





National Grid Infrastructure (NGI)

for scientific computations, collaborative research & its support services

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National Grid Infrastructure (NGI)

National Grid Infrastructure (NGI)

- operated by MetaCentrum NGI (CESNET) since 1996
- MetaCentrum responsible for management and coordination

Distributed infrastructure

NGI integrates medium/large HW centers (clusters, powerful servers, storages) of several universities/institutions

 further integrated into the European Grid Infrastructure (EGI.eu) http://www.metacentrum.cz









Computing clusters

a set of interconnected ("common") computers



(oldschool)







Computing clusters

a set of interconnected ("common") computers



(nowadays)

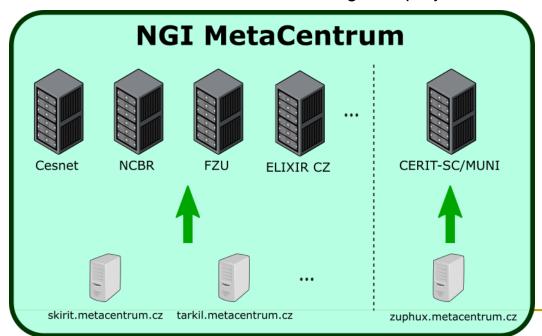






MetaCentrum NGI & Resource integration I.

- MetaCentrum and CERIT-SC
 - MetaCentrum provides <u>own HW resources</u> (CESNET) and <u>integrates resources of external providers</u>
 - CERIT-SC/MUNI is one of them.
 - others are CEITEC/NCBR, FZU, ČVUT, JČU, ZČU, UPOL, MU, TUL, etc.
 - as well as global projects like ELIXIR CZ



+ shared storages and shared SW apps







MetaCentrum NGI & Resource integration II.

- resource owners (usually) have priority access to their HW resources
 - under agreed conditions
 - technically accomplished using specific scheduler queues
 - more details later







MetaCentrum NGI

Available to all academic users from Czech universities, Academy of Science, research institutes, etc.

- commercial bodies just for public research

Offers:

- computing resources
- storage resources
- application programs

After registration, all the resources/services are available free of charge

- without any administrative burden
- users "pay" via publications with acknowledgements
 - → results in user priorities in cases of high load



http://metavo.metacentrum.cz







Meta NGI – basic characteristics

- after registration, all the resources are available without any administrative burden
 - $\rightarrow \sim$ immediately (depending on actual usage)
 - no applications for resources
- user accounts extensions every year
 - validates users' relationship to an academic institution
 - federated infrastructure edulD.cz used for minimalization of users' burden
 - reports of user publications with acknowledgements to MetaCentrum/CERIT-SC
 - used as a proof of infrastructure benefits for Czech research area
- best-effort service







Meta NGI – computing resources available

Computing resources: ca 17900 cores (x86_64)

- nodes with lower number of computing cores:
 2x4-8 cores
- nodes with medium number of comp. cores(SMP nodes): 32-80 cores
- memory (RAM) up to 1 TB per node
- nodes with high number of computing cores:SGI UV2000
 - 384 cores (x86_64), 6 TB of RAM
 - 288 cores (x86_64), 6 TB of RAM
- other "exotic" hardware:
 - nodes with GPU cards, SSD discs, Xeon Phi, etc.



http://metavo.metacentrum.cz/cs/state/hardware.html







Meta NGI – storage resources available

ca 5 PB for operational data

- centralized storage arrays distributed through various cities in the CR
- user quota 1-3 TB on each storage array

ca 22 PB for archival data

- (HSM MAID, tapes)
- "unlimited" user quota

http://metavo.metacentrum.cz/cs/state/nodes







Meta NGI – software available

- ~ 300 different applications (commercial & free/open s.)
 - see http://meta.cesnet.cz/wiki/Kategorie:Aplikace
- development tools
 - GNU, Intel, and PGI compilers, profiling and debugging tools (TotalView, Allinea), ...
- mathematical software
 - Matlab, Maple, Mathematica, gridMathematica, ...
- application chemistry
 - Gaussian 09, Gaussian-Linda, Gamess, Gromacs, ...
- material simulations
 - Wien2k, ANSYS Fluent CFD, Ansys Mechanical, Ansys HPC...
- structural biology, bioinformatics
 - CLC Genomics Workbench, Geneious, Turbomole, Molpro, ...







Meta NGI – grid environment

- batch jobs
 - the computations described by script files
- interactive jobs
 - text & graphical environment
- cloud computing
 - instead of running jobs with computations, users run the whole virtual machines

focused on research computations again (not for webhosting)
Windows & Linux images provided, user-uploaded images also supported
more info later...









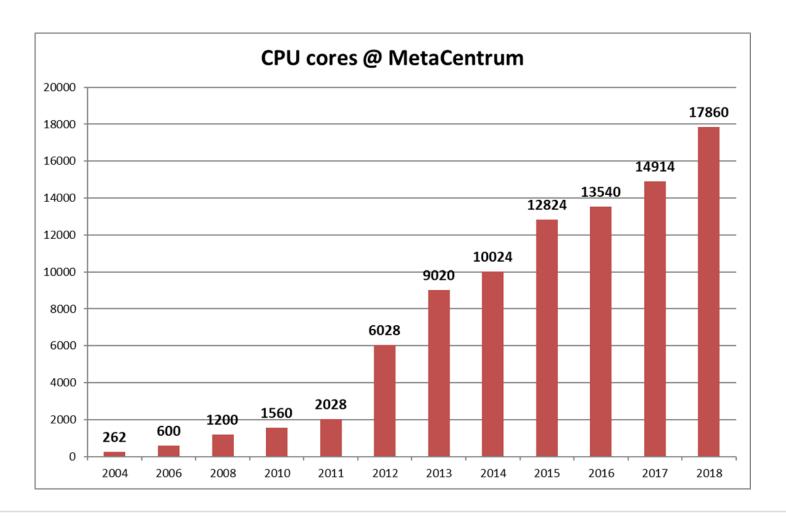
Meta VO in numbers...

- · ca 17860 cores, ca 600 server nodes
- year 2018:
 - **2020** users (31.12.2018)
 - ca 5 mil. of running jobs
 - ca 13700 jobs per day
 - · ca 2500 jobs per user
 - CPU time
 ca 11,4 thousands of CPUyears





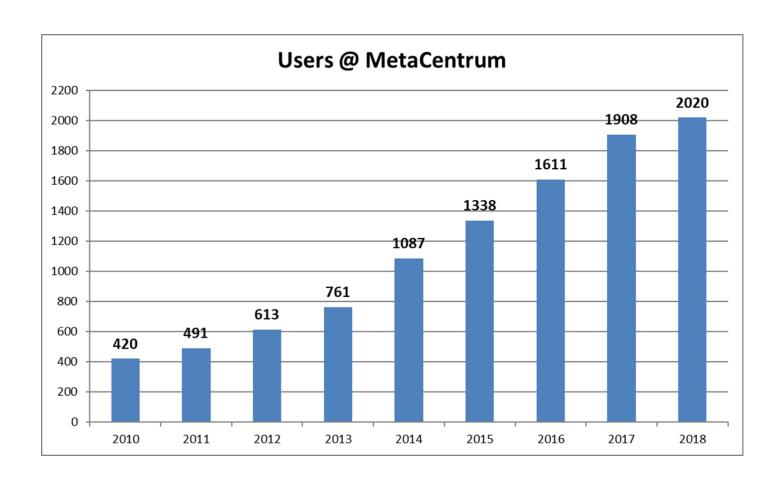








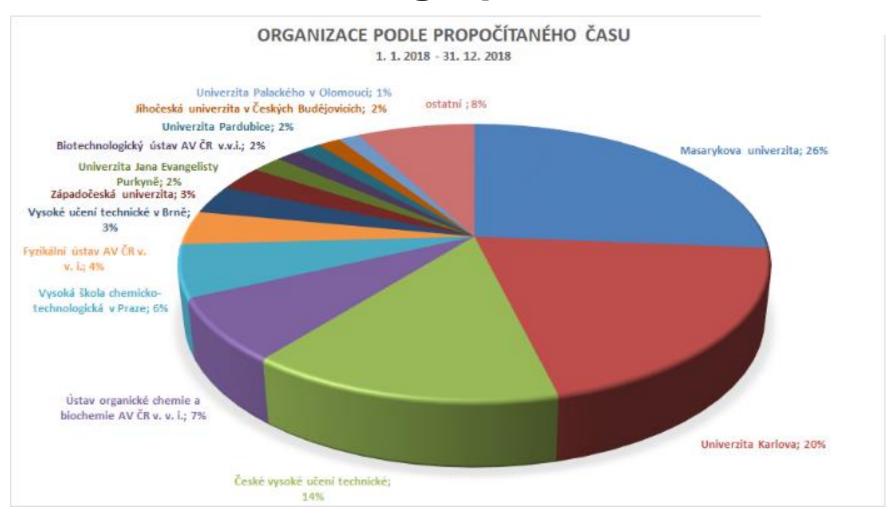








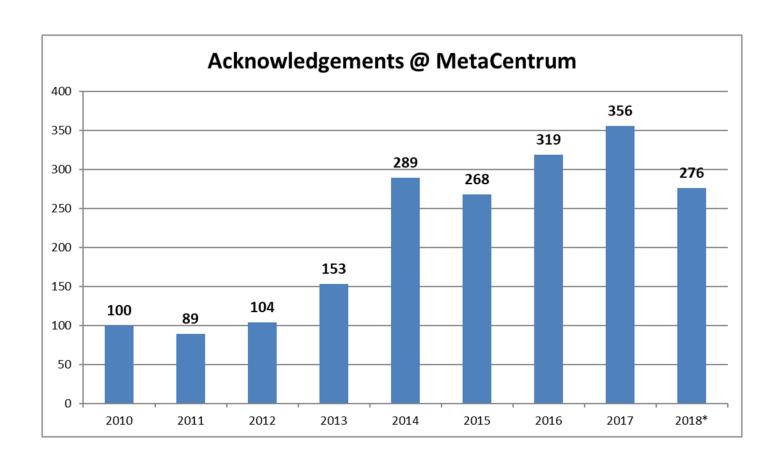


















Meta VO – how to become our user?

- register
 - http://metavo.metacentrum.cz , section "Application"
 - EduID.cz => proves your academic identity using your home institution services (and credentials)
- make yourself familiar with basics of OS Linux
 - http://metavo.metacentrum.cz , section "Documentation"
 - -http://lb.poznejlinux.cz/xhtml/linuxbook.html#id2571419
- compute







Research Infrastructures in CR I.

- IT4innovations (Ostrava)
 - 3344 computing cores ("small" supercomputer/cluster)
 - 24192 computing cores ("big" supercomputer/cluster)
 - attributes:
 - computing time assigned to **research projects**
 - formal application is necessary (evaluation of research and technical readiness + socio-economic factors)
 - public competitions 2x per year
 - if accepted, easier resource access (low number of competitive users)
 - purpose:
 - large (proven) computations using homogeneous infrastructure







Research Infrastructures in CR II.

