

# DarkSUSY 6

## Tutorial

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# DarkSUSY 6 pre-release I

- This is a pre-release. Compared to the final DS 6.0 version (expected later this fall):
  - It is not fully finalized regarding SLHA reading/writing.
  - It is not fully tidied up and commented (output statements, main programs e.g.)
  - It does not have a completely updated manual
  - It is not finalized regarding charged cosmic ray diffusion and the interface to different halo models.
  - We have not yet tested on all compilers. gfortran 5 and 6 should work.
- If you find problems/have questions, e-mail [edsjo@fysik.su.se](mailto:edsjo@fysik.su.se)

Please don't distribute this pre-release version further at this point!

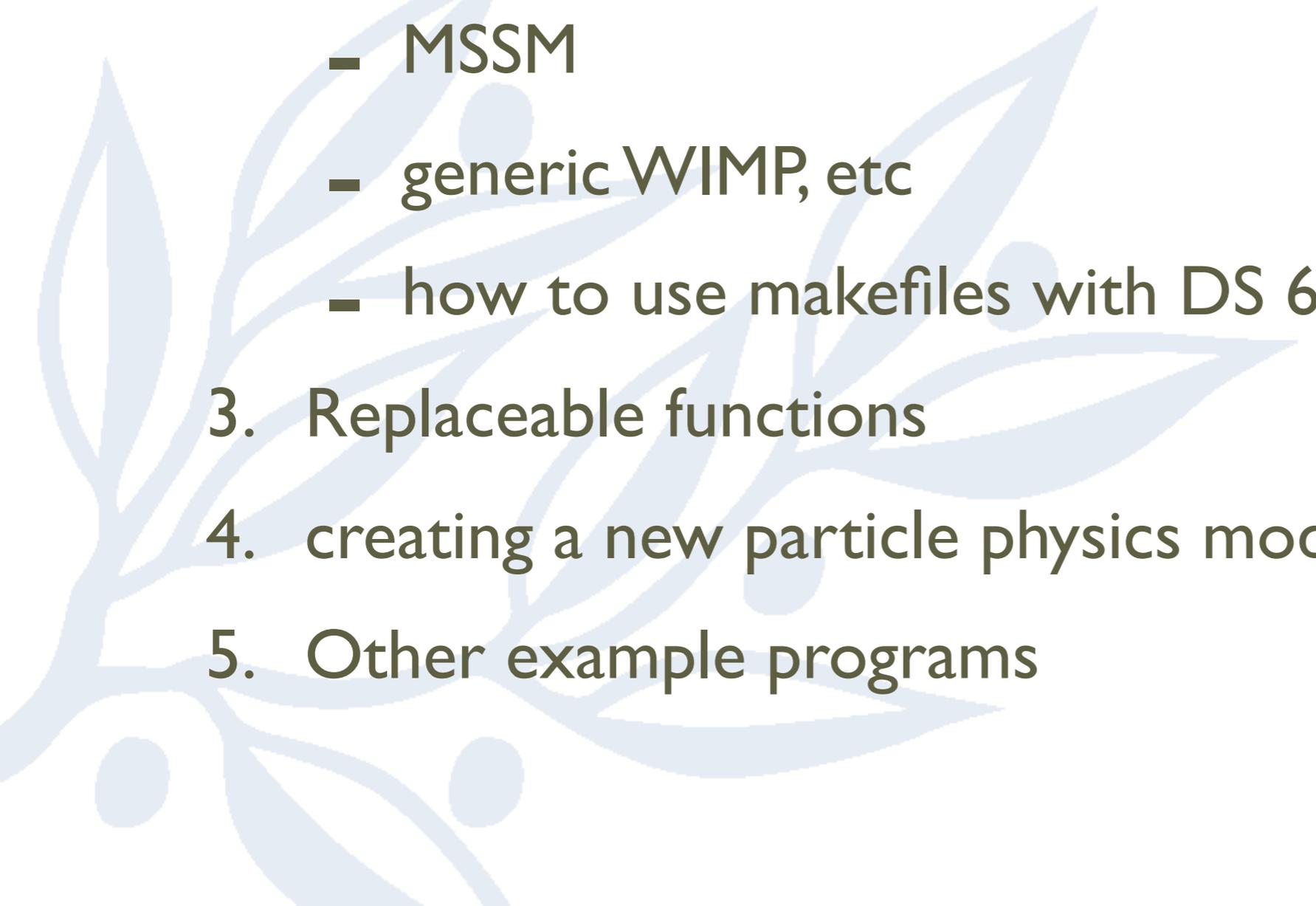
# DarkSUSY 6 pre-release 1

- Download: [www.astroparticle.se/ds/](http://www.astroparticle.se/ds/)
- Unpack it: `tar zxvf darksusy-6.0-pre1.tar.gz`
- Replace `examples/dsmain_wimp.F` with the version on the web page above

```
./configure  
make
```

Please don't distribute this pre-release version further at this point!

# Outline of hands-on

1. dstest program
  2. dsmain\_wimp program
    - MSSM
    - generic WIMP, etc
    - how to use makefiles with DS 6
  3. Replaceable functions
  4. creating a new particle physics module
  5. Other example programs
- 

# I. dstest program

- The dstest program is used to test your installation
- It calculated observables (masses, relic density, direct and indirect rates, ...) and compares with pre-computed values

```
cd examples/test  
./dstest
```

(already compiled with main make, takes about 60 seconds to run)

Output should end with

```
Total number of errors in dstest:      0
```

