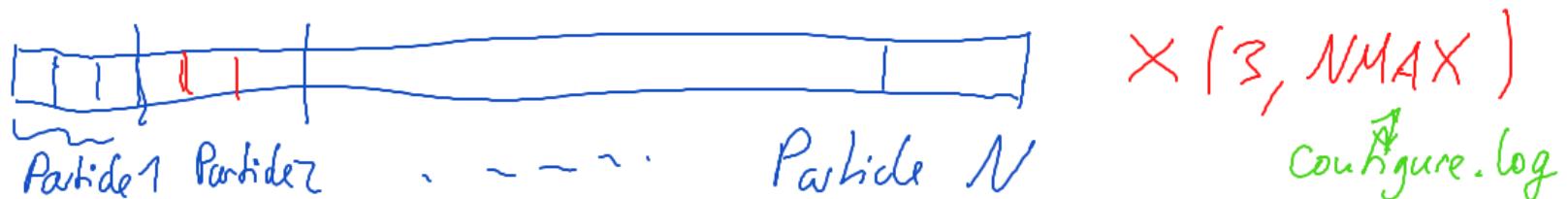


• 1) Data Structure / 2) Code Structure

1) old fashioned; Aarseth: single particles



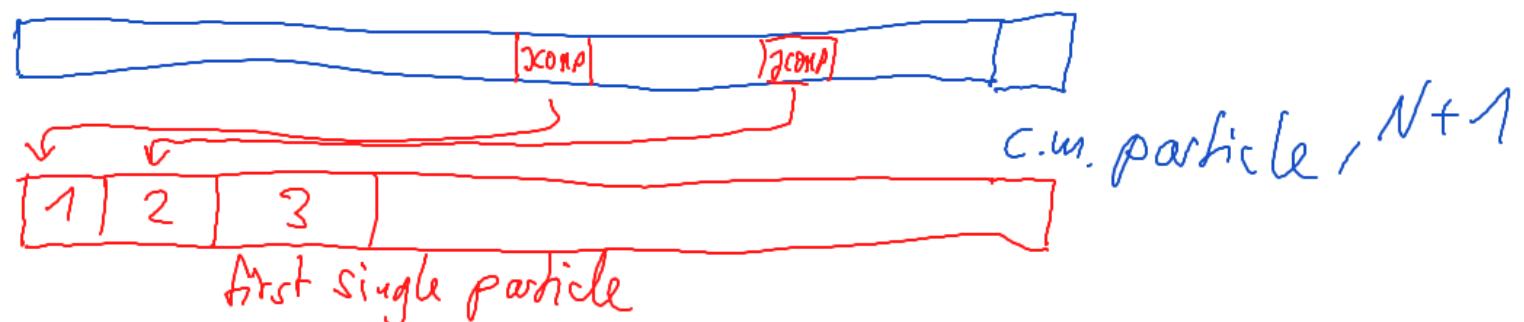
N is number of particles actually used, input par.

$$N \leq N_{MAX}, N < N_{MAX}$$

KS binary is doing:

i_{COMP}, j_{COMP}

NEW
KSREG:





binary first single

KTERM:



single particles

If 10.000 KS Lunes:



IFIRST = 20001

$$N_{TOT} = N + N_{PAIRS}$$

total elements used:

$$IFIRST = 2 \times N_{PAIRS} + 1$$

index of first single part,

find two kinds
of do loops:

DO . 1, N

all particles

DO .. IFIRST, NTOT

particles in Henrite

scheme

How do we find a particle? Variable NAME(I)
is defined for all outputs:

NAME(1) particle initially at pos. 1 ($t=0$)
(most massive particle)

