

Status of TM6

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Outline

Context

Previously in TM6

TM6 Status & Perf

Extra

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Extra

Performance issue

speed

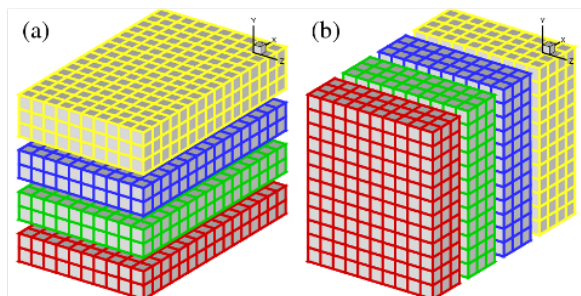
- ▶ fast as standalone, but slow for CGCM (EC-Earth)
 - ▶ decade/day wanted
 - ▶ BUT max nb processors = nb Tracers (**27, 1,..**)

resolution

- ▶ high resolution
 - ▶ very demanding in memory (10 Gb/proc @ 1x1)

Basic idea of MPI: domain decomposition

- ▶ Arrays are split across processors, along any dimension.



- ▶ TM5 **4D MASS** arrays are distributed along either **LEVELS** or **TRACERS**.

The bottleneck

meteo fields are **NOT** distributed... but **COPIED** !!!!

- ▶ every 3h => **FREQUENT** communication
- ▶ 50+ met fields
 - ▶ **HUGE** memory requirement
 - ▶ **HEAVY** communication



MPI profiling

TM5-chem v3, 2-days run, 4 MPI tasks

	% of elapsed time
switching decomposition	3 %
broadcasting meteo	50 %
other MPI comm	2 %
total MPI comm	55 %

Outline

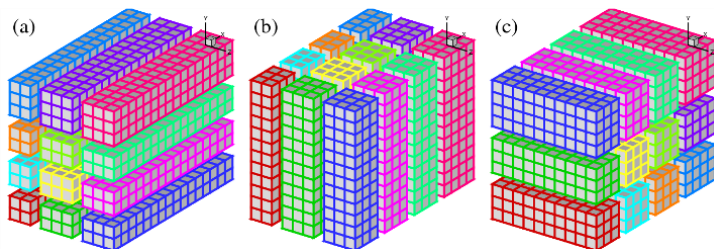
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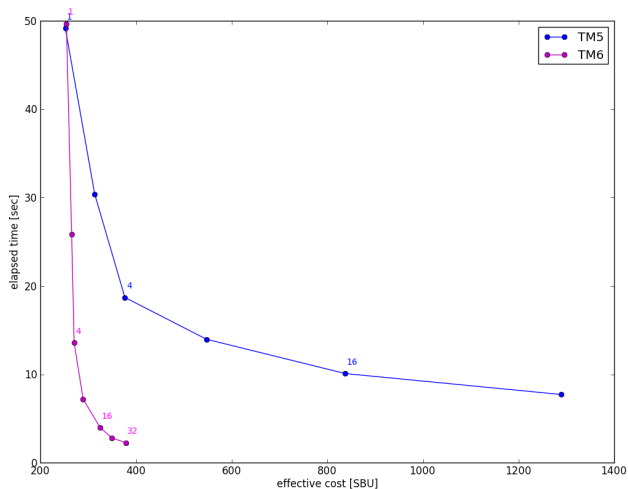
Extra

Revised domain decomposition = (b)



	TM5	TM6
max #processor	27	30x22 = 660 (@6x4) 60x45 = 2700 (@3x2) 180*90=16200 (@1x1)
meteo communication	broadcast	scatter

Performance TM-chemistry



**8x faster
same price**

ToDo list as of last Crete meeting

To test/fix

- ▶ ~~M7, online dust & outputs: mix, station, planeflight~~
- ▶ ~~debug : "1x8" case, "qfltrap=enable:inv" required (EBI)~~

To code & test

- ▶ ~~chunk reading of meteo in netCDF-4~~
- ▶ ~~aerocom & time-series outputs~~
- ▶ ~~EC-Earth proj~~
- ▶ ~~updated chem emissions (Edgar 4.2 + GFED3)~~

Missing features

~~reduced grid ; zoom regions~~

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Porting to ECMWF/c2a (IBM/AIX power7)

Fixed:

- ▶ Pure MPI-2 :
 - ▶ MPI_GET_EXTENT → MPI_TYPE_GET_EXTENT
 - ▶ MPI_TYPE_HVECTOR → MPI_TYPE_CREATE_HVECTOR
- ▶ libs
- ▶ totalview requires ssh

but still issues

- ▶ unexplained frozen runs
- ▶ M7 : crashes w/ 5+ cpus, sedimentation bug

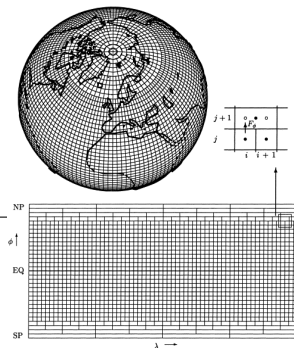
Reduced grid - implementation

first question at every talk!