- National Grid Infrastructure (NGI) MetaCentrum
 - ca 17860 computing cores (including CERIT-SC resources)
 - attributes:
 - computing time available free of charge, without formal applications
 - heterogeneous resources available (including "exotic" ones)
 - resources shared with competitive users (sometimes hard to access)
 - purpose:
 - common smaller to middle -sized computations (larger computations after agreement)
 - preparation of computations/projects for computations at IT4innovations
 (~ technical readiness)
- · CERIT-SC @ ICS MU
 - HW and SW provider (resources available through NGI)
 - main emphasis on services supporting user research



CERIT-SC & NGI



Centre CERIT-SC

A computing and research centre operating at Masaryk University

- long-term history (→ long-term experience in ICT science)
 - CERIT-SC evolved from Supercomputing Center Brno (established in 1994), and
 - participates on the operation of National Grid Infrastructure

Our mission:

http://www.cerit-sc.cz

- production services for computational science
 - high-performance computing clusters
 - large data storage, back-ups and data archives
 - web portals & projects' back-office
- an application of top-level ICT in the science
 - own research in e-infrastructures (know-how)
 - novel forms of infrastructure utilization (experimental usage support)
 - research collaborations with other science areas



Centre CERIT-SC & NGI

CERIT-SC is an important NGI partner

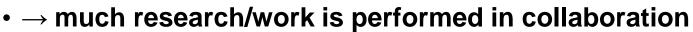
HW & SW resources provider

SMP nodes (2592 cores)
HD nodes (2624 cores)
SGI UV node (288 cores, 6 TB RAM)
SGI UV node (384 cores, 6 TB RAM)

storage capacity (~ 3,5 PB)



CERIT-SC (SCB) established MetaCentrum NGI



http://www.cerit-sc.cz





Fact I. Common HW centers provide just a "dumb" power without any support how to effectively use it

Fact II. Common HW centers do not participate on the users' research <u>aiming to help them</u> with ICT problems

CERIT-SC collaborates with its users:

- to help them effectively use the provided resources
- to help them to cope with their ICT research problems focusing on an application of top-level ICT in the science smaller as well as bigger (= funded) projects



What's the idea?

We focus on <u>intelligent & novel usage forms</u> of the provided infrastructure

- the provided HW/SW resources serve just as a tool for research and development
 - → highly-flexible infrastructure (convenient to experiments) in comparison with NGI resources, the production computations are at the second-level of interest
- the centre aims to be equipped with cutting-edge technologies in order to allow top-level research (both internal & collaborative)
- real research collaboration with our partners
 the collaborations generate new questions/problems for IT
 the collaborations generate novel opportunities for the science
 (we DON'T want to be a common service organization)

How do we fulfill the idea?



How are the research collaborations performed?

- the work is carried via a doctoral/diploma thesis of a FI MU student
- the CERIT-SC staff supervises/consults the student and regularly meets with the research partners
 - the partners provide the expert knowledge from the particular area

Collaborations through (international) projects

CERIT-SC participates on several projects, usually developing IT infrastructure supporting the particular research area

ELIXIR-CZ, BBMRI, Thalamoss, SDI4Apps, Onco-Steer, CzeCOS/ICOS, ... KYPO, 3M SmartMeters in cloud, MeteoPredictions, ...

Strong ICT expert knowledge available:

- long-term collaboration with Faculty of Informatics MU
- long-term collaboration with CESNET
 - → consultations with experts in particular areas



Selected research collaborations

Selected (ongoing) collaborations is

3D tree reconstructions from terrestrial LiDAR scans

- partner: Global Change Research Centre Academy of Sciences of the Czech Republic (CzechGlobe)
- the goal: to propose an algorithm able to perform fully-automated reconstruction of tree skeletons (main focus on Norway spruce trees)
 - from a 3D point cloud
 - scanned by a LiDAR scanner
 - the points provide information about XYZ coordinates
 + reflection intensity
 - the expected output: 3D tree skeleton
- the main issue: overlaps (→ gaps in the input data)

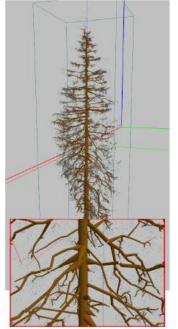
Selected (ongoing) collaborations

3D tree reconstructions from terrestrial LiDAR scans - cont'd

the diploma thesis proposed a novel innovative approach

to the reconstructions of 3D tree models

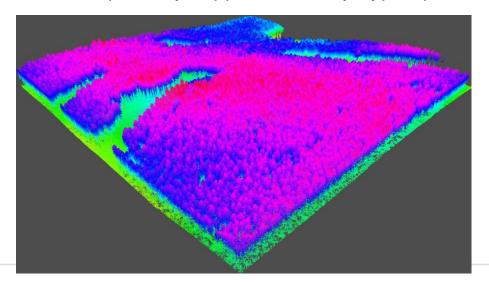
- the reconstructed models used in subsequent research
 - determining a statistical information about the amount of wood biomass and about basic tree structure
 - parametric supplementation of green biomass
 (young branches+ needles) a part of the PhD work
 - importing the 3D models into tools performing various analysis (e.g., DART radiative transfer model)

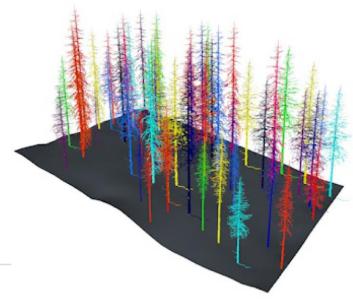


Selected (ongoing) collaborations

3D reconstruction of tree forests from full-wave LiDAR scans

- subsequent work
- the goal: an accurate 3D reconstruction of tree forests scanned by aerial full-waveform LiDAR scans
 - possibly supplemented by hyperspectral or thermal scans, in-situ measurements,...







An algorithm for determination of problematic closures in a road network

- partner: Transport Research Centre, Olomouc
- the goal: to find a robust algorithm able to identify all the road network break-ups and evaluate their impacts
- main issue: computation demands
 - the brute-force algorithms fail because of large state space
 - 2 algorithms proposed able to cope with multiple road closures



- An application of neural networks for filling in the gaps in eddy-covariance measurements
 - partner: CzechGlobe
- Biobanking research infrastructure (BBMRI_CZ)
 - partner: Masaryk Memorial Cancer Institute, Recamo
- Propagation models of epilepsy and other processes in the brain
 - partner: MED MU, ÚPT AV, CEITEC
- Photometric archive of astronomical images
- Extraction of photometric data on the objects of astronomical images
 - 2x partner: partner: Institute of theoretical physics and astrophysics SCI MU
- Bioinformatic analysis of data from the mass spectrometer
 - partner: Institute of experimental biology SCI MU
- Synchronizing timestamps in aerial landscape scans
 - partner: CzechGlobe
- Optimization of Ansys computation for flow determination around a large two-shaft gas turbine
 - partner: SVS FEM
- 3.5 Million smartmeters in the cloud
 - partner: CEZ group, MycroftMind







Additional services available to academic research community







Data services

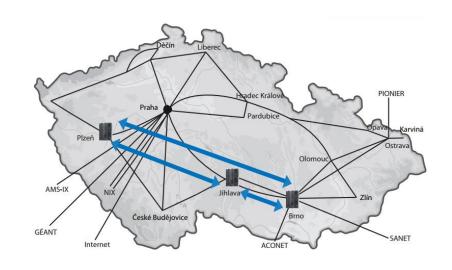
Hierarchical data storages

- 22+ PB of physical capacity
- useful for data archivals, backups, etc.
- various access protocols available

Further end-user services

- FileSender
- OwnCloud

http://du.cesnet.cz









Data Services for end-users

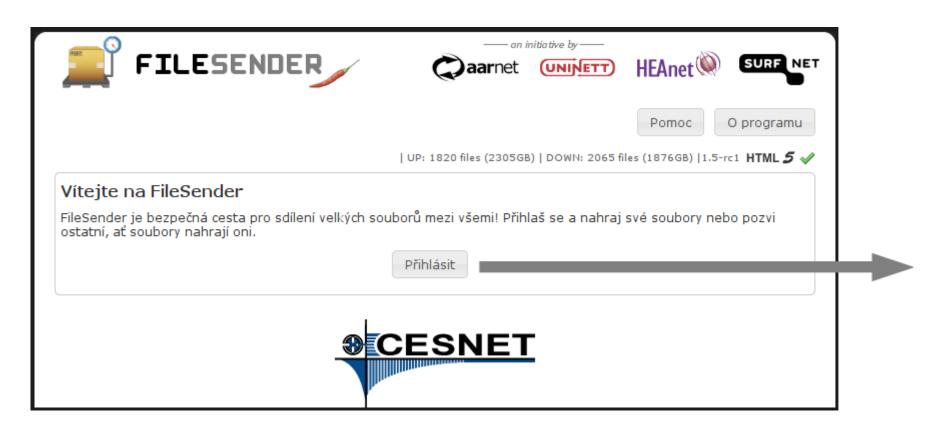
- FileSender file sharing/transfering service
 - web service intended for sending big data files
 - big = current limit is 500 GB
 - http://filesender.cesnet.cz
 - at least one user has to be an authorized infrastructure user
 - federated authentication through edulD.cz
 - authorized user is allowed to upload a file (and send a notification to the receiver)
 - if an authorized user needs to receive data from a non-authorized user, she sends him an invitation link (so he is allowed to use it for uploading the file)







FileSender – example I.









FileSender – example II.











FileSender – example III.

FILE	ESENDER Caarnet	UNINETT	HEAnet SURF NET					
Nahrát nový soubor Vítejte Tomáš Košňar	Pozvánky Mé soubory	Pomoc	O programu Odhlásit					
Nahrát soubor	10.11020 11.07 (200000)							
Příjemce: Odesílatel: Předmět: (volitelné) Zpráva: (volitelné)	tomas.kosnar@cesnet.cz	234	Vložte emailové adresy příjemců Nastavte datum expirace Vyberte soubor Klikněte na Odeslat					
Datum expirace: Vyberte soubor:	18. 04. 2013 Vybrat soubor Soubor nevybrán Souhlasím s podmínkami užití této služby. [Zobrazit/Skrýt]							
**CESNET								







OwnCloud

- cloud storage "like Dropbox"
 - quota: 100 GB / user
 - available through web interface
 - https://owncloud.cesnet.cz/
 - clients for Windows, Linux, OS X
 - clients for smartphones and tablets
 - allows sharing among a group of users
 - data backups every day
 - document versioning
 - calendars and contacts sharing
 - etc.









OwnCloud – example I.









OwnCloud – example II.

