#### PALM group

#### Institute of Meteorology and Climatology, Leibniz Universität Hannover

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Flow chart





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- Machine dependencies





#### PALM Code: General Features





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- Data handling between subroutines is mostly done using FORTRAN90-modules instead of using parameter lists.



Most modules can be found in file .../trunk/SOURCE/modules.f90



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- The serial and parallel (MPI) PALM version is also activated by preprocessor directives.
- The code is splitted into several files, most of them containing just one subroutine, e.g. file "parin.f90" contains "SUBROUTINE parin".
- The code includes an interface which can be used to add your own code extensions. Advantage: These code extensions can be reused (normally) for future PALM releases without requiring any changes.





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# PALM Flow Chart (I)

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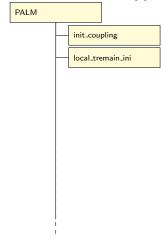
#### PALM Flow Chart (I)

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init\_coupling

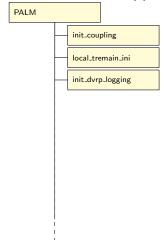






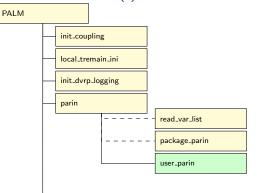






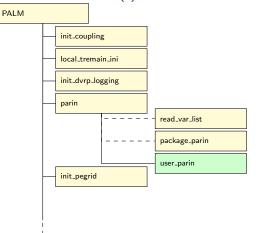








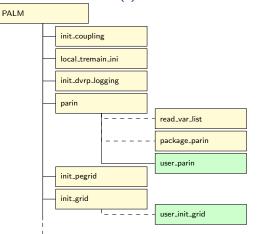








#### PALM Flow Chart (I)







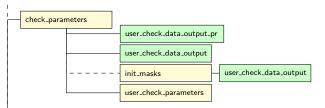
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## PALM Flow Chart (II)

check\_parameters



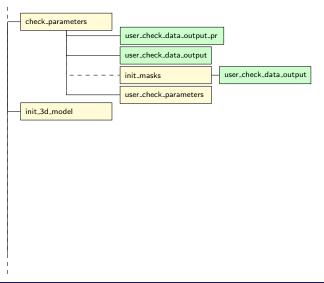








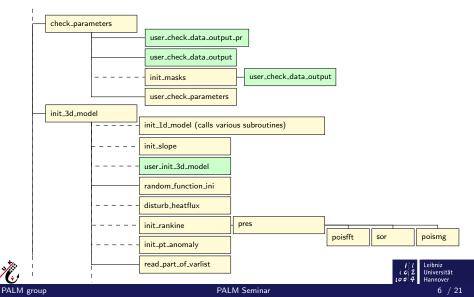
## PALM Flow Chart (II)



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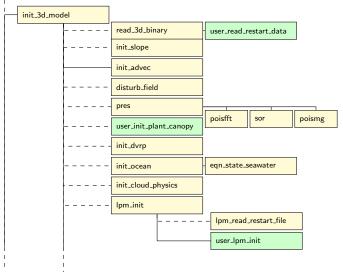
## PALM Flow Chart (III)

init\_3d\_model





## PALM Flow Chart (III)



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PALM Program Structure

## PALM Flow Chart (IV)





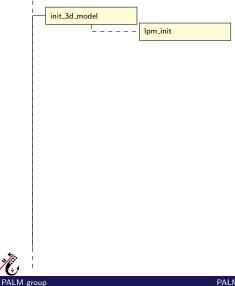
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## PALM Flow Chart (IV)