- ▶ implement case of 'no decomposition along longitudes'

TM5	TM6	TM6 w/ redgrid
27	660 (@6x4)	22 (@6x4)
	2700 (@3x2)	45 (@3x2)
	16200 (@1x1)	90 (@1x1)

Max #processors



Reduced grid // 1-month runs // chemistry w/o M7

- ▶ 3x2 w/ reduced grid
 - ▶ 7 bands (90-74) at each pole
 - ▶ merging [40, 8, 8, 4, 4, 4, 2] cells



- ▶ TM5 → TM6 :
60-70 % speed-up
- ▶ Reduced grid :
30-40 % speed-up

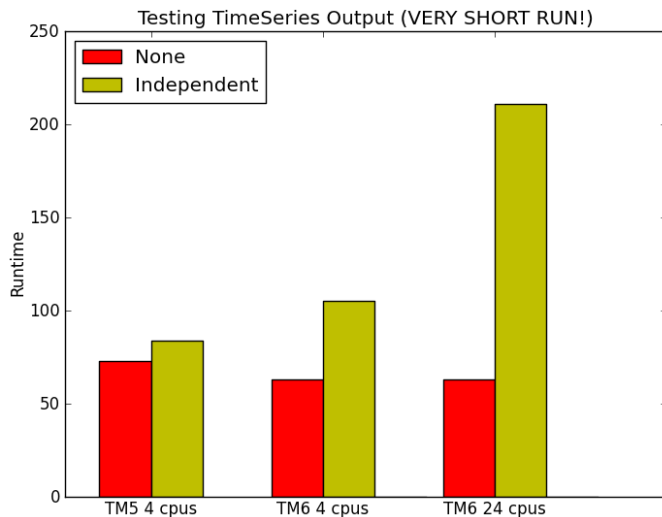
The I/O experiment (1) - Time-Series Output

(former RETRO output)

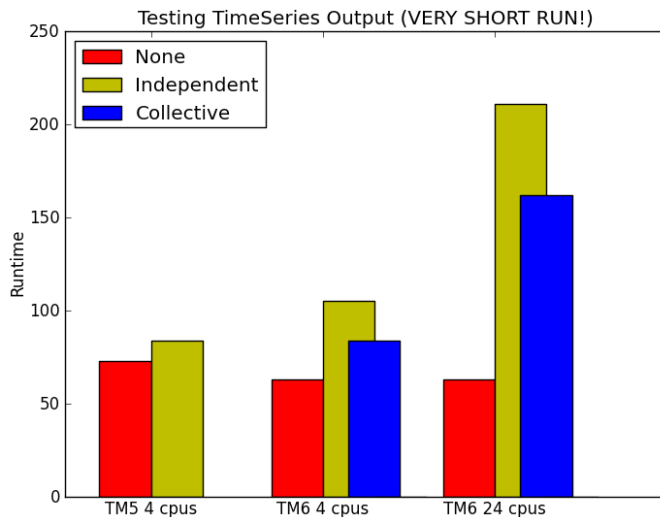
- ▶ TM5
 - ▶ pnetcdf → netcdf4 (MDF)
 - ▶ INDEPENDENT access mode <= **unlimited dimensions**

- ▶ TM6
 - ▶ case 1: stick to INDEPENDENT
 - ▶ case 2: switch to COLLECTIVE access mode
 - ▶ case 3: write once a day (not every time step!)