-	Přihlásit účtem		
	České vysoké učení technické v Praze, Fakulta elektrotechnická	F	
	CESNET	* CESNET	
	Masarykova univerzita	AND SOLVE STATE OF THE SOLVE STA	
	Univerzita Hradec Králové	Univerzita Hradec Králové	
	Univerzita Pardubice	4	
	Západočeská univerzita v Plzni	D •	
	Jiný účet		
	፟	CESNET 1	







OwnCloud – example III.

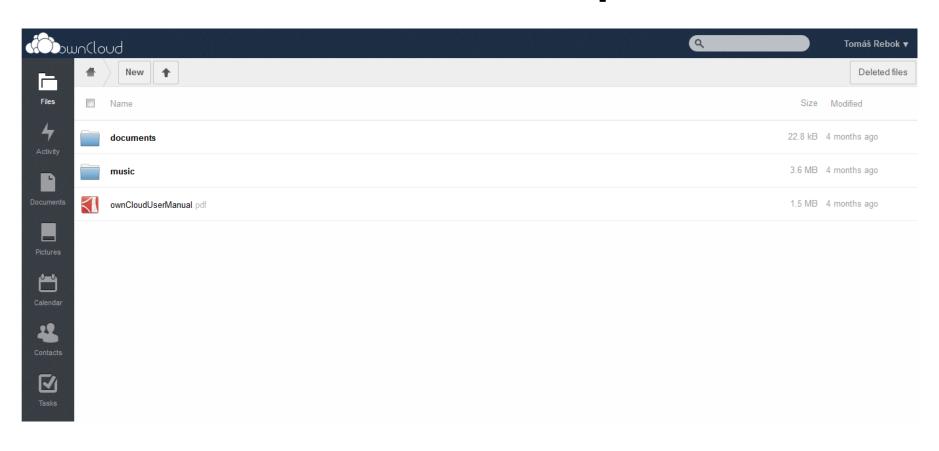
WERSITAS . SISNEY	MASARYKOVA UNIVERZITA Česká republika
	Poskytovatel identit MU
	UČO: 39685 Heslo: •••••••
	Přihlásit Pokusili jste se přistoupit na stránky, které vyžadují autentizaci. Pro přihlášení použijte UČO a sekundární heslo.
	Hosté s guest účtem použijí místo UČO své GuestID.
	<u>Nápověda</u>
	Službu zajišťuje <u>Ústav výpočetní techniky MU</u> .
	<u>English</u>







OwnCloud – example IV.









Remote collaboration support

Support for interactive collaborative work in real-time

- videoconferences
 HD videoconferencing support via H.323 HW/SW equippment
- webconferences
 SD videoconferencing support via Adobe Connect (Adobe Flash)
 http://meetings.cesnet.cz
- special transmissions
 HD, UHD, 2K, 4K, 8K with compressed/uncompressed video transmission (UltraGrid tool)
- IP telephony

Support for offline content access

- streaming
- video archive







http://vidcon.cesnet.cz







Security services

Security incidents handling

- detailed monitoring of possible security incidents
- the users/administrators are informed about security incidents, and
- helped to resolve the incident
- additional services:
 seminars, workshops, etc.

Security teams CSIRT-MU and CESNET-CERTS

– several successes:

e.g., Chuck Norris botnet discovery











Federated identity management

Czech academic identity federation edulD.cz

- provides means for inter-organizational identity management and access control to network services, while respecting the privacy of the users
- users may access multiple applications using just a single password
- service provider administrators do not have to preserve user's credentials and implement authentication
- user authentication is always performed at the home organization, user credenitals are not revealed to the service providers



http://www.eduid.cz







PKI – users and servers certificates

CESNET CA certification authority

- provides the users with TERENA (Trans-European Research and Education Networking Association) certificates
 - usable for electronic signatures as well as for encryption

CESNET CA services:

- issues personal certificates
- issues certificates for servers and services
- certificates registration offices
- certificates certification offices



http://pki.cesnet.cz







Mobility and roaming support

Eduroam.cz

 idea to enable transparent usage of (especially wireless) networks of partner (Czech as well as abroad) institutions



http://www.eduroam.cz



30.1.2019





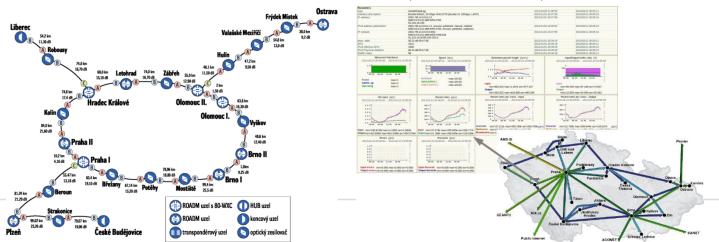
Communication infrastructure and its monitoring

The basis of all the services: high-speed computer network

- 100 Gbps, called CESNET2
- interconnected with pan-european network GÉANT

and its monitoring

- detailed network monitoring (quality issues as well as individual nodes behaviour) available
- automatic detection of various events, anomalies, etc.









Conclusions







Conclusions

CESNET infrastructure:

- computing services (MetaCentrum NGI & MetaVO)
- data services (archivals, backups, data sharing and transfers, ...)
- remote collaborations support servicese (videoconferences, webconferences, streaming, ...)
- further supporting services (...)

Centrum CERIT-SC:

- computing services (flexible infrastructure for production and research)
- services supporting collaborative research
- user identities/accounts shared with the CESNET infrastructure
- The message: "If you cannot find a solution to your specific needs in the provided services, let us know - we will try to find the solution together with you…"











The CERIT Scientific Cloud project (reg. no. CZ.1.05/3.2.00/08.0144) is supported by the *Operational Program Research and Development for Innovations*, priority axis 3, subarea 2.3 *Information Infrastructure for Research and Development*.

http://metavo.metacentrum.cz

http://www.cerit-sc.cz

EUROPEAN REGIONAL DEVELOPMENT FUND

INVESTING IN YOUR FUTURE







Hands-on training for MetaCentrum/CERIT-SC users

Tomáš Rebok

MetaCentrum, CESNET CERIT-SC, Masaryk University rebok@ics.muni.cz







Overview

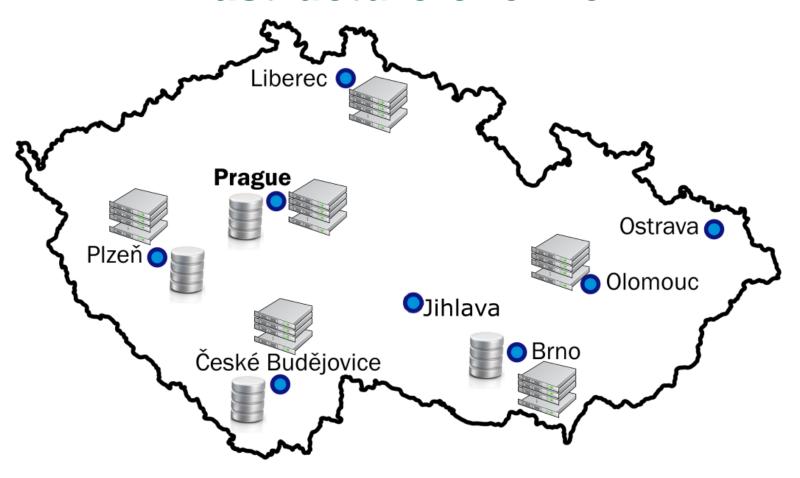
- Introduction
- MetaCentrum / CERIT-SC infrastructure overview
- How to ... specify requested resources
- How to ... run an interactive job
- How to ... use application modules
- How to ... run a batch job
- How to ... determine a job state
- Another mini-HowTos ...
- What to do if something goes wrong?
- Real-world examples
- Appendices







Infrastructure overview

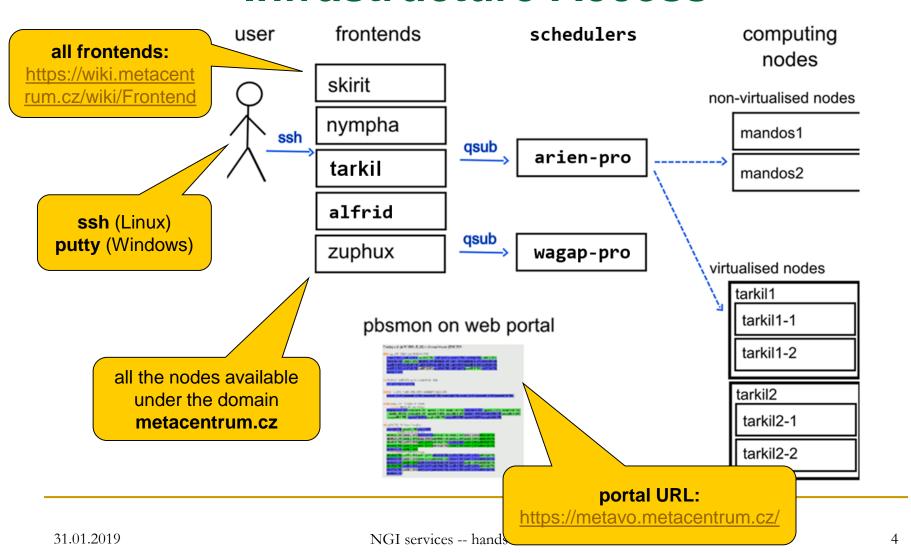








Infrastructure Access

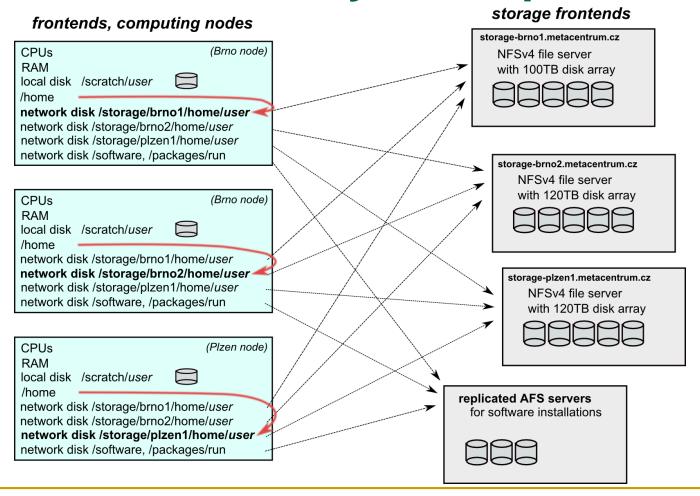








Infrastructure System Specifics









Overview

- Introduction
- MetaCentrum / CERIT-SC infrastructure overview
- How to ... specify requested resources
- How to ... run an interactive job
- How to ... use application modules
- How to ... run a batch job
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- Another mini-HowTos ...
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How to ... specify requested resources I.