[Show code]

## 2. dsmain\_wimp.F

- In examples/ we have the file dsmain\_wimp.F which essentially does what dstest does, but in a more user-friendly way.
- run it with  
`./dsmain_wimp`
- It will ask you which model you want to run:

```
What kind of SUSY model do you want to look at?  
1 = MSSM-7  
2 = cMSSM  
3 = as read from an SLHA2 file
```

# MSSM-7 example

- Pick 1: MSSM-7 and enter (e.g.)  
mu: 1000  
M2: 1000  
MA: 400  
tan( $\beta$ ): 10  
m0: 3000  
At/m0: 0  
Ab/m0: 0
- Then answer 0 to not write out an SLHA file (or something else if you want to)
- Observables are then calculated...

# Output (cut)

```
Calculating omega h^2 without coannihilations, please be patient...
  without coannihilations Oh2 = 0.96585250586039517 0
0
Calculating omega h^2 with coannihilations, please be patient...
  with coannihilations Oh2 = 0.96585250586039517 0
0
  Chemical decoupling (freeze-out) occurred at
  T_f = 22.878440648494614 GeV.

Kinetic decoupling temperature, Tkd = 216.93665213661242 MeV
  The resulting cutoff in the power spectrum corresponds to a mass of
M_cut/M_sun = 2.2908727364927531E-009

dsddset: unrecognized option 'si' 'best'
dsddset: unrecognized option 'sd' 'best'
Calculating DM-nucleon scattering cross sections...
  sigsip (pb) = 8.5855360125101907E-010
  sigsin (pb) = 8.9165540437856185E-010
  sigsdp (pb) = 1.9718211101071476E-007
  sigsdn (pb) = 1.4088315037835129E-007
```

etc

# Which module?

- At the end of the `dsmain_wimp` run we got

```
-----  
The DarkSUSY example program has finished successfully.  
Particle module that was used: MSSM  
-----
```

```
[simply call 'make -B dsmain_wimp DS_MODULE=<MY_MODULE>' if you want to try  
with a different module <MY_MODULE>]
```

