# Femtocode: query system for HEP

# What's a query system?

User asks a question, gets an answer quickly enough to *explore* the data.

Like a Google query, but aggregating HEP data, returning (e.g.) histograms.

Embedded within an analysis script: provides sliced projections of the data for users to fit/plot/analyze in any way they want.

### How it differs from what we do now

Physicists arrange data as sets of files that have to be filtered into progressively smaller sets of files until the final set is small enough for real-time data analysis.

Instead, we propose a service that serves aggregated views of analysis object data on demand.

Must be responsive to requests in realtime: ~1 sec for each scan over a dataset.





## High-level language

User writes expressions that pick apart the structure of objects within arbitrary-length lists, to any depth of nesting.

Higher-order functions like .map, .pairs, .filter, .reduce instead of explicit for loops.

Femtocode query language is distributed in quoted snippets throughout a structured workflow and tree of aggregators.

(See http://histogrammar.org for histogram abstraction.)

workflow = session.source("b-physics")	# pull from a named dataset
.define(goodmuons = "muons.filter(\$1.pt > 5)")	# muons with pt > 5 are good
.filter("goodmuons.size >= 2")	# keep events with at least two
.define(dimuon = """	
<pre>mu1, mu2 = goodmuons.maxby(\$1.pt, 2);</pre>	# pick the top two by pt
energy = mu1.E + mu2.E;	<pre># compute imploded energy/momentum</pre>
px = mu1.px + mu2.px;	
py = mu1.py + mu2.py;	
pz = mu1.pz + mu2.pz;	

## Low-level execution

#### No objects at runtime:

All nested structures are represented as homogeneous arrays.

type	= c	ollec	tion(co	5110	ectio	n(reco	ord(a	a=intege	er, b=re	eal)))
values	= [	[(1,	1.1)]	[]	[(2,	2.2)	(3,	3.3)]]	[[(4,	4.4)]]
becomes										
data[][]@size	= 3	1		0	2				1 1	
data[][]-a data[][]-b	=	1	1.1		2	2.2	3	3.3	4	<b>↓</b> 4.4



```
pt = sqrt(px^{*2} + py^{*2}),
               phi = atan2(py, px),
               eta = \ln((\text{energy} + \text{pz})/(\text{energy} - \text{pz}))/2) # construct a record as output
           # make a bundle of plots
       .bundle(
                                                           # using the variables we've made
           mass = bin(120, 0, 12, "dimuon.mass"),
           pt = bin(100, 0, 100, "dimuon.pt"),
           eta = bin(100, -5, 5, "dimuon.eta"),
           phi = bin(314, 0, 2*pi, "dimuon.phi + pi"),
           muons = loop("goodmuons", "mu", bundle(
                                                          # also make plots with one muon per entry
               pt = bin(100, 0, 100, "mu.pt"),
               eta = bin(100, -5, 5, "mu.eta"),
               phi = bin(314, -pi, pi, "mu.phi")
           ))
pending = workflow.submit()
                                                          # submit the query
pending["mass"].plot()
                                                          # and plot results while they accumulate
pending["muons"]["pt"].plot()
                                                           # (they'll be animated)
blocking = pending.await()
                                                           # stop the code until the result is in
massplot = blocking.plot.root("mass")
                                                           # convert to a familiar format, like ROOT
massplot.Fit("gaus")
                                                           # and use that package's tools
```

#### No runtime errors:

Type system includes min/max ranges of numbers and collection lengths, so that out-of-domain errors are caught at compile-time. **Examples:** "x/y" must be "if y != 0: x/y else: None".

filter("goodmuons.size >= 2") above ensures
that we can assign two of them as mu1, mu2.

#### **Total functional language:**

Every expression that compiles returns a result. (For simplicity, recursion is not allowed.)

#0 #1	<pre>:= **(muons[]-p, 2) := +(#0, 0.011164) loop over muons[]@size</pre>
#2@size	<b>:=</b> \$explodesize(muons[], muons[]) over #2@size
#2	<pre>:= \$explodedata(#1, #2@size, (muons[]))</pre>
#3	<pre>:= \$explodedata(#1, #2@size, (muons[], muons[]))</pre>
#4	:= +(#2, #3)
#5	<pre>:= \$implode(#4, muons[], "max")</pre>
#6	<pre>:= \$implode(#5, None, "max")</pre>

Each non-\$ statement can be a branchless, loopless GPU kernel or all statements can be grouped by @size as CPU loops.



### Performance studies

- Typical analysis asks for ~10 TB of input data.
  Typical query touches < 1% of the *columns*.
- Disk read at 40 MB/sec  $\Rightarrow$  2.6 sec for 1000 disks.
- Typical analysis touches < 10% of the *columns*.
  Therefore, fill 1 TB of a cluster's RAM as cache.
- Cache read at 1 GB/sec  $\Rightarrow$  0.1 sec for 1000 cores.

#### Microbenchmarks to get a sense of scale:

Just adding a constant to an array of 64-bit floats on different systems. The first are disk limited, the last are memory bandwidth limited.

- 0.018 MHzCMSSW EDAnalyzer (C++ event framework)1.5 MHzTTree::Draw in ROOT (HEP analysis toolkit)2.8 MHzminimal disk read and unzip
  - 12 MHz allocating C++ objects on heap, iterating, deleting
  - 31 MHz allocating C++ objects on stack and iterating
- 54 MHz Femtocode loop, current implementation
- 250 MHz minimal single-threaded loop in C (our goal)
- 8 000 MHz same loop on 128 threads in KNL's MCDRAM
- 57 000 MHz equivalent on Tesla P100-SXM2 GPU