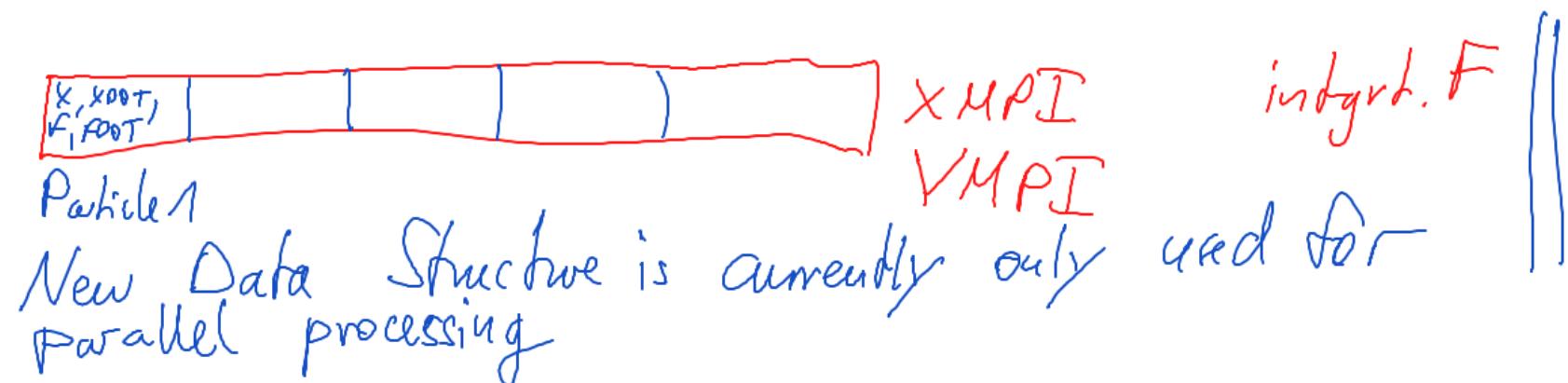
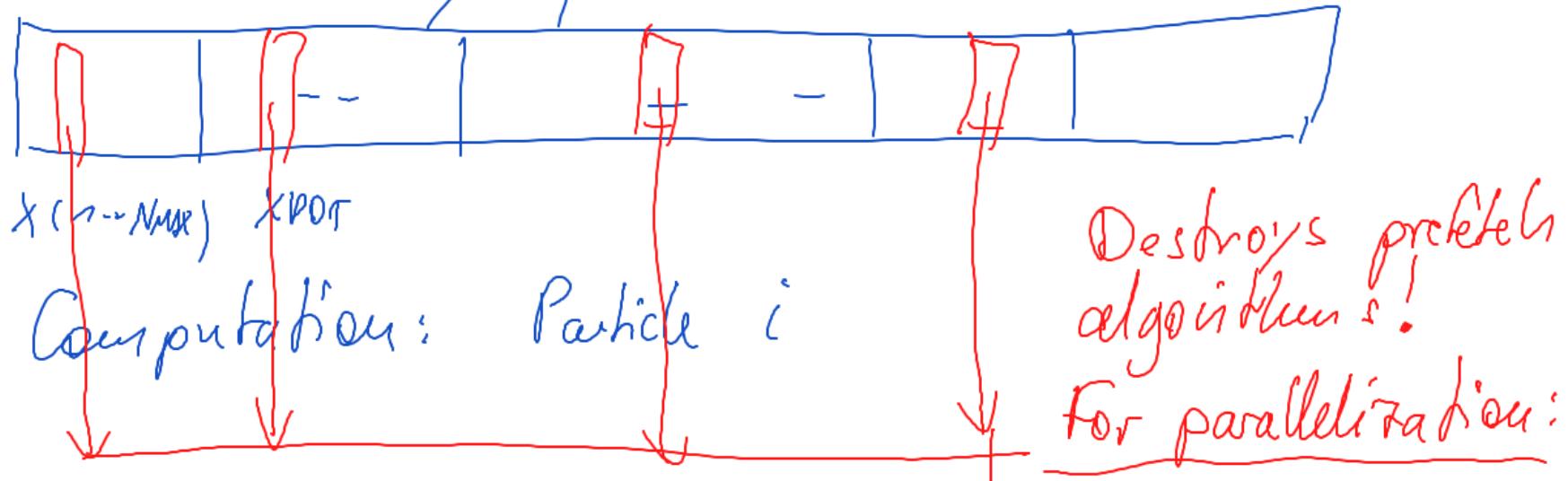
~~IF (I.EQ.1) ...~~

IF (NAME(I).EQ.1) ...