- Try compiling again with

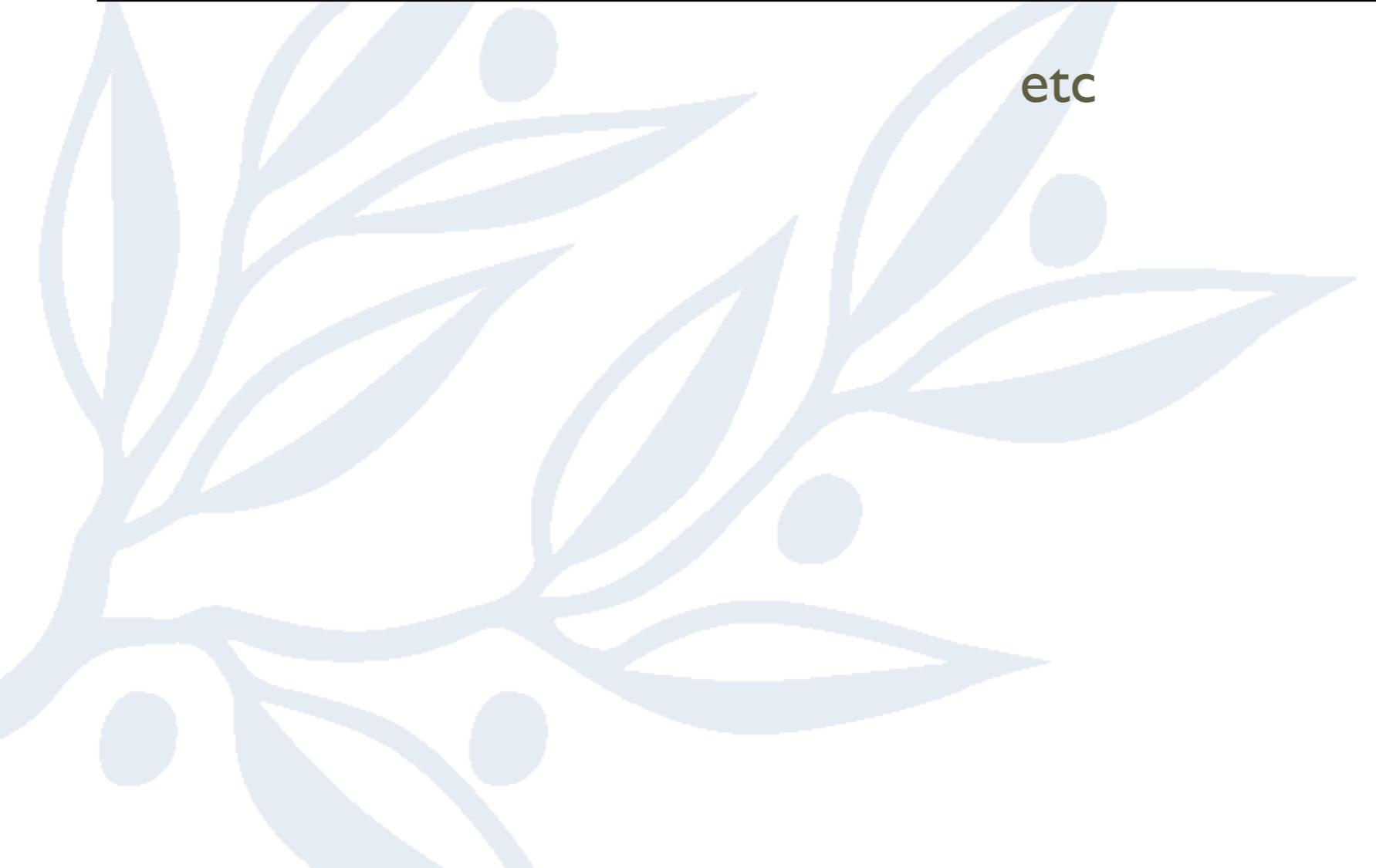
```
make -B dsmain_wimp DS_MODULE=generic_wimp  
./dsmain_wimp
```

- Enter e.g.  
mass: 100  
self-conjugate: 0  
ann cross section: 3e-26  
PDG: 5  
scattering cross section: 1e-42

# Output

```
Calculating omega h^2 without coannihilations, please be patient...  
without coannihilations Oh2 =      8.5782015186659649E-002      0      0  
Chemical decoupling (freeze-out) occurred at  
T_f =      4.4034841137539358      GeV.
```

etc



# Makefiles

- The way we choose which particle physics module to use is when we build our main program, e.g.

```
gfortran -o dsmain_wimp dsmain_wimp.F -lds_core.a -lds_mssm.a
```