- before running a job, one needs to know what resources the job requires
 - and how much/many of them
- for example:
 - number of nodes
 - number of CPUs/cores per node
 - an upper estimation of job's runtime
 - amount of free memory
 - amount of scratch space for temporal data
 - number of requested software licenses
 - etc.
- the resource requirements are then provided to the qsub utility (when submitting a job)
 - the requested resources are reserved for the job by the infrastructure scheduler
 - the computation is allowed to use them
- details about resources' specification: https://wiki.metacentrum.cz/wiki/About_scheduling_system







How to ... specify requested resources II.

Graphical way:

- qsub assembler: https://metavo.metacentrum.cz/pbsmon2/qsub_pbspro
- allows to:
 - graphically specify the requested resources
 - check, whether such resources are available
 - generate command line options for qsub
 - check the usage of MetaVO resources

Textual way:

- more powerful and (once being experienced user) more convenient
- see the following slides/examples →







PBS Professional – the infrastructure scheduler

- a novel scheduling system used in MetaCentrum NGI
 - see advanced information at
 https://wiki.metacentrum.cz/wiki/Prostředí PBS Professional

New term - CHUNK:

- chunk = further indivisible set of resources allocated to a job on a physical node
 - ≈ virtual node
- contains resources, which could be asked from the infrastructure nodes
- for simplicity reasons: chunk = node
 - later, we will generalize...







How to ... specify requested resources III.

Chunk(s) specification:

general format: -1 select=...

Examples:

- 2 chunks/nodes:
 - □ -l select=2
- 5 chunks/nodes:
 - □ -l select=5
- by default, allocates just a single core in each chunk
 - ⇒ should be used together with number of CPUs (NCPUs) specification
- if "-l select=..." is not provided, just a single chunk with a single CPU/core is allocated







How to ... specify requested resources IV.

Number of CPUs (NCPUs) specification (in each chunk):

- general format: -1 select=...:ncpus=...
- 1 chunk with 4 cores:
 - □ -l select=1:ncpus=4
- 5 chunks, each of them with 2 cores:
 - □ -l select=5:ncpus=2

(Advanced chunks specification:)

- general format: -1 select=[chunk_1][+chunk_2]...[+chunk_n]
- 1 chunk with 4 cores and 2 chunks with 3 cores and 10 chunks with 1 core:
 - □ -l select=1:ncpus=4+2:ncpus=3+10:ncpus=1







How to ... specify requested resources V.

Other useful features:

- chunks from just a single (specified) cluster (suitable e.g. for MPI jobs):
 - general format: -1 select=...:cl <cluster name>=true
 - □ e.g., -1 select=3:ncpus=1:cl doom=true
- chunks located in a specific location (suitable when accessing storage in the location)
 - general format: -l select=...:
brno|plzen|praha|...>=true
 - □ e.g., -l select=1:ncpus=4:brno=true
- exclusive node(s) assignment (useful for testing purposes, all resources available):
 - □ *general format:* -1 select=... -1 place=exclhost
 - e.g., -l select=1 -l place=exclhost
- negative specification:
 - general format: -l select=...:<feature>=false
 - □ e.g., -l select=1:ncpus=4:hyperthreading=false
- · ...

A list of nodes' features can be found here: http://metavo.metacentrum.cz/pbsmon2/props







How to ... specify requested resources VI.

Specifying memory resources (default = 400mb):

```
general format: -l select=...:mem=...<suffix>
    e.g., -l select=...:mem=100mb
    e.g., -l select=...:mem=2gb
```

Specifying job's maximum runtime (default = 24 hours):

- it is necessary to specify an upper limit on job's runtime:
- general format: -l walltime=[[hh:]mm:]ss
 - □ **e.g.**, -l walltime=13:00
 - □ **e.g.**, -1 walltime=2:14:30







How to ... specify requested resources VII.

Specifying requested scratch space:

- useful, when the application performs I/O intensive operations OR for long-term computations (reduces the impact of network failures)
- requesting scratch is <u>mandatory</u> (no defaults)
- scratch space specification: -l select=...:scratch_type=...<suffix>
 e.g., -l select=...:scratch local=500mb

Types of scratches:

- scratch_local
- scratch_ssd
- scratch_shared

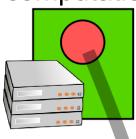






Why to use scratches?

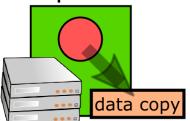
computation



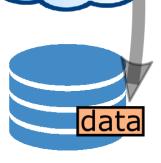
Data processing using central storage

- low computing performance (I/O operations)
- dependency on (functional) network connection
- high load on the central storage

computation







Data processing using scratches

- + highest computing performance
- + resilience to network connection failures
- + minimal load on the central storage











How to use scratches?

- there is a private scratch directory for particular job
 - □ /scratch/\$USER/job \$PBS JOBID directory for (local) job's scratch
 - /scratch.ssd/\$USER/job_\$PBS_JOBID for job's scratch on SSD
 - /scratch.shared/\$USER/job_\$PBS_JOBID for shared job's scratch
 - the master directory /scratch*/\$USER is not available for writing
- to make things easier, there is a SCRATCHDIR environment variable available in the system
 - (within a job) points to the assigned scratch space/location

Please, clean scratches after your jobs

- there is a "clean_scratch" utility to perform safe scratch cleanup
 - also reports scratch garbage from your previous jobs
 - usage example will be provided later







How to ... specify requested resources VIII.

Specifying requested software licenses:

- necessary when an application requires a SW licence
 - the job becomes started once the requested licences are available
 - the information about a licence necessity is provided within the application description (see later)
- general format: -1 <lic name>=<amount>
 - □ **e.g.**, -l matlab=1 -l matlab Optimization Toolbox=4
 - \bullet e.g., -1 gridmath8=20

(advanced) Dependencies among jobs

- allows to create a workflow
 - e.g., to start a job once another one successfully finishes, breaks, etc.
- see qsub's "-w" option (man qsub)
 - □ **e.g.**, \$ qsub ... -W depend=afterok:12345.arien-pro.ics.muni.cz







Resource chunks vs. nodes

select=4:ncpus=2:mem=5g:scratch_local=10g

2x CPU	2x CPU	2x CPU	2x CPU	
5GB RAM	5GB RAM	5GB RAM	5GB RAM	?
10GB HDD	10GB HDD	10GB HDD	10GB HDD	

How do chunks correspond to nodes?

- chunks arrangement option "-1 place=..."
 - -1 place=free: chunks are **free to spread** over available nodes
 - default behaviour
 - -1 place=pack: all chunks will be allocated **on the same node**
 - the node has to have enough resources available
 - -1 place=scatter: each chunk will be allocated on a different node

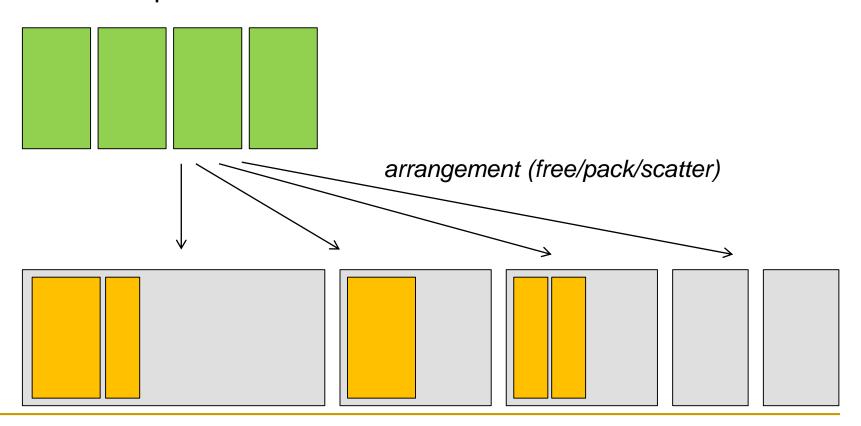






Chunks arrangement

free vs. pack vs. scatter



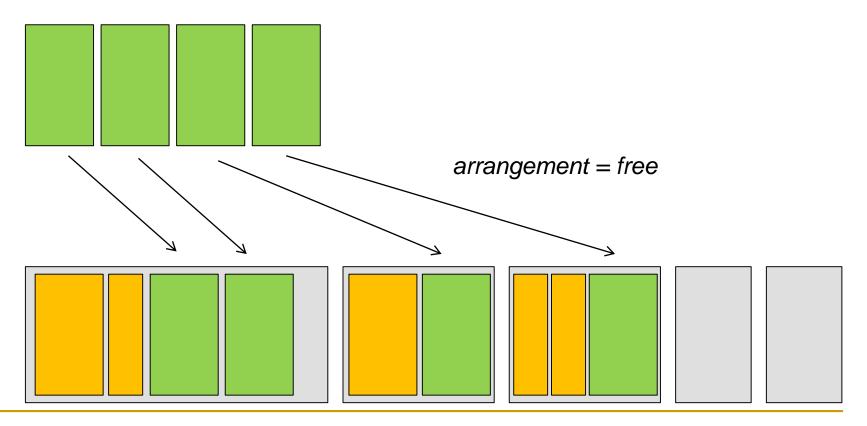






Chunks arrangement

free (default) vs. pack vs. scatter



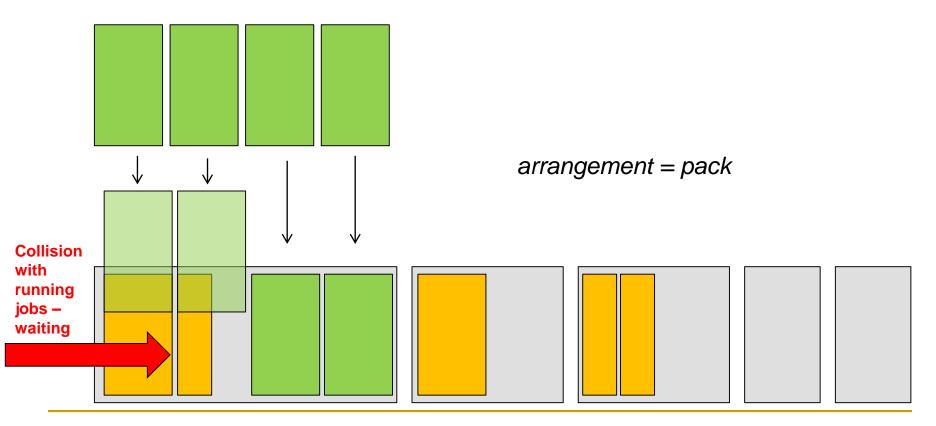






Chunks arrangement

free (default) vs. <u>pack</u> vs. scatter



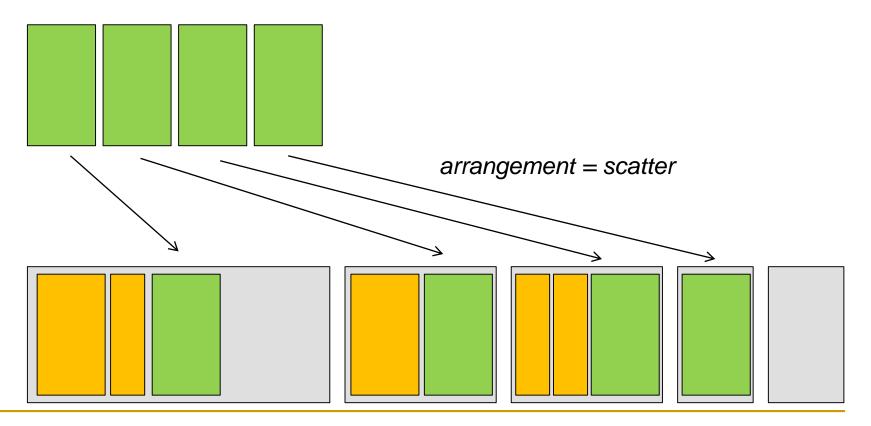






Chunks arrangement

free (default) vs. pack vs. <u>scatter</u>









How to ... specify requested resources IX.

