/elpa/) or accept present
path (Enter):
Checking for ELPA version ...
More than one set of include files in your ELPA-ROOT directory. Pick one (enter line number)
or enter 0 to manually specify a version!:
/home/username/elpa-new release 2025.06.001/test/Fortran/elpa generalized
/home/username/elpa-new release 2025.06.001/test/Fortran/elpa2
/home/username/elpa-new release 2025.06.001/include/elpa-2025.06.001
/home/username/elpa-new release 2025.06.001/include/elpa-2025.06.001/elpa
/home/username/elpa-new release 2025.06.001/elpa
/home/username/elpa-new release 2025.06.001/share/doc/elpa
/home/username/elpa-new release 2025.06.001/src/elpa generalized
/home/username/elpa-new release 2025.06.001/src/elpa2
/home/username/elpa-new release 2025.06.001/src/elpa1
/home/username/elpa-new release 2025.06.001/.fortran dependencies/elpa2 print kernels
3
Is this correct? (Y,n):Y
Please specify the lib-directory of your ELPA installation (e.g. lib or lib64):
lib
```

```
...
Is this correct? (Y,n):Y
...
Please specify the name of your installed ELPA library (e.g. elpa or elpa_openmp): elpa
Your current lib-directory of your ELPA installation is: lib
Is this correct? (Y,n):Y
The current library name of your ELPA installation is: elpa
Is this correct? (Y,n):Y
...
Is this correct? (Y,n): Y

Press RETURN to continue
...

Please specify your parallel compiler options or accept the recommendations (Enter - default)!:

Please specify your parallel OMP_SWITCH (type "del" to blank it) or accept the recommendations (Enter - default)!:

Please specify your MPIRUN command or accept the recommendations (Enter - default)!:

Press RETURN to continue
...
```

Current settings:

Parallel compiler : mpiifx

SCALAPACK LIBS :-lmkl scalapack lp64 -lmkl blacs intelmpi lp64

FFTW_PLIBS : -lfftw3_mpi ELPA OPT : -DELPA

```
-I/home/username/elpa-new release 2025.06.001/include/elpa-2025.06.001/elpa
              -I/home/username/elpa-new release 2025.06.001/include/elpa-2025.06.001/
   modules
        ELPA LIBS
                            : -lelpa -L/home/username/elpa-new release 2025.06.001/lib -Wl,-
   rpath=/home/username/elpa-new release 2025.06.001/lib
        FPOPT(par.comp.options): -O -FR -mp1 -w -prec div -pc80 -pad -ip -DINTEL VML
   -traceback -assume buffered io -I$(MKLROOT)/include
        OMP SWITCH
                               : -qopenmp
        MPIRUN command
                                : mpirun -np NP -machinefile HOSTS EXEC
    parallel execution:
        RP LIBS
                          : $(R LIBS)
   Selection: S
     Press RETURN to continue
      Selection: Q
      Selection: A
   Compile time errors (if any) were:
                                          <= It should be blank here if successful.
   Check file compile.msg in the corresponding SRC * directory for the
   compilation log and more info on any compilation problem.
     Press RETURN to continue
    Please enter the full path of the perl program: /usr/bin/perl
     Press RETURN to continue
   Please enter the full path to your temporary directory: /tmp
     Press RETURN to continue
10. If it is the first time installing WIEN2k on the computer, enter in the terminal:
   username@computername:~/WIEN2k$ ./userconfig lapw
   Specify your prefered editor (default is emacs):
      editor shall be: gedit
```

Set editor to gedit (Y/n) Y

Specify your prefered DATA directory, where your cases should be stored (for /home/username/WIEN2k, just enter RETURN key): DATA directory: /home/username/wiendata

Set DATA directory to /home/username/wiendata (Y/n) Y

••

Specify your prefered scratch directory, where big case.vector files can be stored (Recommended is a local directory (maybe /scratch), not a NFS directory. For your working directory, just enter RETURN key): scratch directory:

Set scratch directory to working directory (Y/n) Y

Specify your program to read pdf files (default is okular) (on some Linux systems use xpdf, evince, pdfstudio, ...):evince

...

Set PDFREADER to evince (Y/n) Y

WIEN2k can use OpenMP parallelization on multicore computers. For details please read the "Parallelization section" of the Usersguide. Your present computer has 6 cores, but more than 4 (8) cores is useless. How many cores do you want to use by default (4):1

Set OMP_NUM_THREADS to 1 (Y/n) Y

!!! The following lines will be added to your .bashrc file if you continue !!! A copy of your current .bashrc will be saved under .bashrc.savelapw!

```
#export LD LIBRARY PATH=.....
  export EDITOR="gedit"
  export SCRATCH=./
  if [ "$WIENROOT" = "" ]; then export WIENROOT=/home/username/WIEN2k; fi
  export W2WEB CASE BASEDIR=/home/username/wiendata
  export STRUCTEDIT PATH=$WIENROOT/SRC_structeditor/bin
  export PDFREADER=evince
  export PATH=$WIENROOT:$STRUCTEDIT PATH:$WIENROOT/SRC IRelast/script-
  elastic:$PATH:.
  export OCTAVE EXEC PATH=${PATH}::
  export OCTAVE PATH=${STRUCTEDIT PATH}::
  ulimit -s unlimited
  alias octave="octave -p $OCTAVE PATH"
    Do you want to continue (Y/n)? Y
    Edit .rhosts file now? (y/N) N
  username@computername:~/WIEN2k$ source ~/.bashrc
11. To setup or start w2web, enter in the terminal:
  username@computername:~/WIEN2k$ w2web
  # w2web starter
  # Copyright (C) 2001 luitz.at
  w2web installer on host computername
  # w2web installer
  # Copyright (C) 2001 luitz.at
  Checking for Installation in /home/username/.w2web/computername
  Creating /home/username/.w2web
  Creating /home/username/.w2web/computername
  conf directory does not exist - creating it.
  logs directory does not exist - creating it.
```

sessions directory does not exist - creating it. tmp directory does not exist - creating it.

Installing w2web files ...
Please answer these questions for proper installation.
Just press enter for the default value of (in brackets).

Please enter the username: [admin] username Please enter the password: [password] password

username:password

Remember these. You will need them when you log in.

Select the port to run on: [7890]

Running on port 7890

Please enter this system's hostname: [computername] localhost Using localhost

Is this your master node?: [y] y Installing... Attempting to start now...

Trying to start /home/username/WIEN2k/SRC_w2web/bin/w2web w2web server started, now point your web browser to http://localhost:7890

done.

- 12. The WIEN2k 24.1 installation is now complete. To test if WIEN2k installed okay or not, run the TiC example at [1]. Additional details of the TiC example are given in the WIEN2k usersguide (section "3 Quick Start" starting on page 13) [2].
- [1] https://github.com/gsabo/WIEN2k-Docs/blob/main/WIEN2k24.1%20TiC%20Example(ifx).pdf
- [2] http://susi.theochem.tuwien.ac.at/reg_user/textbooks/usersguide.pdf