

Introduction to STAR software and makers

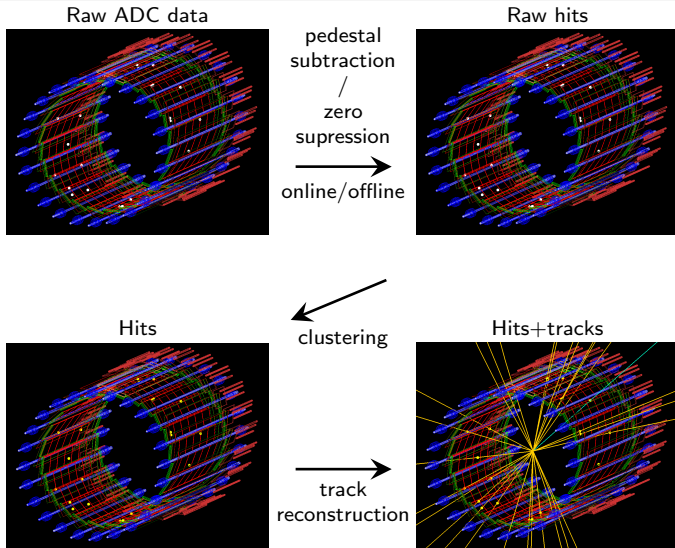
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STAR Collaboration Meeting 2015
Stony Brook University, 1-6.6.2015

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Reconstruction in STAR - Example with IST



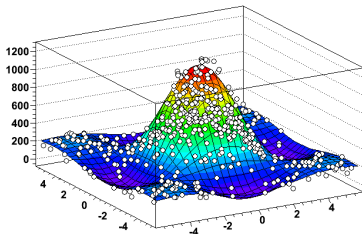
Each operation is done by its corresponding maker (pedestal subtraction, clustering...)

STAR software

Most of STAR software is based on libraries of ROOT data analysis framework.

- ROOT is widely used in Heavy Ion and High Energy Physics community
- STAR ROOT libraries contain classes specific to STAR:
 - Detector geometry classes
 - Event data
 - Specific makers
 - Database classes
- Look for information and class description on the ROOT webpage.

Minuit fit result on the Graph2DErrors points



ROOT webpage

<https://root.cern.ch/drupal/>

Listing 1 : To run STAR ROOT

```
root4star
```

STAR Database

- Contains calibrations, gains, geometry transformations, etc.
- Access to Database information is done via `StMaker::GetDataBase()` method which calls `St_db_Maker`.
- `St_db_Maker` reads DB tables with a timestamp automatically set from the data or with `SdtYYYYMMDD.HHMMSS` command for simulation

DB tables are read in order:

- 1 `./StarDb` - local directory
- 2 `$STAR/StarDb` - from current STAR Library version through AFS
- 3 MySQL database

This allows to substitute official Db entries by a private version in order to check/modify calibration.

More information on STAR Database

<https://drupal.star.bnl.gov/STAR/comp/db>

- STAR software is constantly being updated as software developers make modifications or new detectors are added (HFT)
- Available remotely through an AFS cell
- The versions of STAR Libraries are labeled (listed only few):
 - SLYYx - where YY is the year, and x is a letter, corresponds to data production PYYix
 - DEV - latest development version, may be buggy or unstable
 - PRO - latest production version, default after login
 - NEW - a few weeks old version for full integration testing, more stable than DEV



```
starver SL15c // To select production P15ic:  
stardev // To select DEV version  
starpro // To select production version  
starnew // to select NEW version
```

Listing 2 : Don't use above commands in login scripts!! Use command below instead:

```
if (${STAR_LEVEL} == 0) setenv STAR_LEVEL XXX
```

Library release structure and policy

<https://drupal.star.bnl.gov/STAR/comp/sofi/soft-n-libs/release>

CVS repository

- An archive of all codes and changes to the STAR software
- Source of codes for users

Listing 3 : To get a local copy of code from CVS repository:

```
cvs co <path_to_files>
```

Listing 4 : Compile source code:

```
cons
```

Quick guide to CVS in STAR

<https://drupal.star.bnl.gov/STAR/comp/sofi/tutorials/cvs>

CVS repository online viewer

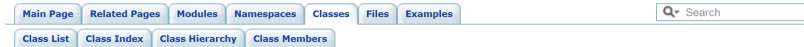
<http://www.star.bnl.gov/cgi-bin/protected/cvsweb.cgi/>

Doxygen

- Automatically generates online code documentation
- Good for quick reference
- Contains links to the source code