Questions and Answers:

- Why is it necessary to specify the resources in a proper number/amount?
 - because when a job consumes more resources than announced, it will be killed by us (you'll be informed)
 - otherwise it may influence other processes running on the node
- Why is it necessary not to ask for excessive number/amount of resources?
 - the jobs having smaller resource requirements are started (i.e., get the time slot) faster
- Any other questions?









How to ... specify requested resources IX.

Questions and Answers:

- Why is it necessary to specify the resources in a proper number/amount?
 - because when a job consumes more resources than announced, it will be killed by us (you'll be informed)
 - otherwise it may influence other processes running on the node
- Why is it necessary not to ask for excessive number/amount of resources?
 - the jobs having smaller resource requirements are started

See more details about PBSpro scheduler:

https://metavo.metacentrum.cz/cs/seminars/seminar2017/presentation-Klusacek.pptx

SHORT guide:

https://metavo.metacentrum.cz/export/sites/meta/cs/seminars/seminar2 017/tahak-pbs-pro-small.pdf







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How to ... run an interactive job I.

Interactive jobs:

- result in getting a prompt on a single (master) node
 - one may perform interactive computations
 - the other nodes, if requested, remain allocated and accessible (see later)
- How to ask for an interactive job?
 - □ add the option "-I" to the qsub command
 - e.g., qsub -I -l select=1:ncpus=4
- Example (valid just for this demo session):
 - □ qsub -I -q MetaSeminar # (-l select=1:ncpus=1)

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How to ... run an interactive job II.

Textual mode: simple

Graphical mode:

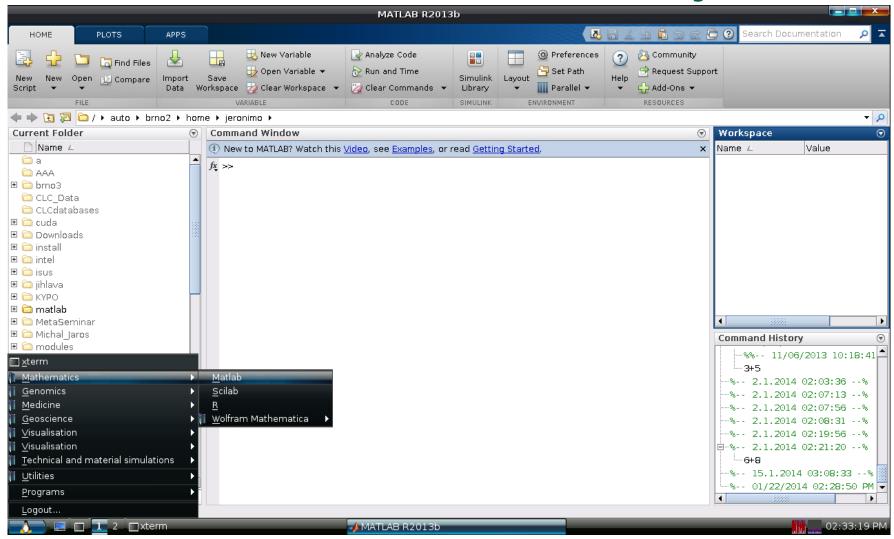
- (preffered) remote desktops based on VNC servers (pilot run):
- available from frontends as well as computing nodes (interactive jobs)
 - module add gui
 - □ gui start [-s] [-g GEOMETRY] [-c COLORS]
 - uses one-time passwords
 - allows to access the VNC via a supported TigerVNC client
 - allows SSH tunnels to be able to connect with a wide-range of clients
 - allows to specify several parameters (e.g., desktop resolution, color depth)
 - gui info [-p] ... displays active sessions (optionally with login password)
 - □ gui traverse [-p] ... display all the sessions throughout the infrastructure
 - gui stop [sessionID] ... allows to stop/kill an active session
- see more info at
 https://wiki.metacentrum.cz/wiki/Remote desktop







How to ... run an interactive job II.









How to ... run an interactive job II.

Graphical mode (further options):

- (fallback) tunnelling a display through ssh (Windows/Linux):
 - connect to the frontend node having SSH forwarding/tunneling enabled:
 - Linux: ssh -X skirit.metacentrum.cz
 - Windows:
 - □ install an XServer (e.g., Xming)
 - set Putty appropriately to enable X11 forwarding when connecting to the frontend node
 - Connection \rightarrow SSH \rightarrow X11 \rightarrow Enable X11 forwarding
 - ask for an interactive job, adding "-x" option to the qsub command
 - \bullet e.g., qsub -I \bullet X -l select=... ...
- (tech. gurus) exporting a display from the master node to a Linux box:
 - export DISPLAY=mycomputer.mydomain.cz:0.0
 - on a Linux box, run "xhost +" to allow all the remote clients to connect
 - be sure that your display manager allows remote connections







How to ... run an interactive job III.

Questions and Answers:

- How to get an information about the other nodes/chunks allocated (if requested)?
 - master node\$ cat \$PBS NODEFILE
 - works for batch jobs as well
- How to use the other nodes/chunks? (holds for batch jobs as well)
 - MPI jobs use them automatically
 - otherwise, use the pbsdsh utility (see "man pbsdsh" for details) to run a remote command
 - if the pbsdsh does not work for you, use the ssh to run the remote command
- Any other questions?







How to ... run an interactive job III.

Questions and Answers:

How to get an information about the other nodes/chunks allocated

Hint:

- there are several useful environment variables one may use
 - \$ set | grep PBS
- e.g.:
 - PBS_JOBID ... job's identificator
 - PBS_NUM_NODES, PBS_NUM_PPN ... allocated number of nodes/processors
 - PBS_O_WORKDIR ... submit directory
 - ...











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How to ... use application modules I.

Application modules:

- the modullar subsystem provides a user interface to modifications of user environment, which are necessary for running the requested applications
- allows to "add" an application to a user environment
- getting a list of available application modules:
 - \$ module avail
 - \$ module avail matl
 - https://wiki.metacentrum.cz/wiki/Kategorie:Applications
 - provides the documentation about modules' usage
 - besides others, includes:
 - information whether it is necessary to ask the scheduler for an available licence
 - information whether it is necessary to express consent with their licence agreement







How to ... use application modules II.

Application modules:

- loading an application into the environment:
 - □ \$ module add <modulename>
 - □ e.g., module add maple
- listing the already loaded modules:
 - □ \$ module list.
- unloading an application from the environment:
 - \$ module del <modulename>
 - e.g., module del openmpi
- Note: An application may require to express consent with its licence agreement before it may be used (see the application's description). To provide the aggreement, visit the following webpage: https://metavo.metacentrum.cz/cs/myaccount/licence.html
- for more information about application modules, see
 https://wiki.metacentrum.cz/wiki/Application_modules







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Preparation before batch demos

Copy-out the pre-prepared demos:

\$ cp -rH /storage/brno2/home/jeronimo/MetaSeminar/latest \$HOME

Text editors in Linux:

- experienced users: vim <filename>
 - very flexible, feature-rich, great editor...
- common users: mcedit <filename>









Preparation before batch demos

Copy-out the pre-prepared demos:

\$ cp -rH /storage/brno2/home/jeronimo/MetaSeminar/latest \$HOME

Text editors in Linux:

- experienced users: vim <filename>
 - very flexible, feature-rich, great editor...
- common users: mcedit <filename>
 - □ easy to remember alternative: pico <filename> ◎









How to ... run a batch job I.

Batch jobs:

- perform the computation as described in their startup script
 - the submission results in getting a job identifier, which further serves for getting more information about the job (see later)
- How to submit a batch job?
 - add the reference to the startup script to the qsub command
 - □ e.g., qsub -1 select=3:ncpus=4 <myscript.sh>
- Example (valid for this demo session):
 - qsub -q MetaSeminar -l select=1:ncpus=1 myscript.sh
 - results in getting something like "12345.arien-pro.ics.muni.cz"







for

How to ... run a batch job I.

B Hint:

- create the file myscript.sh with the following content:
 - \$ vim myscript.sh

```
#!/bin/bash
```

my first batch job
uname -a

- see the standard output file (myscript.sh.o<JOBID>)
 - \$ cat myscript.sh.o<JOBID>
 - qsub -q MetaSeminar -l select=1:ncpus=1 myscript.sh
 - results in getting something like "12345.arien-pro.ics.muni.cz"







How to ... run a batch job II.

Startup script skelet: (non IO-intensive computations)

use just when you know, what you are doing...

```
#!/bin/bash

DATADIR="/storage/brno2/home/$USER/" # shared via NFSv4
cd $DATADIR

# ... load modules & perform the computation ...
```

further details – see

https://wiki.metacentrum.cz/wiki/How_to_compute/Requesting_resources







How to ... run a batch job III.

Recommended startup script skelet: (IO-intensive computations or long-term jobs)

#!/bin/bash # set a handler to clean the SCRATCHDIR once finished trap 'clean scratch' EXIT TERM # if temporal results are important/useful # trap 'cp -r \$SCRATCHDIR/neuplna.data \$DATADIR && clean scratch' TERM # set the location of input/output data # DATADIR="/storage/brno2/home/\$USER/" DATADIR="\$PBS O WORKDIR" # prepare the input data cp \$DATADIR/input.txt \$SCRATCHDIR # go to the working directory and perform the computation cd \$SCRATCHDIR # ... load modules & perform the computation ... # copy out the output data # if the copying fails, let the data in SCRATCHDIR and inform the user cp \$SCRATCHDIR/output.txt \$DATADIR || export CLEAN SCRATCH=false







How to ... run a batch job IV.

Using the application modules within the batch script:

- module add SW
 - □ e.g., "module add maple"
- include the initialization line ("source ...") if necessary:
 - i.e., if you experience problems like "module: command not found", then add source /software/modules/init before "module add" sections

Getting the job's standard output and standard error output:

- once finished, there appear two files in the directory, which the job has been started from:
 - < < job_name > . o < jobID > ... standard output
 - <job name>.e<jobID>... standard error output
 - □ the <job_name> can be modified via the "-N" qsub option







How to ... run a batch job V.