init\_3d\_model

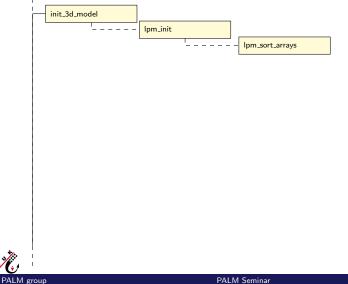








## PALM Flow Chart (IV)

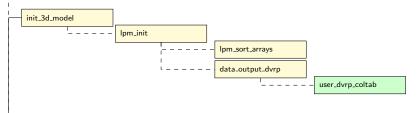


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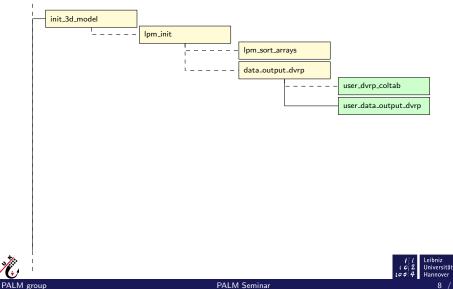


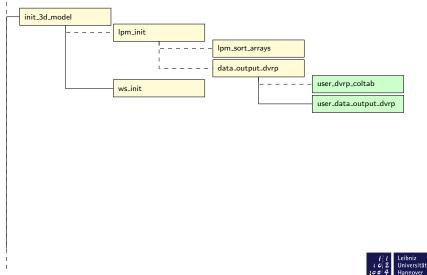


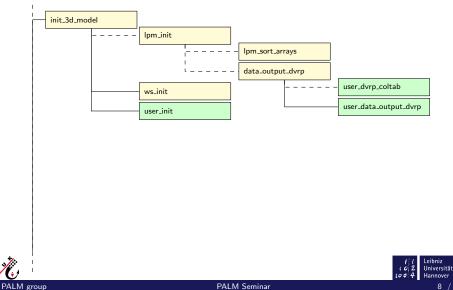


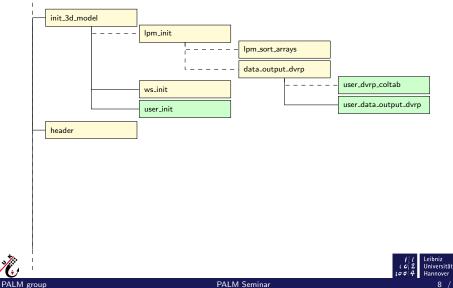


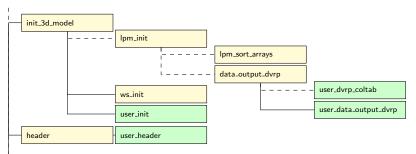






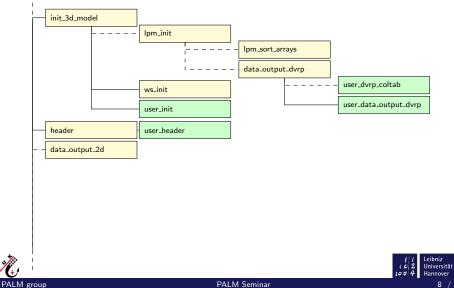




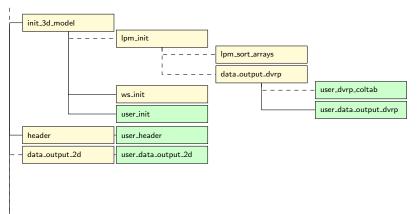








# PALM Flow Chart (IV)

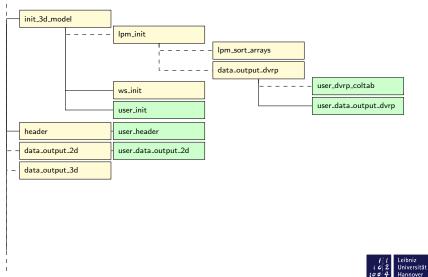






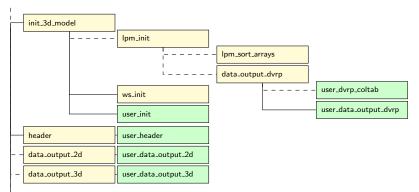
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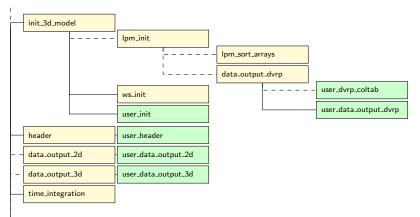


