UrQMD (Fortran) calls ROOT (C++)

Fan Si

Status

- Generally successful (but with some cases to be determined / developed)
 - Define C++ functions, and let them called by Fortran at proper locations
 - Define C++ extern struct to link Fortran quantities (commons)
- Location: ui:/ustcfs/HICUser/fsi/urqmd
- Original UrQMD: /ustcfs/HICUser/fsi/urqmd/urqmd
 - Change: pythia6409.f -> pytahi6428.f (final pythia6), GNUmakefile updated
 - Change: new gfortran option -std=legacy in mk/Linux.mk

New Fortran & C++ codes: /ustcfs/HICUser/fsi/urqmd/curqmd
c_*.f, curqmd.[cxx,h], GNUmakefile, run.sh

What to store (UrqmdDst)

colum	m#	contents
0		"E" (only in <u>016</u>)
1		# of collisions
2		# of elastic collisions
3		# of inelastic collisions
4		# of Pauli-blocked collisions
5	N	# of decays
6	N	# of produced <i>hard</i> baryon resonances
7		# of produced <i>soft</i> baryon resonances
8		# of baryon resonances produced via a decay of another resonance

- Entry for each time step of events (few)
- Additional branch

• b

- time: event level (instead of track level)
- Npart: Aproject + Atarget #(pptype%100==0)
- Ntracks
- spin
- iso3_old1 & iso3_old2

0 013	_01 4	_015	_O16	contents
0013	0011	1	0010	ind: index of particle (see CTOption (56))
1	1	2	1	<i>t</i> : time of particle
2	2	3	2 🗸	r_x : x coordinate
3	3	4	3	r_y : y coordinate
4	4	5	4	r_{z} : z coordinate
5	5	6	5	<i>E</i> : energy of particle
6	6	7	6 1	p_x : x momentum component
7	7	8		p_{u} : y momentum component
8	8	9	8 1	p_{z} : z momentum component
9	9	10	9 🗸	m: mass of particle
10	10	11		ityp: particle-ID
11	11	12		iso3: $2 \cdot I_3$ (see Section 1.2)
12	12	13	12	
13	13	14		parent collision number (see Table 10)
14	14	15	14	N _{coll} number of collisions
		16	V	S : strangeness
15	15		15 🗸	parent process type (see Table 11)
		17	V	history information (debugging only)
16			\checkmark	$t^{\rm fr}$: freeze-out time of particle
17				$r_x^{\rm fr}$: freeze-out x coordinate
18				r_{y}^{fr} : freeze-out y coordinate
19				$r_z^{\rm fr}$: freeze-out z coordinate
20			v	\tilde{E}^{fr} : freeze-out energy of particle
21			\checkmark	$p_x^{\rm fr}$: freeze-out momentum x component
22				p_{y}^{fr} : freeze-out momentum y component
23				$p_z^{\rm fr}$: freeze-out momentum z component
	16*			τ_{dec} decay time of particle
	17*			τ_{form} formation time of particle
	18*			R_{σ} cross section reduction factor
	19*			unique particle number (not ID!)
			16*	ityp ₁ ^{old} : particle-ID of parent particle # 1
			17*	$ityp_2^{old}$: particle-ID of parent particle # 2

What to store (UrqmdCollisionDst)

colu	mn#	format	contents
1		(i8)	number of ingoing particles $N_{\rm in}$
2		(i8)	number of outgoing particles $N_{\rm out}$
3	1	(i4)	process ID (see Table 11)
4	N	(i7)	collision/entry counter
5	N	(f8.3)	collision time t_{coll} in fm/c
6	N	(e12.4)	center of mass energy of the collision \sqrt{s} in GeV
7		(e12.4)	total cross-section of the collision σ_{tot} in mbarn
8		(e12.4)	partial cross-section of the actual sub-process σ_i in mbarn
9		(e12.4)	Baryon density at collision point $\rho_{\rm B}$ in units of ρ_0