Job attributes specification:

in the case of batch jobs, the requested resources and further job information (job attributes in short) may be specified either on the command line (see "man qsub") or directly within the script:

by adding the "#PBS" directives (see "man qsub"):

```
#PBS -N Job_name
#PBS -l select=2:ncpus=1:mem=320kb:scratch_local=100m
#PBS -m abe
#
< ... commands ... >
```

- the submission may be then simply performed by:
 - □ \$ qsub myscript.sh
 - if options are provided both in the script and on the command-line, the command-line arguments override the script ones







How to ... run a batch job VI. (complex example)

```
#!/bin/bash
#PBS -1 select=1:ncpus=2:mem=500mb:scratch local=100m
#PBS -m abe
# set a handler to clean the SCRATCHDIR once finished
trap "clean scratch" EXIT TERM
# set the location of input/output data
DATADIR="$PBS O WORKDIR"
# prepare the input data
cp $DATADIR/input.mpl $SCRATCHDIR
# go to the working directory and perform the computation
cd $SCRATCHDIR
# load the appropriate module
module add maple
# run the computation
maple input.mpl
# copy out the output data (if it fails, let the data in SCRATCHDIR and inform the user)
cp $SCRATCHDIR/output.gif $DATADIR || export CLEAN SCRATCH=false
```







How to ... run a batch job VII.

Questions and Answers:

- Should you prefer batch or interactive jobs?
 - definitely the batch ones they use the computing resources more effectively
 - use the interactive ones just for testing your startup script, GUI apps, or data preparation

Any other questions?









How to ... run a batch job VIII.

Example:

Create and submit a batch script, which performs a simple Maple computation, described in a file:

```
plotsetup(gif, plotoutput=`myplot.gif`,
    plotoptions=`height=1024,width=768`);
plot3d( x*y, x=-1..1, y=-1..1, axes = BOXED, style =
    PATCH);
```

- process the file using Maple (from a batch script):
 - hint: \$ maple <filename>







How to ... run a batch job VIII.

Example:

Create and submit a batch script, which performs a simple Maple computation, described in a file:

```
plotsetup(gif, plotoutput=`myplot.gif`,
    plotoptions=`height=1024,width=768`);
plot3d( x*y, x=-1..1, y=-1..1, axes = BOXED, style =
    PATCH);
```

- process the file using Maple (from a batch script):
 - hint: \$ maple <filename>

Hint:

• see the solution at /storage/brno2/home/jeronimo/MetaSeminar/latest/Maple







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How to ... determine a job state I.

Job identifiers

- every job (no matter whether interactive or batch) is uniquely identified by its identifier (JOBID)
 - e.g., 12345.arien-pro.ics.muni.cz
- to obtain any information about a job, the knowledge of its identifier is necessary
 - how to list all the recent jobs?
 - graphical way PBSMON: http://metavo.metacentrum.cz/pbsmon2/jobs/allJobs
 - frontend\$ qstat (run on any frontend)
 - □ to include finished ones, run \$ qstat -x
 - how to list all the recent jobs of a specific user?
 - graphical way PBSMON: https://metavo.metacentrum.cz/pbsmon2/jobs/my
 - frontend\$ qstat -u <username> (again, any frontend)
 - □ to include finished ones, run \$ qstat -x -u <username>







How to ... determine a job state II.

How to determine a job state?

- graphical way see PBSMON
 - list all your jobs and click on the particular job's identifier
 - http://metavo.metacentrum.cz/pbsmon2/jobs/my
- textual way qstat command (see man qstat)
 - □ brief information about a job: \$ qstat JOBID
 - informs about: job's state (Q=queued, R=running, E=exiting, F=finished, ...), job's runtime, ...
 - complex information about a job: \$ qstat -f JOBID
 - shows all the available information about a job
 - useful properties:
 - exec_host -- the nodes, where the job did really run
 - resources_used, start/completion time, exit status, ...
 - necessary to add "-x" option when examining already finished job(s)







How to ... determine a job state III.

Hell, when my jobs will really start?

- nobody can tell you ☺
 - the God/scheduler decides (based on the other job's finish)
 - we're working on an estimation method to inform you about its probable startup
- check the queues' fulfilment:
 http://metave.metacentrum.or/en/etate/ich
 - http://metavo.metacentrum.cz/cs/state/jobsQueued
 - the higher fairshare (queue's AND job's) is, the earlier the job will be started
- stay informed about job's startup / finish / abort (via email)
 - by default, just an information about job's abortation is sent
 - o when submitting a job, add "-m abe" option to the qsub command to be informed about all the job's states
 - or "#PBS -m abe" directive to the startup script







How to ... determine a job state IV.

Monitoring running job's stdout, stderr, working/temporal files

- 1. via ssh, log in directly to the execution node(s)
 - how to get the job's execution node(s)?
- to examine the working/temporal files, navigate directly to them
 - logging to the execution node(s) is necessary -- even though the files are on a shared storage, their content propagation takes some time
- to examine the stdout/stderr of a running job:
 - navigate to the /var/spool/pbs/spool/ directory and examine the files:
 - \$PBS_JOBID.OU for standard output (stdout e.g., "1234.arien-pro.ics.muni.cz.OU")
 - \$PBS_JOBID.ER for standard error output (stderr e.g., "1234.arien-pro.ics.muni.cz.ER")

Job's forcible termination

- \$ qdel JOBID (the job may be terminated in any previous state)
- during termination, the job turns to E (exiting) and finally to F (finished) state







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Another mini-HowTos ...

- how to use privileged resources?
 - if your institution/project integrates HW resources, a defined group of users may have priority access to them
 - technically accomplished using scheduler queues
 - a job has to be submitted to the particular queue
 - □ qsub -l select=... -l walltime=... -q PRIORITY QUEUE script.sh
 - e.g., ELIXIR CZ project integrates a set of resources
 - priority queue "elixir_2w" available for ELIXIR CZ users
 - moving jobs between scheduler queues
 - from priority queue to default queue
 - □ qmove default JOBID
 - from default queue(s) to a priority queue
 - □ qmove elixir 2w JOBID







Another mini-HowTos ...

- how to make your SW tool available within MetaVO?
 - commercial apps:
 - assumption: you own a license, and the license allows the application to be run on our infrastructure (nodes not owned by you, located elsewhere, etc.)
 - once installed, we can restrict its usage just for you (or for your group)
 - open-source/freeware apps:
 - you can compile/install the app in your HOME directory
 - OR you can install/compile the app on your own and ask us to make it available in the software repository
 - compile the application in your HOME directory
 - prepare a modulefile setting the application environment
 - inspire yourself by modules located at /packages/run/modules-2.0/modulefiles
 - □ test the app/modulefile
 - \$ export MODULEPATH=\$MODULEPATH:\$HOME/myapps
 - □ see https://wiki.metacentrum.cz/wiki/How to install an application
 - OR you can ask us for preparing the application for you







Another mini-HowTos ...

- how to ask for nodes equipped by GPU cards?
 - determine, how many GPUs your application will need (-1 ngpus=X)
 - consult the HW information page: http://metavo.metacentrum.cz/cs/state/hardware.html
 - determine, how long the application will run (if you need more, let us know)
 - gpu queue ... maximum runtime 1 day
 - qpu long queue ... maximum runtime 1 week
 - Note: GPU Titan V available through gpu titan queue (zuphux.cerit-sc.cz)
 - make the submission:
 - \$ qsub -l select=1:ncpus=4:mem=10g:ngpus=1 -q gpu_long -l walltime=4d ...
 - specific GPU cards by restricting the cluster: qsub -l select=...:cl doom=true ...
 - do not change the CUDA_VISIBLE_DEVICES environment variable
 - it's automatically set in order to determine the GPU card(s) that has/have been reserved for your application
 - general information: https://wiki.metacentrum.cz/wiki/GPU_clusters







How to ask for nodes equipped with Xeon Phi?

phi[1-6].cerit-sc.cz

- new cluster purchased by CERIT-SC
 - available through "phi" queue (PBS Pro) on zuphux.cerit-sc.cz frontend

```
zuphux$ qsub -q phi -l select=...
```

- the newest generation of Xeon Phi (7210 Knights Landing)
- see more details at https://metavo.metacentrum.cz/export/sites/meta/cs/seminars/seminar2017/meta-xeonphi-17.pdf

Installation specifics:

- /storage/brno3-cerit is the only central storage directly mounted
 - all the rest storages available through SCP only







Central storages of phi.cerit-sc.cz cluster

Central storages not available through NFS

i.e. /storage/XXX/home/<username>

- technical reasons
- data storages available through SCP
 - besides brno3-cerit, which is mounted directly
 - in most cases, just a minor change in your scripts

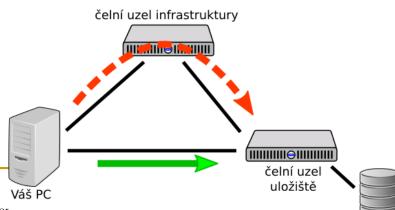
NFS sharing (most clusters)	SCP sharing (phi[1-6].cerit-sc.cz)
DATADIR="/storage/brno2/home/ <username>/example"</username>	DATADIR="storage-brno2.metacentrum.cz:~/example"
cp -R \$DATADIR/mydata \$SCRATCHDIR	scp –R \$DATADIR/mydata \$SCRATCHDIR







- how to transfer large amount of data to computing nodes?
 - copying through the frontends/computing nodes may not be efficient (hostnames are storage-XXX.metacentrum.cz)
 - XXX = brno2, brno3-cerit, plzen1, budejovice1, praha1, ...
 - \rightarrow connect directly to the storage frontends (via **SCP** or **SFTP**)
 - \$ sftp storage-brno2.metacentrum.cz
 - \$ scp <files> storage-plzen1.metacentrum.cz:<dir>
 - etc.
 - use FTP only together with the Kerberos authentication
 - otherwise insecure









how to get information about your quotas?