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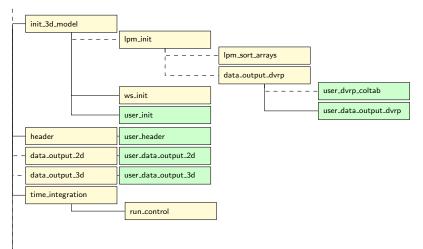








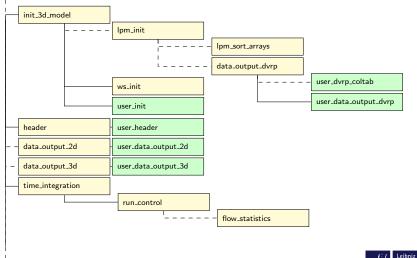








# PALM Flow Chart (IV)





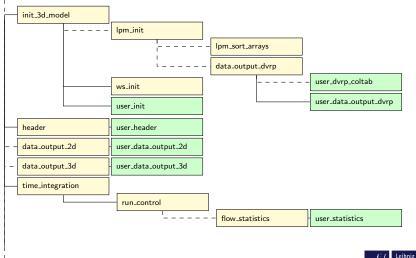
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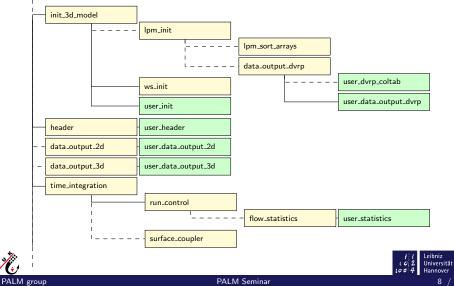


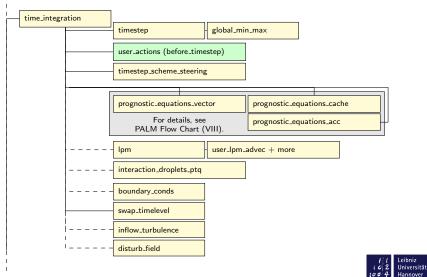


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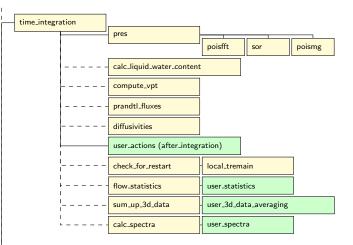
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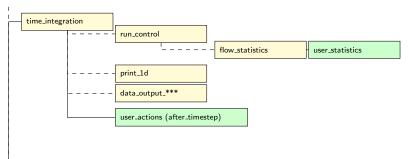
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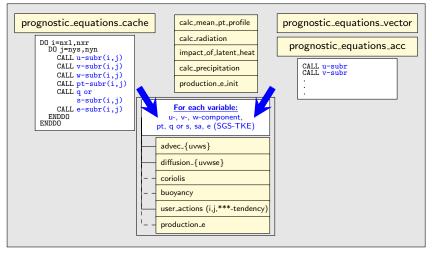




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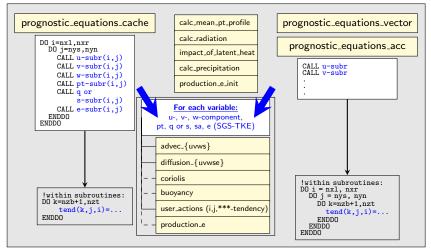








## PALM Flow Chart (VIII)





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	time_integration	
	i	
	write_3d_binary	write_var_list
$\vdash$	lpm_write_restart_file	
$\square$	header	user_header
$\square$	user_last_actions	
$\vdash$	cpu_statistics	
1		





#### Important Variables and Their Declaration





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3D-arrays of prognostic variables are named Ψ, and Ψ<sub>p</sub> for time level t, and t + Δt, respectively, with Ψ = u, v, w, pt, q, e, sa, u\_p, v\_p, ...