- This can be made more flexible with makefiles,

```
dscheckmod :
    test `ls ../lib/ | grep libds_${DS_MODULE}.a` || { echo ERROR: Module $
{DS_MODULE} does not exist, or is not compiled; exit 1;}

dsmain_wimp : DS_MODULE = $(shell sed -n '1p' dsmain_wimp.driver)

dsmain_wimp : dscheckmod makefile dsmain_wimp.F
    printf "#define MODULE_CONFIG MODULE_"$(DS_MODULE)"\n" > module_compile.F
    printf "$(LIB)/libds_core_user.a\n"$(LIB)" /libds_core.a\n"$(LIB)" /libds_"$
(DS_MODULE)"_user.a\n"$(LIB)" /libds_"$(DS_MODULE)".a" > module_link.txt
    $(ADD_SCR) libds_tmp.a module_link.txt
    $(FF) $(FOPT) $(INC) $(INC_MSSM) -L$(LIB) -o dsmain_wimp dsmain_wimp.F \
libds_tmp.a $(shell if [ "x$(DS_MODULE)" = "xmssm" ]; then printf "%s" " $
(AUX_LIB_MSSM)"; fi)
    rm -f module_compile.F
    rm -f module_link.txt
    rm -f libds_tmp.a
```

# dsmain\_wimp.F

- dsmain\_wimp.F is a good starting point for your own program. If you want to use it as a starting point,
  - make a copy out of it
  - modify examples/makefile.in to copy-paste the lines about dsmain\_wimp.F and modify to your liking
  - run ./configure in the DS root
  - make and be happy!

# Some details of dsmain\_wimp.F

- In dsmain\_wimp we have code blocks of this type

```
#if MODULE_CONFIG == MODULE_generic_wimp
  subroutine dspnterparameters
    [more code for this module]
#endif
```

- This is how dsmain\_wimp.F performs model-specific setup.
- We could as well have prepared one separate main program for each particle physics module if we preferred (the makefile is then a bit simpler as well, see e.g. examples/aux/makefile)

# 3. Replaceable functions

- If you want to modify an existing DarkSUSY function or subroutine, **DON'T!**
- Instead create your own version of the routine and link to that one instead.
- You can either just create your own version and link to it (before the DS library is linked to), or
- Use the script `scr/make_replaceable.f` to make a `user_replaceable` function for you, for which the makefiles are already set up to work

# Replaceable function example

- As an example, we will look at the source term for DM annihilation in the galactic halo

$$\mathcal{S}_2(E_f) = \frac{1}{N_\chi m_\chi^2} \sum_i \sigma_i v \frac{dN_i}{dE_f},$$