- by default, all the users have quotas on the storage arrays (per array)
 - may be different on every array
- to get an information about your quotas and/or free space on the storage arrays
 - textual way: log-in to a MetaCentrum frontend and see the "motd" (information displayed when logged-in)
 - graphical way:
 - your quotas: https://metavo.metacentrum.cz/cs/myaccount/kvoty
 - free space: http://metavo.metacentrum.cz/pbsmon2/nodes/physical

how to restore accidentally erased data

- □ the storage arrays (⇒ including homes) are regularly backed-up
 - several times a week
- □ → write an email to <u>meta@cesnet.cz</u> specifying what to restore







- how to secure private data?
 - by default, all the data are readable by everyone
 - □ → use common Linux/Unix mechanisms/tools to make the data private
 - r,w,x rights for user, group, other
 - e.g., chmod go= <filename>
 - see man chmod
 - □ use "¬R" option for recursive traversal (applicable to directories)
- how to share data among working group?
 - ask us for creating a common unix user group
 - user administration will be up to you (GUI frontend is provided)
 - use common unix mechanisms for sharing data among a group
 - see "man chmod" and "man chgrp"
 - see https://wiki.metacentrum.cz/wikiold/Sdílení dat ve skupině







- how to use SGI UV2000 nodes? (ungu, urga .cerit-sc.cz)
 - because of their nature, these nodes are not by default used
 by common jobs
 - to be available for jobs that really need them
 - to use these nodes, one has to submit the job to a specific queue called "uv"
 - \$ qsub -l select=1:ncpus=X:mem=Yg -q uv
 -l walltime=Zd ...
 - to use a specific UV node, submit e.g. with
 \$ qsub -q uv -l select=1:ncpus=X:cl_urga=true ...
 - □ for convenience, submit from zuphux.cerit-sc.cz frontend







- how to run a set of (managed) jobs?
 - some computations consist of a set of (managed) sub-computations
 - optional cases:
 - the computing workflow is known when submitting
 - specify dependencies among jobs
 - qsub's "-w" option (man qsub)
 - □ in case of many parallel subjobs, use "job arrays" (qsub's "-J" option)
 - see https://www.pbsworks.com/pdfs/PBSUserGuide13.0.pdf, page 209
 - the computing workflow depends on result(s) of subcomputations
 - □ run a master job, which analyzes results of subjobs and submits new ones
 - the master job should be submitted to a node dedicated for lowperformance (controlling/re-submitting) tasks
 - available through the "oven" queue
 - qsub -q oven -l select=1:ncpus=... control script.sh







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What to do if something goes wrong?

- 1. check the MetaVO/CERIT-SC documentation, application module documentation
 - whether you use the things correctly
- 2. check, whether there haven't been any infrastructure updates performed
 - visit the webpage http://metavo.metacentrum.cz/cs/news/news.jsp
 - one may stay informed via an RSS feed
- 3. write an email to meta@cesnet.cz, resp. support@cerit-sc.cz
 - your email will create a ticket in our Request Tracking system
 - identified by a unique number → one can easily monitor the problem solving process
 - please, include as good problem description as possible
 - problematic job's JOBID, startup script, problem symptoms, etc.







Overview

- Introduction
- MetaCentrum / CERIT-SC infrastructure overview
- How to ... specify requested resources
- How to ... run an interactive job
- How to ... use application modules
- How to ... run a batch job
- How to ... determine a job state
- Another mini-HowTos ...
- What to do if something goes wrong?
- Real-world examples
- Appendices







Real-world examples

Examples:

- Maple
- Gaussian + Gaussian Linda
- Gromacs (CPU + GPU)
- Matlab (parallel & GPU)
- Ansys CFX
- OpenFoam
- Echo
- R Rmpi

demo sources:

/storage/brno2/home/jeronimo/MetaSeminar/latest

command: cp -rH /storage/brno2/home/jeronimo/MetaSeminar/latest \$HOME











Projekt CERIT Scientific Cloud (reg. no. CZ.1.05/3.2.00/08.0144) byl podporován operačním programem *Výzkum a vývoj pro inovace*, 3 prioritní osy, podoblasti 2.3 *Informační infrastruktura pro výzkum a vývoj.*

www.cesnet.cz

www.metacentrum.cz

www.cerit-sc.cz

EUROPEAN REGIONAL DEVELOPMENT FUND

INVESTING IN YOUR FUTURE







Overview

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Appendices

- Common mistakes in computations
- How to deal with parallel/distributed computations?
- Other computing possibilities
 - MetaCloud
 - Hadoop (MapReduce)
 - Specialized frontends Galaxy, Chipster, ...







Common mistakes in computations







Common mistakes in computations

Feel free to use the infrastructure – if something crashes, it's our fault. ©







Big data transfers

Do not copy higher amounts of data through frontends

- slower transfer
- frontends load







Big data transfers

Do not copy higher amounts of data through frontends

- slower transfer
- frontends load

Data could be copied directly through storage frontends

SCP, WinSCP
 /storage/brno2 -> storage-brno2.metacentrum.cz
 /storage/brno3-cerit -> storage-brno3-cerit.metacentrum.cz

- - -

 https://wiki.metacentrum.cz/wiki/Working_with_data/Direct_a ccess_to_data_storages







Big data transfers

Do not copy higher amounts of data through

frontends

- slower transfer
- frontends load

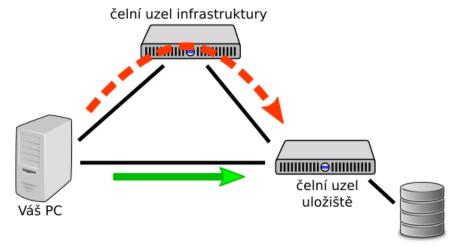
Data could be copied directly through storage frontends

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. . .

 https://wiki.metacentrum.cz/wiki/Working_with_data/Direct_a ccess_to_data_storages









Computations and central storages

Do not run computations that compute over data located at central storages

- especially the I/O-intensive ones
 - increses central storage load and makes the computation slower

Compute over local copies in scratch directories

- benefits:
 - faster computations
 - computations do not rely on the availability of the central storage
- usage:
 - \$ qsub -I select=1:ncpus=4:scratch_local=1gb ...
 cp /storage/.../home/<username>/mydata \$SCRATCHDIR/mydata
 cd \$SCRATCHDIR
 <compute>
 cp \$SCRATCHDIR/results /storage/.../home/<username>/results
 - ...:scratch_shared=Xgb ... shared scratch (distributed computations)
 - ...:scratch_ssd=Xgb ... local scratch SSD disks







Data in scratches

Clean the scratches once computations finish

- scratch data could be seen similarly as data in RAM memory
 - once a computation finishes, the data should be cleaned
- scratches are automatically cleaned by us
 - usually after 2 weeks a computation finishes

Clean scratch after your computations

- "clean_scratch" utility
- usage (in script file): trap 'clean_scratch' TERM EXIT

cp results /storage/... || export CLEAN_SCRATCH=false

- if the central storage is not available (the results could not be copied out), the data will remain in the particular scratch directory
 - o user is informed about (non-)useful scratch cleanup
 - user is informed about scratches not correctly cleaned after previous computations







Central storages overused

Central (working) storages are not infinitely large ⊗

/storage/<CITY>

Clean/move currently unnecessary data

- possibilities:
 - delete unnecessary data
 - move actually unnecessary data to archival storages
 see https://wiki.metacentrum.cz/wiki/Archival_data_handling







Huge jobs' outputs and data in /tmp

Computing nodes restrict the amount of data a user is able to store to local disks (outside the scratch space) = 1 GB quota

- influences the /tmp direktory (temporal files)
- influences the huge job's outputs (stdout, stderr)

Store bigger amounts of data to scratches

- forwarding the temporal directory
 many applications follow the system variable TMPDIR
 - usage: export TMPDIR=\$SCRATCHDIR
- forwarding the stdout/stderr of an application
 - myapp ... 1>\$SCRATCHDIR/stdout 2>\$SCRATCHDIR/stderr
- checking the status of your local quota and a list of files occupying it (once being informed by email)
 - utility \$ check-local-quota has to be run on the particular node (with exhausted local quota)







Non-effective computations

Be aware of the resource usage effectivity of your jobs

- a request for multiple CPUs/cores will not make a single-processor (single-thread) computation parallel (= it won't be faster)
 - just a single CPU will be used
- many applications significantly vary between the number of CPUs used throughout a computation
 - higher number of CPUs might be used just for a short time of the computation

Observing the computation usage of (not only) CPUs:

- during a computation:
 - log-in to the computation node (SSH) and use standard Linux tools (top, htop, ...)
- after a computation:
 - see the list of jobs at the MetaCentrum portal (https://metavo.metacentrum.cz/cs/myaccount/myjobs.html) the non-effective jobs have red background color







Infiniband

Distributed jobs might run ineffectively because of slow communication channel

- the inter-process communication using standard network services (Ethernet) is slow
- Infiniband specialized low-latency interconnect for fast interprocess communication in distributed computations

Most of our clusters are equipeed with Infiniband

- considerably accelerates the performance of distributed (MPI) computations
 - the Infiniband availability is automatically detected computations started always in the same way: mpirun myapp
 - if the Infiniband is not available, Ethernet is used as a fallback
- request:
 - \$ qsub –l select=... –l place=group=infiniband script.sh







Many short-term jobs processing

Group/gather short-term jobs

- e.g., the ones running less than few minutes
 - startup overhead may be a significant part of the whole processing time results in wasting resources

Run more computations within a single job

- possibilities:
 - serial computations run inside a single job process data1 process data2

. . .

- parallel computations run inside a single job (necessary to ask for enough CPUs)
 - o pbsdsh
 - o parallel







Computations on frontends

Do not run computations on frontends

- neither for computations nor for complex results analyses
 - increased frontend load results in limitation of its services (and usually frontend crash)
- frontend's primary job is jobs' preparation and very simple and short-term computations

Use interactive jobs

- request:
 - \$ qsub -I -l select=...
- usage possibilities:
 - textual mode
 - graphical mode VNC access \$ module add gui

 - \$ qui start
 - see https://wiki.metacentrum.cz/wiki/Remote_desktop







Interactive jobs

Minimize the time lags in interactive jobs

- especially the time between job startup and your work (starting computations)
 - -> in the time lag, the resources are wasted

Stay informed about your job's startup

- request:
 - \$ qsub -m ab -I -l select=...
 will send you an email once the job begins
 (,,-m abe" also in the case of job's finish)
- these options could be also used in batch jobs
 but be aware of running too many jobs with this option set!
 - overloads your mailbox
 - may blacklist our mailservers at external mail providers ©







Cloud nodes

Be aware about your VMs running

- even the unused VMs/nodes (but running) consume infrastructure resources
 - -> results in wasting resources, which somebody could use

Terminate/Suspend unused VMs

 we'll regularly inform you about your VMs running in case of no response (= time extension), the VMs are terminated







How to deal with parallel/distributed computations?







How to ... run a parallel/distributed computation I.

Parallel jobs (OpenMP):

- if your application is able to use multiple threads via a shared memory, ask for a single node with multiple processors
 - \$ qsub -l select=1:ncpus=...
- make sure, that before running your application, the OMP_NUM_THREADS environment variable is appropriately set
 - otherwise, your application will use all the cores available on the node
 - → and influence other jobs...
 - usually, setting it to NCPUs is OK
 - \$ export OMP NUM THREADS=\$PBS NUM PPN







How to ... run a parallel/distributed computation II.

