

WIEN2k 24.1 Install with OneAPI (ifx)

December 21, 2024

Linux Operating System Version (lsb_release -a): Ubuntu 24.04.1 LTS

Fortran Compiler Version (ifx -v): ifx version 2025.0.4

WIEN2k Version (cat \$WIENROOT/WIEN2k_VERSION): 24.1 (Release 1/8/2024)

The Ubuntu 24.04.1 LTS was already installed, but if that is needed it should be available at:

<https://ubuntu.com/download/desktop>

Operating system packages that were 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 tcsh 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 xcrysdn
```

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

```
...
```

```
XCrySDEN_TOPDIR=/usr/bin
```

```
XCrySDEN_LIB_BINDIR=/usr/lib
```

```
export XCrySDEN_TOPDIR XCrySDEN_LIB_BINDIR
```

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

Of note, for XCrySDen to work with Wayland, togIOpt(accum) needs set to false in the /usr/share/xcrysdn/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/2025-0/apt-005.html>] 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.
Setting up intel-oneapi-tcm-1.2 (1.2.0-589) ...
Setting up intel-oneapi-dpcpp-debugger-2025.0 (2025.0.0-663) ...
Setting up intel-oneapi-dal-common-2025.0 (2025.0.1-9) ...
Setting up intel-oneapi-tbb-2022.0 (2022.0.0-402) ...
Setting up intel-oneapi-common-oneapi-vars (2025.0.1-15) ...
Setting up intel-pti-0.10 (0.10.0-284) ...
Setting up intel-oneapi-umf-0.9 (0.9.1-6) ...
Setting up intel-oneapi-dpcpp-ct-2025.0 (2025.0.1-17) ...
Setting up intel-oneapi-compiler-dpcpp-eclipse-cfg-2025.0 (2025.0.4-1519) ...
Setting up intel-oneapi-mkl-classic-include-2025.0 (2025.0.1-14) ...
Setting up intel-oneapi-libdpsd-devel-2022.7 (2022.7.1-14) ...
Setting up intel-oneapi-compiler-shared-common-2025.0 (2025.0.4-1519) ...
Setting up intel-oneapi-dev-utilities-eclipse-cfg-2025.0 (2025.0.0-599) ...
Setting up intel-oneapi-ippcp-2025.0 (2025.0.0-615) ...
Setting up intel-oneapi-icc-eclipse-plugin-cpp-2025.0 (2025.0.4-1519) ...
Setting up intel-oneapi-mpi-2021.14 (2021.14.1-5) ...
Setting up intel-oneapi-compiler-cpp-eclipse-cfg-2025.0 (2025.0.4-1519) ...
Setting up intel-oneapi-ippcp-devel-2025.0 (2025.0.0-615) ...
Setting up intel-oneapi-dev-utilities-2025.0 (2025.0.0-599) ...
Setting up intel-oneapi-tbb-devel-2022.0 (2022.0.0-402) ...
Setting up intel-oneapi-dev-utilities (2025.0.0-599) ...
Setting up intel-oneapi-mpi-devel-2021.14 (2021.14.1-5) ...
Setting up intel-oneapi-ccl-2021.14 (2021.14.0-505) ...
Setting up intel-oneapi-ippcp-devel (2025.0.0-615) ...
Setting up intel-pti-dev-0.10 (0.10.0-284) ...
Setting up intel-oneapi-openmp-2025.0 (2025.0.4-1519) ...
```