• Not created or stored if storeCollision = false

- Entry for each collision of events (many)
- Additional branch

• b

- Ntracks
- spin
- iso3_old1 & iso3_old2

O ⁰¹³	O ⁰¹⁴	O ⁰¹⁵	_O1 6	contents
		1		ind: index of particle (see CTOption (56))
1	1	2	1	t: time of particle
2	2	3	2 🗸	r_x : x coordinate
3	3	4	3 🗸	r_y : y coordinate
4	4	5	4 🗸	r_z : z coordinate
5	5	6	5	E: energy of particle
6	6	7	6 🗸	p_x : x momentum component
7	7	8	7 🗸	p_y : y momentum component
8	8	9	8 🗸	p_z : z momentum component
9	9	10	9 🗸	m: mass of particle
10	10	11	10	ityp: particle-ID
11	11	12	11 🗸	iso3: $2 \cdot I_3$ (see Section 1.2)
12	12	13	12	ch : charge of particle
13	13	14	13	parent collision number (see Table 10)
14	14	15	14	N _{coll} number of collisions
		16		S: strangeness
15	15		15 🗸	parent process type (see Table 11)
		17		history information (debugging only)
16				$t^{\rm fr}$: freeze-out time of particle
17				$r_x^{\rm fr}$: freeze-out x coordinate
18				$r_y^{\rm fr}$: freeze-out y coordinate
19				$r_z^{\rm fr}$: freeze-out z coordinate
20				$E^{\rm fr}$: freeze-out energy of particle
21				p_x^{fr} : freeze-out momentum x component
22				p_y^{fr} : freeze-out momentum y component
23				$p_z^{\rm fr}$: freeze-out momentum z component
	16*			τ_{dec} decay time of particle
	17*			τ_{form} formation time of particle
	18*			R_{σ} cross section reduction factor
	19*		\checkmark	unique particle number (not ID!)
			16*	ityp11 : particle-ID of parent particle # 1
			17*	ityp21 : particle-ID of parent particle # 2

What is new

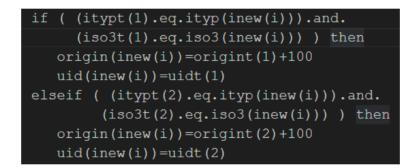
- Can store spin, iso3_old
- Can store parent information (iType_old, iso3_old) during final output
- Can store f13 information @ each time step (original f13 only stores final)
- Can set quiet output to command screen
- Can skip empty event (Ncoll + Ndecays == 0)
- Can set random seed event-by-event

How to store parent information

• Integer (4 bytes): origin (origin definition)

- = ppType + 100*(# scatters) + 1000*(|iType_old1| + 1000*|iType_old2|)
- 0 <= ppType < 100, 0 <= |iType_old| < 1000

• No iso3_old or sign of iType_old



ityp	nucleon	ityp	delta	ityp	lambda	ityp	sigma	ityp	xi	ityp	omega
1	N_{938}	17	Δ_{1232}	27	Λ_{1116}	40	Σ_{1192}	49	Ξ_{1317}	55	Ω_{1672}
2	N_{1440}	18	Δ_{1600}	28	Λ_{1405}	41	Σ_{1385}	50	Ξ_{1530}		
3	N_{1520}	19	Δ_{1620}	29	Λ_{1520}	42	Σ_{1660}	51	Ξ_{1690}		
4	N_{1535}	20	Δ_{1700}	30	Λ_{1600}	43	Σ_{1670}	52	Ξ_{1820}		
5	N_{1650}	21	Δ_{1900}	31	Λ_{1670}	44	Σ_{1775}	53	Ξ_{1950}		
6	N_{1675}	22	Δ_{1905}	32	Λ_{1690}	45	Σ_{1790}	54	Ξ_{2025}		
7	N_{1680}	23	Δ_{1910}	33	Λ_{1800}	46	Σ_{1915}				
8	N_{1700}	24	Δ_{1920}	34	Λ_{1810}	47	Σ_{1940}				
9	N_{1710}	25	Δ_{1930}	35	Λ_{1820}	48	Σ_{2030}				
10	N_{1720}	26	Δ_{1950}	36	Λ_{1830}						
11	N_{1900}			37	Λ_{1890}						
12	N_{1990}			38	Λ_{2100}						
13	N_{2080}			39	Λ_{2110}						
14	N_{2190}										
15	N_{2200}										
16	N_{2250}										

Table 1: Baryon-itypes used in UrQMD. Antibaryons carry a negative sign.