Distributed jobs (MPI):

 if your application consists of multiple processes communicating via a message passing interface, ask for a set of nodes (with arbitrary number of processors)

```
$ qsub -1 select=...:ncpus=...
```

- make sure, that before running your application, the appropriate openmpi/mpich2/mpich3/lam module is loaded into the environment
 - \$ module add openmpi
 - then, you can use the mpirun/mpiexec routines
 - \$ mpirun myMPIapp
 - it's not necessary to provide these routines neither with the number of nodes to use ("-np" option) nor with the nodes itself ("--hostfile" option)
 - the computing nodes are automatically detected by the openmpi/mpich/lam







How to ... run a parallel/distributed computation III.

Distributed jobs (MPI): accelerating their speed I.

- to accelerate the speed of MPI computations, ask just for the nodes interconnected by a low-latency Infiniband interconnection
 - all the nodes of a cluster are interconnected by Infiniband
 - there are several clusters having an Infiniband interconnection
 - mandos, minos, hildor, skirit, tarkil, nympha, gram, luna, manwe (MetaCentrum)
 - zewura, zegox, zigur, zapat (CERIT-SC)
- submission example:

```
$ qsub -1 select=4:ncpus=2 -1 place=group=infiniband MPIscript.sh
```

- starting an MPI computation using an Infiniband interconnection:
 - □ in a common way: \$ mpirun myMPIapp
 - the Infiniband will be automatically detected
 - is the Infiniband available for a job? check using \$ check-IB

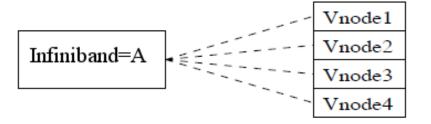


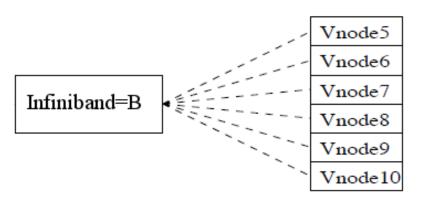


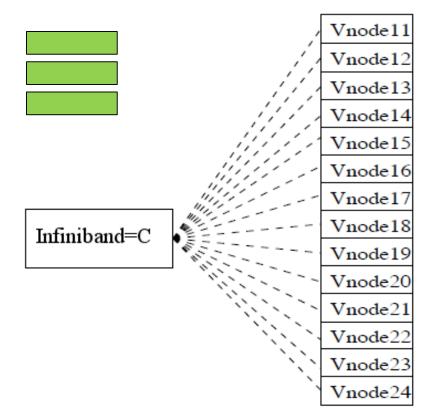


Chunks grouping

- accelerating distributed jobs
- -l place=group=infiniband







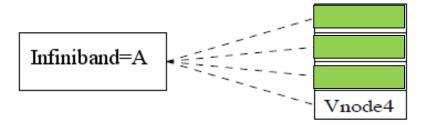


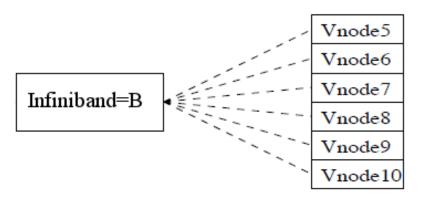


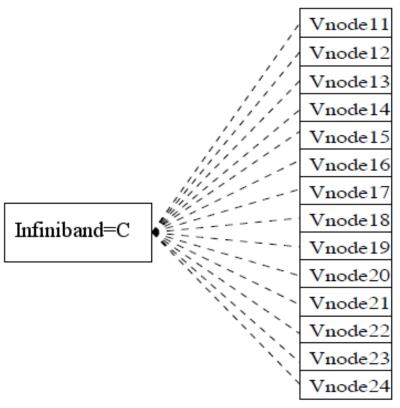


Chunks grouping

- accelerating distributed jobs
- -1 place=group=infiniband







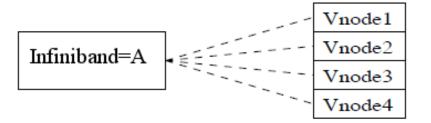


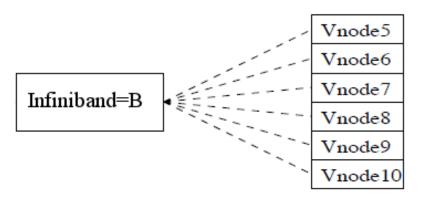


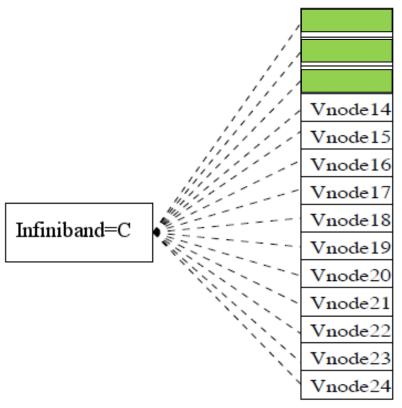


Chunks grouping

- accelerating distributed jobs
- -1 place=group=infiniband









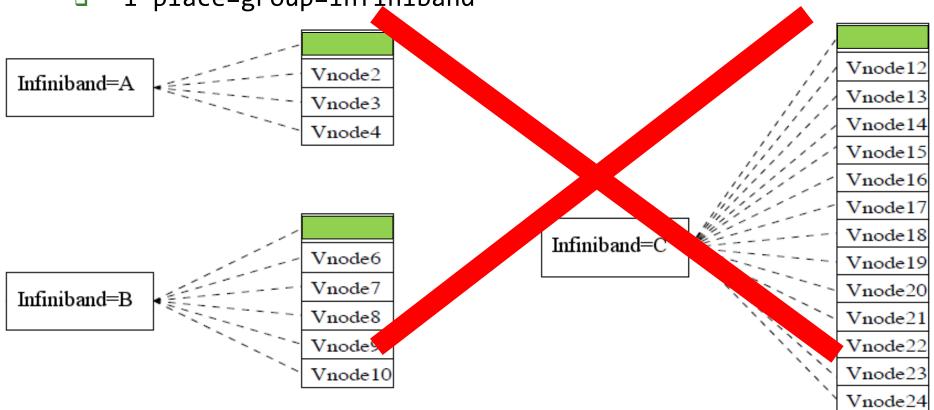




Chunks grouping

accelerating distributed jobs

-l place=group=infiniband









How to ... run a parallel/distributed computation IV.

Questions and Answers:

- Is it possible to simultaneously use both OpenMP and MPI?
 - Yes, it is. But be sure, how many processors your job is using
 - appropriately set the "-np" option (MPI) and the OMP_NUM_THREADS variable (OpenMP)
 - OpenMPI: a single process on each machine (mpirun -pernode ...) being threaded based on the number of processors (export OMP_NUM_THREADS=\$PBS_NUM_PPN)

Any other questions?









Other computing possibilities

Cloud computing - MetaCloud







Grid vs. Cloud computing

Grid computing suitable for:

- long-term and/or large-scale computations
 - (primarily batch processing)
- applications not requiring special OSs (features)
 - pre-installed or users' ones

Cloud computing suitable for:

- applications requiring special environment (OS) and/or features
 - one can run various operating systems (incl. Windows OS) and/or application equipment
 - administrator/root access is provided







Cloud computing

How to compute?

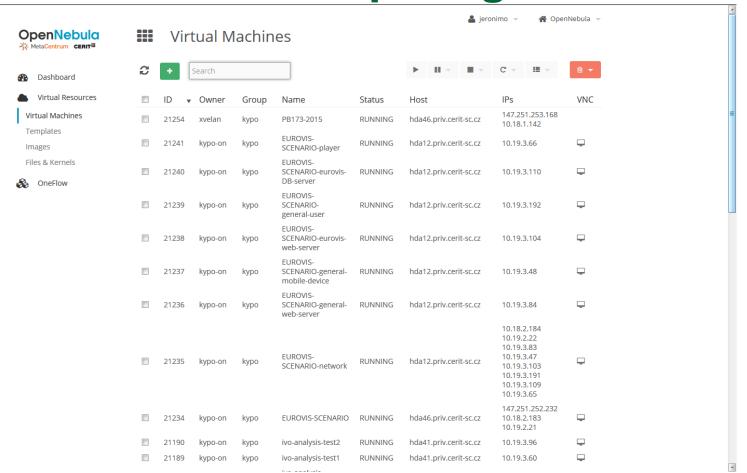
- additional registration to MetaCloud group required
 - your SSH key is needed to access the VMs
 - https://perun.metacentrum.cz/fed/registrar/?vo=meta&group=metacloud
- OpenNebula GUI for deployed nodes management
 - https://cloud.metacentrum.cz/
- interact via:
 - cloud/VM console
 - internal OS services (SSH, VNC, Rdesktop, ...)







Cloud computing









Basic terminology

- template a description of the node ("computer") you are asking from the cloud
 - o speficies requested CPUs, memory, disk storage, network interfaces, etc.
 - predefined ones & own ones
- image a storage space ("HDD")
 - equipped with an OS or not
 - persistent & non-persistent (default; data are lost when destroying the VM)
- network security group a set of firewall rules
 - restricts the access to the VM
- instance a running node/computer created based on some template receipt
- VM console a VNC connection to the VM ("computer screen")







Common usage

Common operations with MetaCloud:

- see <u>video tutorial</u>
- □ for advanced use, see MetaCloud documentation
 - e.g., creating your own template (duplicate existing one) or disk image

Warning: Please, be aware of the VMs you are running.

And if not used, suspend or terminate them...

- every 3 months, we'll recommend you your running VMs
 - if not explicitly renewed/extended in the defined time period, the VMs will be terminated







Other computing possibilities

Hadoop computing







Hadoop / MapReduce computing

Hadoop:

- an open-source framework for distributed storage and distributed processing of large volumes of data
 - large data blocks splitted and distributed amongst nodes
 - a MapReduce-based algorithm (= data processing code) is distributed over the distributed blocks and processed in parallel

Suitable for:

- huge datasets to be processed
 - but NOT suitable for arbitrary data processing one can imagine
 - just for the processing meeting the MapReduce programming model
 - e.g., counting the number of times worlds of times world of times







Other computing possibilities

Specialized frontends - Galaxy, Chipster, ...







Specialized frontends/environments

Suitable for:

- user communities with well-defined processing needs
 - workload & computing pipeline orchestrators available via GUI
 - usually adapted to user needs and/or because of interoperability with our infrastructure
 - in background, the proper computing method is used (grid, cloud, etc.)

How to compute?

- Galaxy (docs https://wiki.metacentrum.cz/wiki/Galaxy)
 - ELIXIR RepeatExplorer Galaxy https://repeatexplorer-elixir.cerit-sc.cz
 - MetaCentrum Galaxy https://galaxy.metacentrum.cz
- Chipster (docs https://wiki.metacentrum.cz/wiki/Chipster)
 - http://chipster.metacentrum.cz:8081
- basic details are provided in the leaflet