This code is in `src_models/generic_wimp/dscrsource.f`

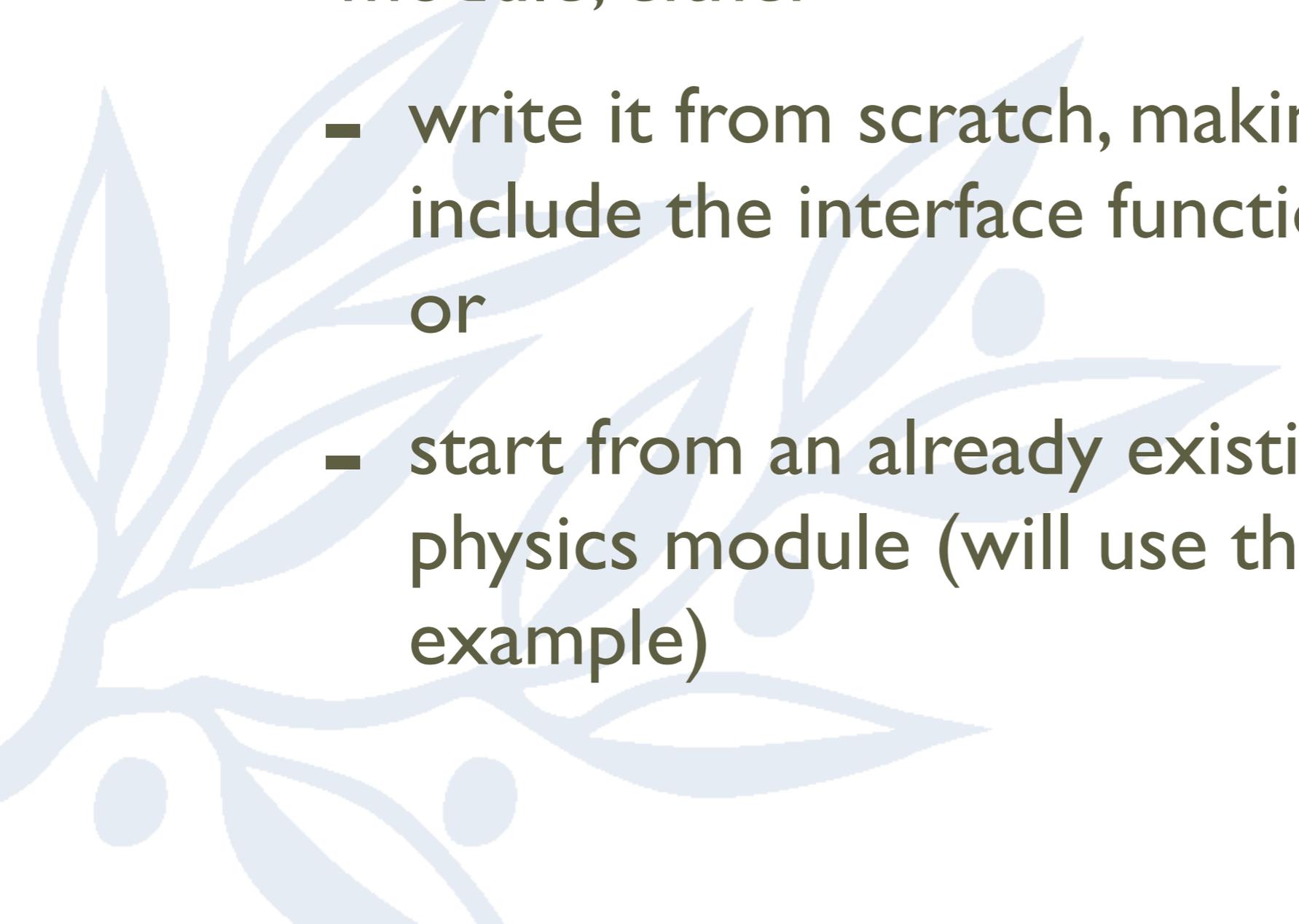
- Let's add a boost factor from substructures

$$\mathcal{S}_2(E_f) = \frac{1}{N_\chi m_\chi^2} \sum_i \sigma_i v \frac{dN_i}{dE_f} *B$$

# Replaceable function (cont)

- In the root directory, type  
`scr/make_replaceable.pl src_models/  
generic_wimp/cr/dscrsource.f`
- This will give you a new file  
`src_models/generic_wimp/  
user_replaceables/dscrsource.f`
- Modify it, configure and make again (in the  
root), then  
`make -B dsmain_wimp DS_MODULE=generic_wimp`  
in examples and run `dsmain_wimp`

# 4. Creating a new particle physics module

- To create a completely new particle physics module, either
    - write it from scratch, making sure to include the interface functions you need, or
    - start from an already existing particle physics module (will use this as an example)
- 

# Particle physics modules

- In `src_models` we currently have
  - `mssm` - Minimal Supersymmetric Standard Model
  - `silveira_zee` - Scalar singlet model
  - `generic_wimp` - a generic annihilating WIMP model
  - `generic_decayingDM` - a generic decaying dark matter model
  - `empty` - an empty model with just the basic set of interface functions for a 'fresh' start
- If you add one and want others to use it, please let us know and we can add it to the distribution (or point to your preferred download page)

# Simple example, extend generic wimp

- Create a new module by typing (in the root directory)

```
scr/make_module.pl generic_wimp extended_wimp
```

- Then type  

```
./configure  
make
```
- You then have a new module `extended_wimp` in `src_models`
- It is right now identical to `generic_wimp`, but you can now modify it to your liking

You need to have `autoconf` installed for this to work

# Helpful tools

- The `extended_wimp` is automatically included in the build system, but when/if you start adding files you need to tell the build system. To help you, we have a few scripts
  - `scr/makemf.pl <directory>` - adds all source files in the given `<directory>` to the relevant makefiles, or rather makefile.in's (without argument it adds source files in all directories in `src/` and `src_models`)
  - `scr/preconfig.pl` - adds source files AND new directories to the build system and updates both the configure script and makefiles