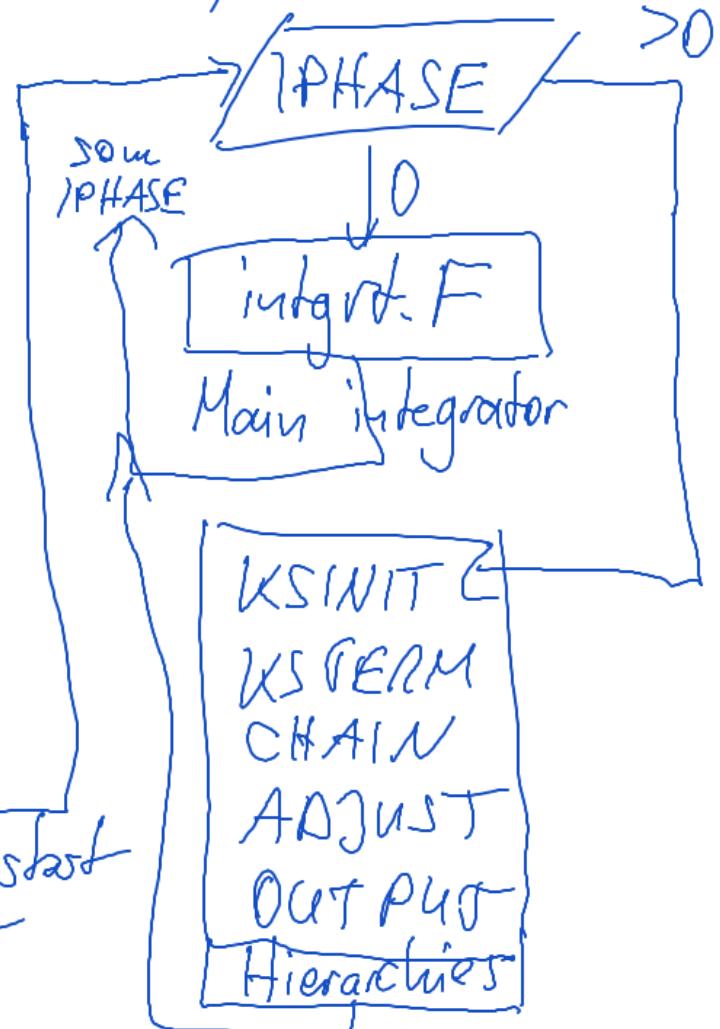
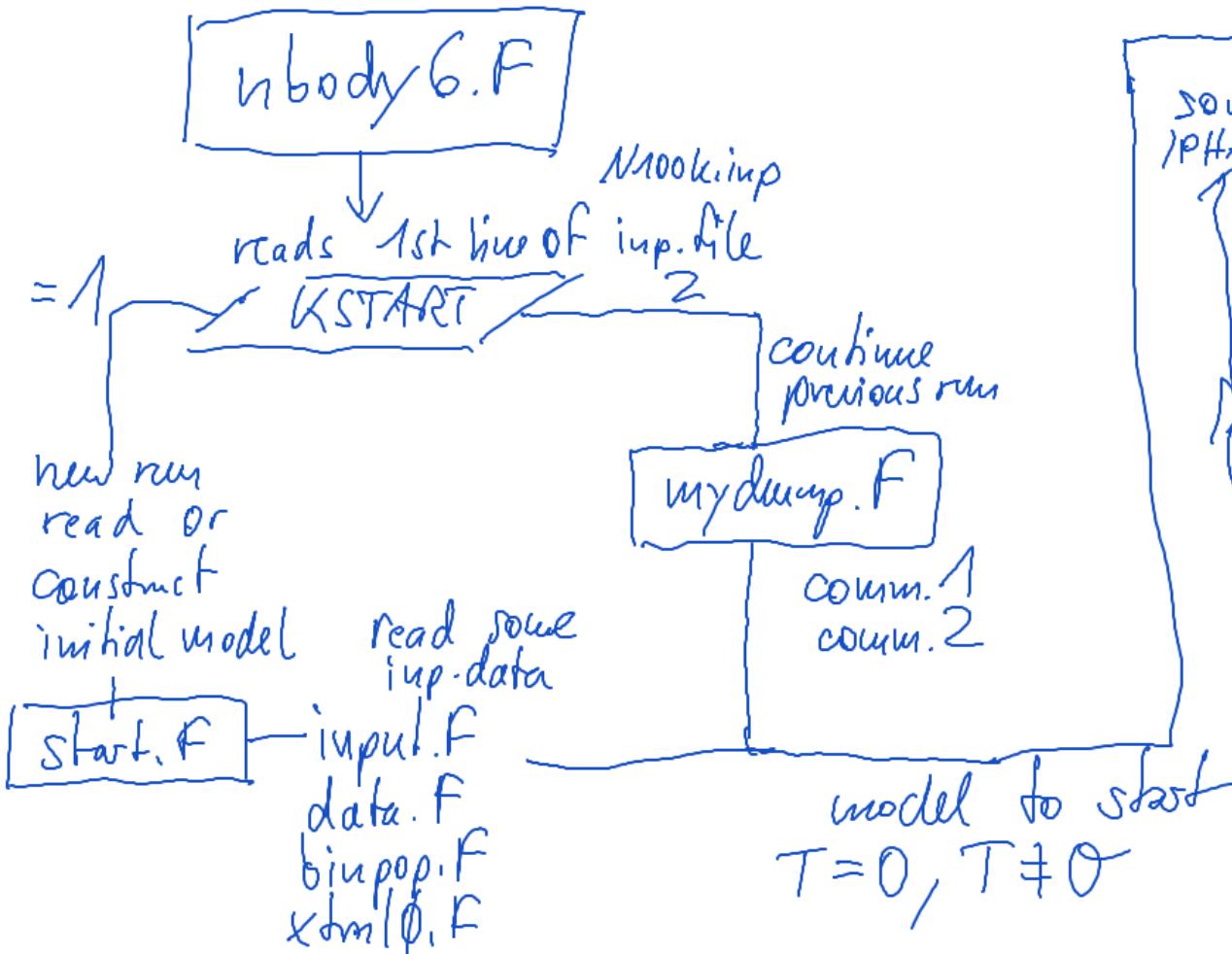
Fortran Style COMMON BLOCKS | include/common6.h