Doxygen webpage

<http://www.star.bnl.gov/webdata/dox/html/index.html>



StMcEvent Class Reference

Event data structure to hold all information from a Monte Carlo simulation. This class is the interface to access all information from the GEANT simulations in STAR. From here one can get all vertices, tracks, hits in the event. [More...](#)

Inheritance diagram for StMcEvent:



[List of all members.](#)

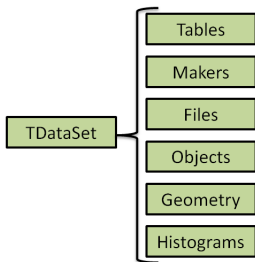
Public Member Functions

	StMcEvent (g2t_event_st *)
virtual void	Browse (TBrowser *b)
virtual bool	IsFolder () const
int	operator== (const StMcEvent &) const
int	operator!= (const StMcEvent &) const
unsigned long	eventGeneratorEventLabel () const

Dataset

Dataset is a named list of dataset objects which allows to organize themselves in a tree. These objects can be almost any data!

Figure : Dataset can be a set of any of these:



TDataSet

- TDataSet (base class) is a part of basic ROOT
- Unique names for datasets are important, because they are accessed by their name.
- Datasets form a basis of STAR software for:
 - Data analysis
 - Simulation
 - Reconstruction

Listing 5 : Run an example chain and check it yourself!

```
root4star -l 'bfc.C(1,1,"gstar")'
```

Figure : 1. Open a TBrowser and see a list of datasets

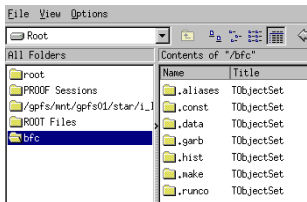


Figure : 2. Click on .make and check a list of makers

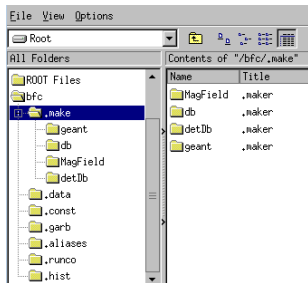


Figure : 3. Click on a maker for a list of it's datasets

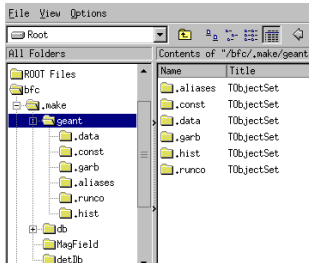
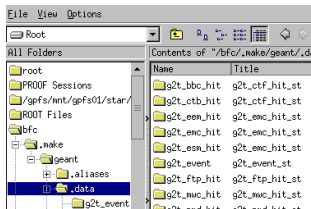


Figure : 4. Click on .data to see a list of data tables



Maker

- Maker is an owner of a dataset.
- StMaker (base class) is also a TDataSet itself.
- Makes operations on data:
 - Clustering
 - Applying corrections
 - Reconstruction
 - Data analysis
 - More...

Contents of a maker

- .make - list of child makers
- .data - list of owned datasets
- .hist - list of histograms created in StMaker::Init();
- .const - collection of constants used by the maker

Listing 6 : User defined functions - overwrite them with your own code

```
virtual Int_t Init(); // Called at the beginning, initialize histograms etc...
virtual Int_t InitRun(); // Called at the start of a new run
virtual Int_t Make(); // Called for every event, do your analysis here, fill histograms
virtual Int_t Clear(); // Used for clearing collections after each Make();
virtual void Finish(); // Finish
virtual Int_t FinishRun(Int_t oldrunnumber); // Executed at the end of a run
```

Listing 7 : Get methods

```

virtual TDataSet *GetData(const char *name, const char *dir=".data") const;
virtual TDataSet *GetDataSet(const char *logInput) const {return FindDataSet(logInput);}
virtual TDataSet *DataSet(const char *logInput) const
    {return GetDataSet(logInput)};
virtual TDataSet *GetInputDS (const char *logInput) const
    {return GetDataSet(logInput)};

virtual TDataSet *GetDataBase(const char *logInput, const TDateTime *td=0);
virtual TDataSet *GetInputDB(const char *logInput)
    {return GetDataBase(logInput)};