Setting up intel-oneapi-dpcpp-ct (2025.0.1-17) ...
Setting up intel-oneapi-mkl-sycl-include-2025.0 (2025.0.1-14) ...
Setting up intel-oneapi-compiler-shared-runtime-2025.0 (2025.0.4-1519) ...
Setting up intel-oneapi-compiler-shared-2025.0 (2025.0.4-1519) ...
Setting up intel-oneapi-compiler-fortran-common-2025.0 (2025.0.4-1519) ...
Setting up intel-oneapi-mpi-devel (2021.14.1-5) ...
Setting up intel-oneapi-tbb-devel (2022.0.0-402) ...
Setting up intel-oneapi-mkl-core-2025.0 (2025.0.1-14) ...
Setting up intel-oneapi-compiler-dpcpp-cpp-common-2025.0 (2025.0.4-1519) ...
Setting up intel-oneapi-mkl-cluster-2025.0 (2025.0.1-14) ...
Setting up intel-pti-dev (0.10.0-284) ...
Setting up intel-oneapi-mkl-core-devel-2025.0 (2025.0.1-14) ...
Setting up intel-oneapi-ccl-devel-2021.14 (2021.14.0-505) ...
Setting up intel-oneapi-ipp-2022.0 (2022.0.0-808) ...
Setting up intel-oneapi-ccl-devel (2021.14.0-505) ...
Setting up intel-oneapi-compiler-dpcpp-cpp-runtime-2025.0 (2025.0.4-1519) ...
Setting up intel-oneapi-mkl-sycl-rng-2025.0 (2025.0.1-14) ...
Setting up intel-oneapi-compiler-fortran-runtime-2025.0 (2025.0.4-1519) ...
Setting up intel-oneapi-mkl-sycl-data-fitting-2025.0 (2025.0.1-14) ...
Setting up intel-oneapi-compiler-fortran-2025.0 (2025.0.4-1519) ...
Setting up intel-oneapi-mkl-sycl-vm-2025.0 (2025.0.1-14) ...
Setting up intel-oneapi-dal-2025.0 (2025.0.1-9) ...
Setting up intel-oneapi-dnnl-2025.0 (2025.0.1-6) ...
Setting up intel-oneapi-ipp-devel-2022.0 (2022.0.0-808) ...
Setting up intel-oneapi-mkl-sycl-dft-2025.0 (2025.0.1-14) ...
Setting up intel-oneapi-mkl-sycl-blas-2025.0 (2025.0.1-14) ...
Setting up intel-oneapi-mkl-cluster-devel-2025.0 (2025.0.1-14) ...
Setting up intel-oneapi-dpcpp-cpp-2025.0 (2025.0.4-1519) ...
Setting up intel-oneapi-compiler-dpcpp-cpp-2025.0 (2025.0.4-1519) ...
Setting up intel-oneapi-mkl-sycl-stats-2025.0 (2025.0.1-14) ...
Setting up intel-oneapi-ipp-devel (2022.0.0-808) ...
Setting up intel-oneapi-mkl-sycl-lapack-2025.0 (2025.0.1-14) ...
Setting up intel-oneapi-compiler-dpcpp-cpp (2025.0.4-1519) ...
Setting up intel-oneapi-mkl-sycl-sparse-2025.0 (2025.0.1-14) ...
Setting up intel-oneapi-mkl-classic-devel-2025.0 (2025.0.1-14) ...
Setting up intel-oneapi-dnnl (2025.0.1-6) ...
Setting up intel-oneapi-mkl-sycl-2025.0 (2025.0.1-14) ...
Setting up intel-oneapi-dnnl-devel-2025.0 (2025.0.1-6) ...
Setting up intel-oneapi-dnnl-devel (2025.0.1-6) ...
Setting up intel-oneapi-mkl-sycl-devel-2025.0 (2025.0.1-14) ...
Setting up intel-oneapi-mkl-devel-2025.0 (2025.0.1-14) ...
Setting up intel-oneapi-dal-devel-2025.0 (2025.0.1-9) ...
Setting up intel-oneapi-dal-devel (2025.0.1-9) ...
Setting up intel-oneapi-mkl-devel (2025.0.1-14) ...
Setting up intel-oneapi-hpc-toolkit (2025.0.1-46) ...
Setting up intel-hpckit (2025.0.1-46) ...

```
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 oneAPI environment ...
  bash: BASH_VERSION = 5.2.21(1)-release
  args: Using "$@" for setvars.sh arguments: intel64
:: advisor -- latest
:: ccl -- latest
:: compiler -- latest
:: dal -- latest
:: debugger -- latest
:: dev-utilities -- latest
:: dnnl -- latest
:: dpcpp-ct -- latest
:: dpl -- latest
:: ipp -- latest
:: ippcp -- latest
:: mkl -- latest
:: mpi -- latest
:: pti -- latest
:: tbb -- latest
:: umf -- latest
:: vtune -- latest
:: oneAPI environment initialized ::
username@computername:~$ ifx -v
ifx version 2025.0.4
username@computername:~$ icx -v
Intel(R) oneAPI DPC++/C++ Compiler 2025.0.4 (2025.0.4.20241205)
...
username@computername:~$ mpiifx -v
mpiifx for the Intel(R) MPI Library 2021.14 for Linux*
Copyright Intel Corporation.
ifx version 2025.0.4
```

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
=====
```

```
...
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/:
```

```
fftw3.f  fftw3.h  fftw3l-mpi.f03  fftw3-mpi.h  
fftw3.f03  fftw3l.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:~/elpa-release_2024_05_001$ cd ~
```

```
username@computername:~$ wget
```

```
https://gitlab.mpcdf.mpg.de/elpa/elpa/-/archive/release_2024_05_001/elpa-  
release_2024_05_001.tar.bz2
```

```
...
```

```
username@computername:~$ tar xvf elpa-release_2024_05_001.tar.bz2
```

```
...
```

```
username@computername:~$ cd elpa-release_2024_05_001/
```

```
username@computername:~/elpa-release_2024_05_001$ ./autogen.sh
```

```
...
```

```
username@computername:~/elpa-release_2024_05_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_FCFLAGS="-L$MKLROOT/lib/intel64  
-lmkl_scalapack_lp64 -lmkl_intel_lp64 -lmkl_sequential -lmkl_core -lmkl_blacs_intelmpi_lp64  
-lpthread -lm -I$MKLROOT/include/intel64/lp64" --prefix=$HOME/elpa-release_2024_05_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: 20241105