- 3D-arrays of prognostic variables are named Ψ, and Ψ<sub>p</sub> for time level t, and t + Δt, respectively, with Ψ = u, v, w, pt, q, e, sa, u\_p, v\_p, ...
- They are by default declared as  $\Psi(z,y,x)$  or  $\Psi(k,j,i)$ , e.g.

```
u(nzb:nzt+1,nysg:nyng,nxlg:nxrg)
```

#### with

nysg = nys - nbgp, nyng = nyn + nbgp nxlg = nxl - nbgp, nxrg = nxr + nbgp nzb, nzt (bottom, top) nys, nyn (south, north) nxl, nxr (left, right)

as the index limits of the (sub-)domain. nbgp is the number of ghost points which depends on the advection scheme (nbgp = 3 for the default Wicker-Skamarock scheme).



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nxl = 0; nxr = nx
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 For speed optimization, most of the 3D-variables are declared as pointers, e.g.

```
REAL(wp), DIMENSION(:,:,:), POINTER ::
u, u_p, v, v_p, ...
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► A pointer free version can be activated with preprocessor-option -D\_\_nopointer.





### Some Other Frequently Used Variables

variable	index bounds	meaning	comment
zu	nzb:nzt+1	heights of the scalar (u,v) grid levels	zu(0)=-zu(1)
ZW	nzb:nzt+1	heights of the w grid level	zw(0)=0
dzu	1:nzt+1	vertical grid spacings be- tween scalar grid levels	dzu(k)=zu(k)-zu(k-1)
ddzu	1:nzt+1	inverse of grid spacings	ddzu(k)=1.0/dzu(k)
dx		grid spacing along ×	
ddx		inverse of dx	ddx=1.0/dx
current_timestep_number		timestep counter	
simulated_time		simulated time in seconds	



## Preprocessor Directives (I)

 Preprocessor directives are special lines in the code which allows to compile alternative parts of the code depending on so-called **define-string** switches.

Code example:

```
#if defined( __nopointer )
REAL(wp), DIMENSION(:,:,:), ALLOCATABLE, TARGET :: e, e_p, ...
#else
REAL(wp), DIMENSION(:,:,:), POINTER :: e, e_p, ...
#endif
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#endif
```

If now the compiler is called e.g.

ifort -cpp -D \_\_nopointer ... (other options)

then the line containing "...,  $\tt allocatable, target :: ...."$  is compiled. If the compiler call is

```
ifort -cpp ... (other options)
```

the line containing "..., POINTER :: ..." is compiled.



## Preprocessor Directives (II)

The preprocessor directives require to include the compiler option "-cpp" in any way. Otherwise, the compilation will give error messages. The option has to be given in the configuration file .mrun.config in the %cpp\_options line. Different compilers may require different options!





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- Define-string switches can be combined using logical AND / OR operators && / ||.

```
#if defined ( __abc && __def )
```



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► To switch between the serial and the parallel code:

\_\_parallel

This switch is set by mrun-option "-K parallel".



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## Preprocessor Directives (IV)

In the PALM code, define-string switches are additionally used for following reasons:

- To enable usage of special software packages which are not included in the compilation process by default
  - \_\_dvrp\_graphics 3D visualization system (currently out of order)
  - \_\_spectra calculation and output of power spectra

Switches are activated with mrun-option "-p",

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- To enable special features
  - \_\_openacc activates call of external routines required for OpenACC programming
  - \_\_netcdf, \_\_netcdf4, \_\_netcdf\_parallel NetCDF I/O with different NetCDF versions



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## Preprocessor Directives (V)

Preprocessor directives are also used for string replacement in the code.

Example:

A compiler call with preprocessor option

```
ifort -cpp -Dabcd=efgh
```

will replace all strings "abcd" in the code with "efgh" before the code is compiled.

This is used in PALM to change the MPI\_REAL datatypes (which are 4 byte long by default), to 8 bytes. The respective cpp-directives are given in the configuration file .mrun.config.:



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