```

Listing 8 : Add methods

```

virtual void AddData (TDataSet *data, const char *dir=".data");
virtual void AddConst(TDataSet *data=0){AddData(data, ".const");}
virtual void AddHist(TH1 *h, const char *dir=0);
virtual void AddGarb (TDataSet *data=0){AddData(data, ".garb");};
virtual void AddRunco (TDataSet *data=0){AddData(data, ".runco");};
virtual void AddRunco (Double_t par, const char *name, const char *comment);
void AddRunCont (TDataSet *data=0){AddRunco(data);}; //alias

```

Chain

- A special StMaker (“Top maker”) and a TDataSet as well
- Contains an ordered list of makers
- Makers are added to the chain automatically in order they are created
- Allows to combine makers as components of a complex analysis, reconstruction or simulation

Listing 9 : Example MuDST analysis chain

```

StChain *chain = new StChain;
StMuDstMaker *muDstMk = new StMuDstMaker(0, 0, "", file_list.data(), ".MuDst.root", 200, "MuDst"); // reads MuDST
St_db_Maker *dbMk = new St_db_Maker("StarDb", "MySQL:StarDb"); // DB maker
StEEmcDbMaker *eemcDb = new StEEmcDbMaker("EEmcDbMaker"); // Additional EEMC DB maker

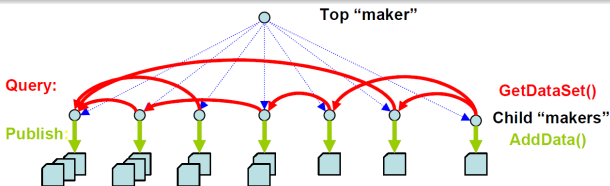
StAnalysisMaker *analysis = new StAnalysisMaker(muDstMk); // Pass StMuDstMaker pointer to your analysis maker

Bool_t status = chain->Init(); // Call Init() for all makers
if (status != kStOK) {
cout << "ERROR: Chain initialisation failed (error " << status << "), not running!" << endl;
return 0;
}
status = chain->EventLoop(); //Process all events in the chain
if ((status != kStOK) && (status != kStEOF)) cout << "WARNING: EventLoop returned error " << status << endl;
status = chain->Finish(); // Call Finish() for all makers, write output data
if (status != kStOK) cout << "WARNING: Chain finished with error " << status << endl;
delete chain;

```

Communication between makers

- Provider consumer model
 - Makers are both providers and consumers of data
 - The order of communication is fixed (providers before consumers) and is controlled by the order of makers instantiation
- 1 Maker “consumes” dataset found in a chain by its name
 - 2 Maker fills the output dataset and “provides” it to the chain



Each maker can **query** the data **published** by the other makers.

PoS ACAT08 040 (2008)

Listing 10 : Communication functions

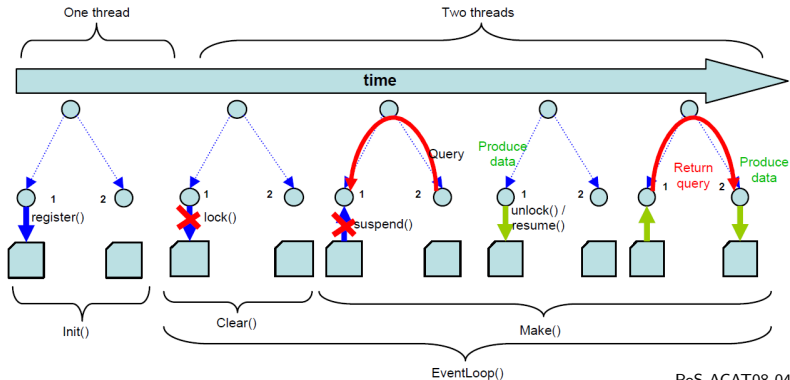
```
virtual TDataSet* GetDataSet(const char *logInput) const // Get dataset by its name
virtual void AddData(TDataSet *data, const char *dir=".data") // Publish data
```

Threads

- Not implemented currently due to limited parallelization - but potentially available
- Makers can run in different threads to take advantage of multi-core CPUs
- `StMaker::Make()` function executed in parallel
- Pipeline-like execution

Pipeline-like execution

- Maker 1 finishes `Make()` for event 1 and passes data to the maker 2
- Maker 1 executes `Make()` for event 2, while maker 2 executes `Make()` for event 1
- ...