This version of ELPA support the minimal autotuning version: 20171201

The current autotune version is: 20241105

...

```
username@computername:~/elpa-release_2024_05_001$ make
```

...

```
username@computername:~/elpa-release_2024_05_001$ make install
```

...

```
username@computername:~/elpa-release_2024_05_001$ ls ~/elpa-release_2024_05_001/lib
```

```
~/elpa-release_2024_05_001/include/elpa-2024.05.001/elpa
```

```
/home/username/elpa-release_2024_05_001/include/elpa-2024.05.001/elpa:
```

```
elpa_configured_options.h elpa_generated_c_api.h elpa.h
```

```
elpa_constants.h          elpa_generated.h      elpa_simd_constants.h
```

```
elpa_explicit_name.h      elpa_generic.h        elpa_version.h
```

```
/home/username/elpa-release_2024_05_001/lib:
```

```
libelpa.a libelpa.la libelpa.so libelpa.so.19 libelpa.so.19.3.0 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:~/fftw-3.3.10$ 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/gsabowien2k-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):

Y

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/src/.libs/libxc.a
/home/username/libxc-7.0.0/lib/libxc.a
2
...
    Press RETURN to continue
...
    Do you want to automatically search for FFTW installations? (Y,n):
Y
    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
...
    Do you want to use a FFTW version from the list above? (Y,n):
Y
    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
...
    Selection: 0
        Compiler options=-O0 -FR -mp1 -w -prec_div -pc80 -pad -ip -DINTEL_VML -traceback
        -assume buffered_io -I$(MKLROOT)/include
...
    Current settings:
    M  OpenMP switch:      -qopenmp
    O  Compiler options:   -O0 -FR -mp1 -w -prec_div -pc80 -pad -ip -DINTEL_VML
    -traceback -assume buffered_io -I$(MKLROOT)/include
    L  Linker Flags:       $(FOPT) -L$(MKLROOT)/lib/$(MKL_TARGET_ARCH) -lpthread
    -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):
y
...
Do you want to automatically search for ELPA installations? (Y,n):
Y
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-release_2024_05_001
Finding the required ELPA files in /home/username/elpa-release_2024_05_001 ....

More than one version of ELPA found:
/home/username/elpa-release_2024_05_001/.libs
/home/username/elpa-release_2024_05_001/lib

Pick one (enter line number) or enter 0 to manually specify a path!:
2
Present root directory of ELPA is: /home/username/elpa-release_2024_05_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-release_2024_05_001/src/elpa1
/home/username/elpa-release_2024_05_001/src/elpa2
/home/username/elpa-release_2024_05_001/src/elpa_generalized

```

/home/username/elpa-release_2024_05_001/.fortran_dependencies/elpa2_print_kernels
/home/username/elpa-release_2024_05_001/test/Fortran/elpa2
/home/username/elpa-release_2024_05_001/test/Fortran/elpa_generalized
/home/username/elpa-release_2024_05_001/elpa
/home/username/elpa-release_2024_05_001/share/doc/elpa
/home/username/elpa-release_2024_05_001/include/elpa-2024.05.001
/home/username/elpa-release_2024_05_001/include/elpa-2024.05.001/elpa
9

...
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

...
These options derive from your chosen Settings:

ELPAROOT: /home/username/elpa-release_2024_05_001/
ELPA_VERSION: 2024.05.001
ELPA_LIB: lib
ELPA_LIBNAME: elpa
Is this correct? (Y,n): Y

Press RETURN to continue

...
Please specify your parallel compiler options or accept the recommendations (Enter - default)!:
-O0 -FR -mp1 -w -prec_div -pc80 -pad -ip -DINTEL_VML -traceback -assume buffered_io -I\$(MKLROOT)/include

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-release_2024_05_001/include/elpa-2024.05.001/elpa
-I/home/username/elpa-release_2024_05_001/include/elpa-2024.05.001/modules
ELPA_LIBS              : -lelpa -L/home/username/elpa-release_2024_05_001/lib -Wl,-
rpath=/home/username/elpa-release_2024_05_001/lib
FPOPT(par.comp.options): -O0 -FR -mp1 -w -prec_div -pc80 -pad -ip -DINTEL_VML
-traceback -assume buffered_io -I$(MKLRROOT)/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 preferred editor (default is emacs):
editor shall be: gedit

Set editor to gedit (Y/n) Y

Specify your preferred 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 preferred 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 !

added by WIEN2k: BEGIN

```
alias lsi="ls -aslp *.in*"
alias lso="ls -aslp *.output*"
alias lsd="ls -aslp *.def"
alias lsc="ls -aslp *.clm*"
alias lss="ls -aslp *.scf* */*.scf"
alias lse="ls -aslp *.error"
alias LS="ls -alsp |grep /"
alias pslapw="ps -ef |grep "lapw""
```

```

alias cdw="cd /home/username/wiendata"
if [ "$OMP_NUM_THREADS" = "" ]; then export OMP_NUM_THREADS=1; fi
#export LD_LIBRARY_PATH=.....
export EDITOR="gedit"
export SCRATCH=./
if [ "$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](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