X, XDOT, X ϕ , X ϕ DOT	F1, FR, F1DOT, F1DOT
X _p V _p	last x _c last v _c
	a _{yr} a _{rg} a _{yr} ^e , a _{yr} ^o

Memory, not nice



2) Code Structure (compare Manual)



intgrt

$\rightarrow \text{TBLOCK} = \text{TIME}$

1) Advance KS pairs, chains

MP 2) Check for new KS pairs

④ 3) Do conversion step for irregular particles due Open MP Case

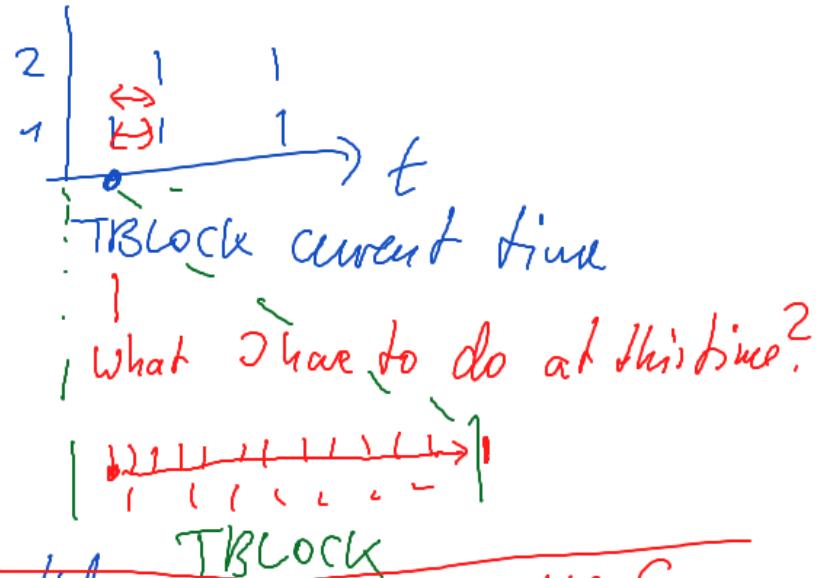
MP Are there any regular steps due?

④ 4) Do conversion step for regular particles due GPU

\rightarrow 5) Advance all stellar evolution / mass loss

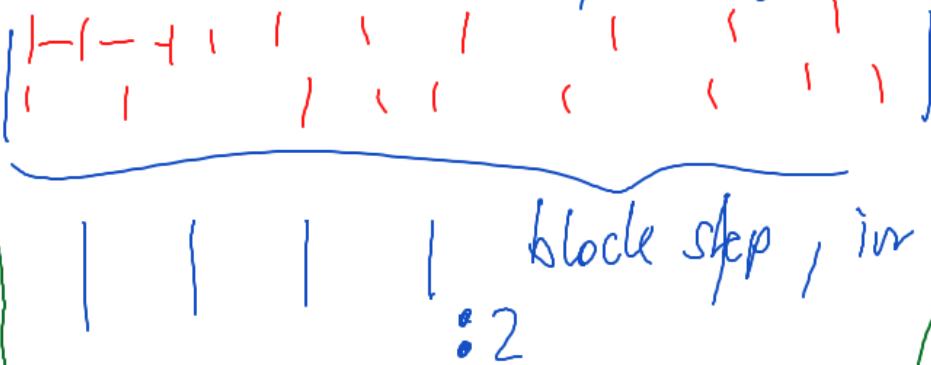
↓ Final Step: Check for Kstem, Adjust \rightarrow jump out with IPHASE (e.g. check)

$\text{TBLOCK} = \text{TBLOCK} + \Delta t$
(sorted list of particles in line, to find next particles for corr. step.)

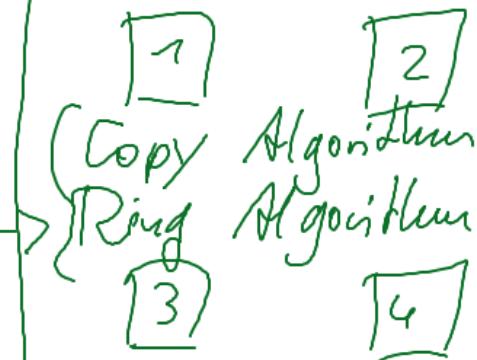


Things to do:

1) Binary parallelization, KS integration
(idea by Aavseth + Sp, Wang)



2) Domain Decomposition
Nodes



Full copy of all data; parallelization of working loops
many particles with small steps

Dorland, Heusendorf, Merritt
Makino 2003
 2002 (GRAPE)

integrate binaries in parallel on smaller block level