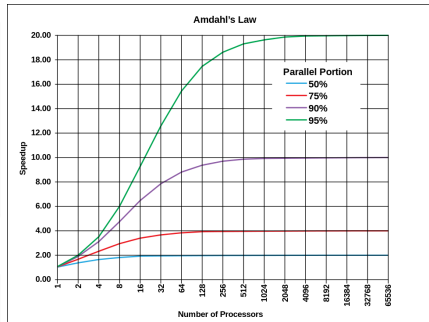
Amdahl's Law

$$S(N) = \frac{1}{(1-P) + \frac{P}{N}}, \text{ where:}$$

- $S(N)$ - speedup compared to executing a serial code on a single core, in fact a **theoretical limit**
- P - fraction of the code, which can be made parallel
- N - number of processors used

Think about efficiency!

- Embarassingly parallel mode - we run independent serial jobs (1 job = 1 core) over multiple data (CPU use efficiency is 98%! in case of reconstruction)
- Now with 60% parallel code run on 16 cores, we get a maximum speedup of 2.28, but we occupy 16 cores for that! The efficiency is 14% compared to embarassingly parallel.
- More resources used, so efficiency drops!
- Tracking can be parallelized, but it takes 60% of the time.
- Parallel threads implementation abandoned in favor of high efficiency.



$E(N)$ - efficiency compared to embarassingly parallel

Theoretical maximum values for 16 cores

P	$S(16)$	$E(16)$
99%	13.9	87%
95%	9.14	57%
60%	2.28	14%

Reconstruction requires running a few makers per subsystem(detector):



Total number: ~ 100 makers

- Easy with Big Full Chain(BFC), `StBFChain` class
- Just choose options you need(subsystems and corrections) and go
- BFC can run reconstruction for data or simulation with makers in predefined, fixed order
- Option and maker dependencies form a tree eg:
istHit option runs `StIstHitMaker`,
which requires **istCluster**, which runs `StIstClusterMaker`,
which also requires **istRaw** and runs `StIstRawHitMaker` and **istDB**, running `StIstDbMaker`
This forms IST reconstruction chain (but no tracking!) and ensures no maker is missing (in most cases)

More details on BFC here:

https://drupal.star.bnl.gov/STAR/system/files/BFC_HQPC_LK_2015_2_12.pdf

After setting options, BFC:

- 1 Creates ordered list of makers
- 2 Loads shared libraries
- 3 Runs

If no options are set it will print the list of available options

Listing 11 : How to run BFC?

```
root4star bfc.C
root -l bfc.C
```

Listing 12 : BFC run arguments

```
root4star -l -b -q bfc.C(First, Last, *Chain, *infile) | tee bfc.log

bfc(Int_t First, Int_t Last, // first and last event
    const Char_t *Chain, // BFC options, separated by spaces or commas, case insensitive, order
                        // irrelevant, put "-" in front to turn off
    const Char_t *infile, // input file .daq (data) or .fzd (simulation)
    const Char_t *outfile, // output file
    const Char_t *TreeFile)
}
```

List of all BFC options

<http://www.star.bnl.gov/cgi-bin/protected/cvsweb.cgi/~checkout~/StRoot/StBFChain/doc/index.html>

MuDST analysis

- Write a macro similar to one shown in Listing 9
- Create a chain, StMuDstMaker and your own maker in this macro
- Pass file list to the StMuDstMaker
- The macro is used to run the analysis chain
- Write your own analysis maker and put it in StRoot/StXxxMaker
- Compile (cons) and run macro when finished

Listing 13 : In the Make() function of your maker

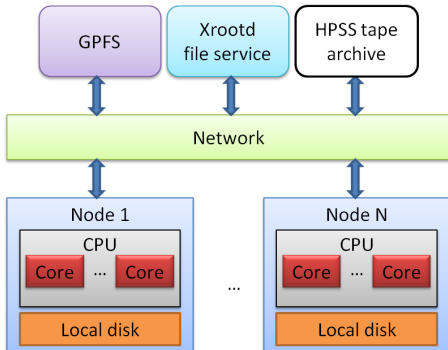
```
StMuDst* mMuDst = mMuDstMaker->muDst(); // Get StMuDst
StMuEvent* muEvent = mMuDst->event(); // Get StMuEvent
mMuDst->primaryVertex()->position(); // Get primary vertex position
mMuDst->primaryVertex()->refMult(); // Get reference multiplicity
TObjArray* tracks = mMuDst->primaryTracks(); // Create a TObjArray containing the primary
    tracks
TObjArrayIter GetTracks(tracks); // Create an iterator to step through the tracks
StMuTrack* track; // Pointer to a track
while((track = (StMuTrack*)GetTracks.Next())) // Main loop for iterating over tracks
    {histogram[1]->Fill(track->pt());} // Do analysis, fill histograms
```