You need to have `autoconf` installed for this to work

# Main program

- You can e.g. use your new module with `dsmain_wimp` (or any other main program you choose)
- For `dsmain_wimp`, you need it to be aware of your new module by adding lines of this type:

```
#if MODULE_CONFIG == MODULE_extended_wimp  
[add your code here]  
#endif
```

This can be done by e.g. copy-pasting the corresponding `generic_wimp` lines and replace `generic_wimp` with `extended_wimp`

# 5. Other main programs

- In examples/aux we have a few example programs for other typical calculations, e.g.
    - the program to calculate the relic density in the Silveira Zee model
    - the program to calculate the relic density in the generic wimp model
  - we will add more examples and a description later
- will look at this code

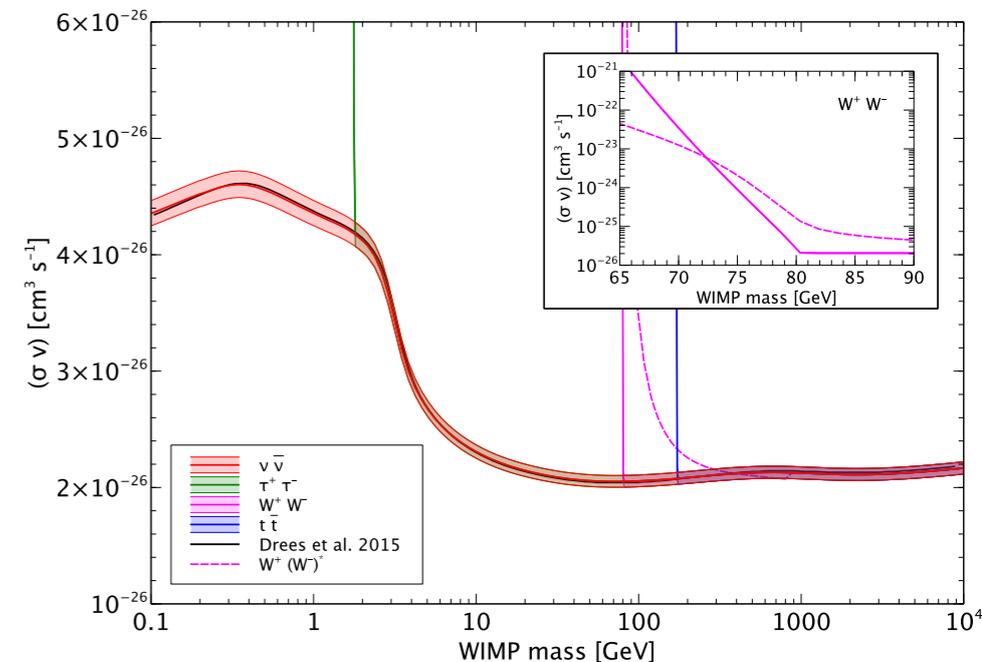
# generic\_wimp\_oh2

- This is the example program that creates the figure on relic density

```
cd examples/aux  
make generic_wimp_oh2  
./generic_wimp_oh2
```

Creates an output file `generic_wimp_oh2-planck-sigmav-thr.dat` that can e.g. be plotted

- It scans through the mass range, and for each mass makes a binary search in  $\sigma v$  to find the Planck measurement  $\pm 2$  sigma
- The default setup takes about 11.5 min to run, change 'f=1.1' to 'f=1.3' in line 40 and 'fth=1.02' to 'fth=1.1' on line 41 to speed it up for the tutorial (takes 3m20s on my laptop)



# Comment

- The default in `generic_wimp` is to use a sharp cut-off in  $W_{\text{eff}}$  when  $m_X < m_{\text{final}}$
- We can use an effective model with an off-shell final state particle, i.e.  $XX \rightarrow W^+ W^{*-}$
- An implementation of this is in `examples/aux/user_replaceables/dsanwx.f`
- Just compile replacing the regular `dsanwx.f` with this new one to test it:  
`make generic_wimp Oh2_threshold`

# Conclusions

- DarkSUSY 5 publically available
- DarkSUSY 6 is much more modular and include other improvements. Pre-release 1 available now. Expect full version later this fall
- When comparing different signals, it is crucial to perform these calculations in a consistent framework, with e.g. a tool like DarkSUSY

ευχαριστώ!



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