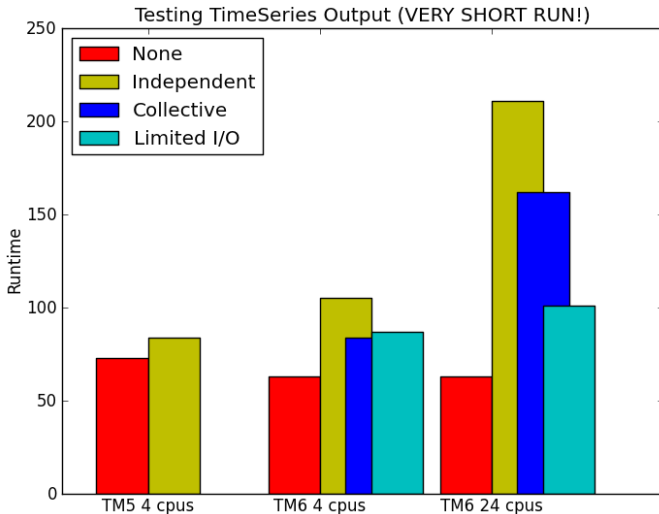
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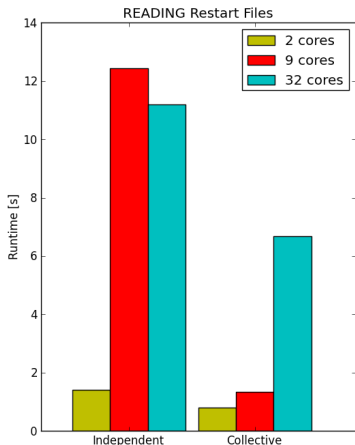
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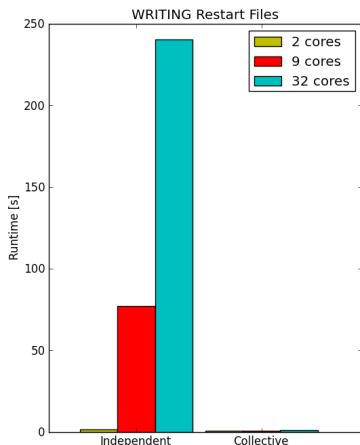
The I/O experiment (2) - Does netCDF4 w/ parallel I/O scale?



READING restart

- ▶ collective **faster** than independent (1.5-9x)
- ▶ **time increases w/ nb cores**
 - ▶ impact for meteo (**must** account for the scatter time saved)

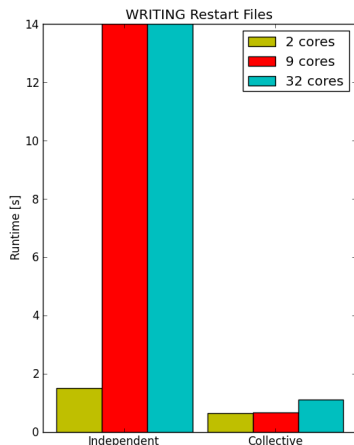
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WRITING restart

- ▶ collective **really faster**
(2.3-110x)

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WRITING restart

- ▶ collective **really faster** (2.3-110x)
- ▶ writing time : no increase with nb cores

I/O next steps - Optimize

- ▶ Time-Series Output
 - ▶ test w/ longer runs
 - ▶ one file /month & /tracer instead of /day & all tracers?
 - ▶ **quilting** : asynchronized I/O for MPI (eg, WRF)?
 - ▶ file splitting?
- ▶ Read/Write restart
 - ▶ file splitting
 - ▶ quilting
- ▶ Meteo Input
 - ▶ switch to parallel reading

"As-fast-as-you-can" experiment @1x1

one node (50 Gb, 32 procs max), one day sim, no reduced grid

Model	Regions	Resources	Runtime	Cost (SBU)
TM6	global 1x1	32 procs	1h 4mn	643
TM5	global 1x1	6 procs	11h 25mn	6850
TM5	global 3x2 + euro 1x1	<i>broken</i>	-	

- ▶ zooming broken in TM5 in 3 places:
 - ▶ when nudging of CH4 emissions
 - ▶ in photolysis
 - ▶ latitudinal decomposition
 - ▶ solar zenith angle

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SUCCESS!

- ▶ 10.6 x cheaper
- ▶ 10.6 x faster!
- ▶ 90.6% speedup



NEXT

- ▶ fix M7
- ▶ couple to EC-Earth
- ▶ optimize reduced grid
- ▶ optimize time-series
- ▶ read netCFD meteo in parallel

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Reduced grid - 1-month runs by numbers

- ▶ Chemistry w/o M7
 - ▶ optimized (-O3 -qstrict)
 - ▶ full chemistry (but w/o m7)
 - ▶ w/o time-series output
 - ▶ 3x2 w/ reduced grid:
 - ▶ 7 bands (90-74) at each pole
 - ▶ merging [40, 8, 8, 4, 4, 4, 2] cells

	w/o redgrid	w/ redgrid	speed-up
TM5	23799	14422	39%
TM6	7909	5401	32%
speed-up	67%	63%	77%

The I/O experiment (2) - by numbers

▶ Reading restart

	8x4 cpus	3x3 cpus	2x1 cpus
coll.	6.35	1.70	0.84
	7.00	0.99	0.75
ind.	10.99	12.30	1.33
	11.40	12.57	1.51

▶ Writing restart

	8x4 cpus	3x3 cpus	2x1 cpus
coll.	1.01	0.75	0.64
	1.23	0.59	0.65
ind.	237.19	73.28	1.51
	243.57	81.26	1.50