MuDST tutorial

<http://rnc.lbl.gov/~jthomas/public/MuDSTutorial2011.pdf>
<http://www.star.bnl.gov/public/comp/train/tut/MuDSTutorial06/>

RHIC Computing Facility (RCF)

- Made of multiple nodes connected by network (657 nodes, 15,576 cores)
- Each node has CPU with multiple cores, so a few jobs can run on the same node (1 core = 1 job)
- Provides resources to analyze massive amounts of data in parallel on different nodes
- Nodes accessed through a gatekeeper `rssh.rhic.bnl.gov`



Remote desktop NX server

<https://www.racf.bnl.gov/docs/services/nx>

Nodes have access to:

- Local disk, the fastest - use these!
- General Parallel File System (GPFS), disk, slower access than local disk
- High Performance Storage System (HPSS), tape archive, the slowest but stores all data
- Files using Xrootd file server, which redirects connections to the physical location of the file

RCF nodes monitoring table

<http://www.star.bnl.gov/cgi-bin/protected/nova/showMachines.pl>

File catalog

- Allows to locate files and to create file lists
- Very flexible, can return number of events or calculate total number

Listing 14 : Usage

```
get_file_list.pl [-all] -keys keyword[,keyword,...] \  
  [-cond keyword=value[,keyword=value,...]] [-start #] [-limit #] [-delim $St] \  
  [-onefile] [-o outputfile]
```

Listing 15 : Example - get MuDST files from Run 14

```
get_file_list.pl -keys 'path,filename' -cond 'trgsetupname=AuAu_200_production_low_2014,  
  available=1,filetype=daq_reco_MuDst,sanity=1,filename~st_physics_14,storage=local' -  
  onefile -delim '/'
```

File catalog manual:

<https://drupal.star.bnl.gov/STAR/comp/sofi/filecatalog/user-manual>

STAR Scheduler - SUMS

- Used for submitting batch jobs to RCF cluster
- Automatically distributes input files between multiple jobs

Listing 16 : Typical XML job description file

```
<?xml version="1.0" encoding="utf-8" ?>
<job minFilesPerProcess="14" maxFilesPerProcess="20" filesPerHour="3" fileListSyntax="xrootd">
<command> starver SL11d
  cons
  root4star -q -b RunTrackSampleMaker.C\{1e8,\"$FILELIST\", \"$SCRATCH\", \"$JOBID\" \}
</command>
<SandBox installer="ZIP"> <Package> <File>file:./RunTrackSampleMaker.C</File>
  <File>file:./StRoot</File> </Package> </SandBox>
<input URL="catalog:star.bnl.gov?production=P11id,trgsetupname=pp500_production_2011,filetype=daq_reco_MuDst,sanity
  =1,storage!=HPSS" nFiles="all"/>
<stdout URL="file:./output/stdout/$JOBID.out"/>
<stderr URL="file:./output/stderr/$JOBID.err"/>
<output fromScratch="*.root" toURL="file:/star/institutions/lbl/lkosarz/Upsilon/nSigmaCode/output/output/" />
</job>
```

```
star-submit jobs.xml // Submit jobs described in jobs.xml file
```

Scheduler manual and FAQ:

<http://www.star.bnl.gov/public/comp/Grid/scheduler/>
<http://www.star.bnl.gov/public/comp/Grid/scheduler/faq.html>

Good practice

- Test your code locally in interactive mode on a small data sample.
- Once tested submit jobs to RCF.
- Write output of jobs locally to avoid unnecessary traffic over the network.
- Turn off unused MuDST branches to speed up analysis.

Listing 17 : Turn of MuDST branches in the run macro

```
mudst_mk->SetStatus("*",0); // Turn off all MuDST branches
mudst_mk->SetStatus("MuEvent",1); // Turn on MuEvent branch
mudst_mk->SetStatus("PrimaryTracks",1); // Turn on Primary Tracks
```

Turning off MuDST branches

http://www.star.bnl.gov/public/comp/train/tut/MuDstTutorial/select_branches.html

STAR coding and naming standards - **Update in preparation!** - see Jerome's talk

<https://drupal.star.bnl.gov/STAR/comp/sofi/soft-n-libs/standards>
<https://drupal.star.bnl.gov/STAR/blog/jerome/c11-and-star-coding-standard-committee-charges>

More tutorials

<https://drupal.star.bnl.gov/STAR/comp/sofi/tutorials>

- Presented STAR software basics and crucial topics
- Good start for juniors
- Included useful links
- Everything you need to start your analysis

Thank you!

BACKUP