ityp	0^{-+}	ityp	1	ityp	0++	ityp	1++	ityp	charmed
101	π	104	ρ	111	a_0	114	a_1	133	D
106	K	108	K^*	110	K_0^*	113	K_1^*	134	D^*
102	η	103	ω	105	f_0	115	f_1	135	J/Ψ
107	η'	109	ϕ	112	f_0^*	116	f'_1	136	χ_c
ityp	1+-	ityp	2^{++}	ityp	(1)*	ityp	(1)**	137	Ψ'
122	b_1	118	a_2	126	ρ_{1450}	130	ρ_{1700}	138	D_s
121	K_1	117	K_2^*	125	K_{1410}^{*}	129	K_{1680}^{*}	139	D_s^*
123	h_1	119	f_2	127	ω_{1420}	131	ω_{1662}		
124	h'_1	120	f_2'	128	ϕ_{1680}	132	ϕ_{1900}		

ppType = origin%100
scatters = origin/100%10

	2 1 10	318 3.486 0	.1665E+01 0.198	3E+02 0.1983E+0	2 0.5118E+00								
714	0.34855015E+01	-0.31837811E+01	0.30769649E+01	-0.22947721E+01	0.62539188E+00	-0.15451485E+00	0.31808005E+00	-0.49701232E+00	0.13800000E+00	101 -2 -1	277	1 0	1001015
233	0.34855015E+01	-0.26863519E+01	0.33516045E+01	-0.23172907E+01	0.44498825E+01	0.47432728E+00	-0.64771840E+00	-0.42751707E+01	0.93800002E+00		267	4 0	1017030
233	0.34855015E+01	-0.29350665E+01	0.32142847E+01	-0.23060314E+01	0.50752743E+01	0.31981242E+00	-0.32963835E+00	-0.47721830E+01	0.16654542E+01	4 -1 0	318	50	1101010

How to store parent information

- Integer (4 bytes): origin (new definition)
 - $\circ = ppType + 100*(\# scatters) + 1000*((iso3_old1+3)+7*(t_iType_old1+100))$
 - $+ 140000*((iso3_old2+3)+7*(t_iType_old2+100))$
 - 0 <= ppType < 100, 0 <= iso3_old+3 <= 6
 - t_iType_old = iType_old; if(>=100) -=40; if(<=-100) +=40
 - $0 \le t_iType_old+100 \le 200$

• origin < 1.96e9 (4-byte integer max ~ 2.1e9)

- !!!Also affect *.txt output
- ppType = origin%100
- # scatters = origin/100%10
- If ppType does not contain # scatters, can be stored 1-byte

ppType = origin%1000;
if(origin<1000)
<pre>iType_old1 = iso3_old1 = iType_old2 = iso3_old2 =</pre>
else{
iso3_old1 = origin/1000%7-3;
iType_old1 = origin/7000%200-100;
if(iType_old1>=60)
iType_old1 += 40;
if(iType old1<=-60)
iType old1 -= 40;
if(origin<1400000)
iType old2 = iso3 old2 = 0;
else{
iso3 old2 = origin/1400000%7-3;
iType old2 = origin/9800000-100;
if (iType old2>=60)
iType old2 $+=$ 40;
if (iType old2<=-60)
iType old2 -= 40;
}

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6	N_{1675}	22	Δ_{1905}	32	Λ_{1690}	45	Σ_{1790}	54	Ξ_{2025}		
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8	N_{1700}	24	Δ_{1920}	34	Λ_{1810}	47	Σ_{1940}				
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11	N_{1900}			37	Λ_{1890}						
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		-						_	
ityp	0^{-+}	ityp	$1^{}$	ityp	0++	ityp	1++	ityp	charmed
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102	η	103	ω	105	f_0	115	f_1	135	J/Ψ
107	η'	109	ϕ	112	f_0^*	116	f'_1	136	χ_c
ityp	1+-	ityp	2^{++}	ityp	(1)*	ityp	(1)**	137	Ψ'
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124	h'_1	120	f'_2	128	ϕ_{1680}	132	ϕ_{1900}		

How to set seed for random generator

- Random seed generator
 iseed is actual seed when used
 - Changed after each use
 - ranseed is seed of iseed
 - 1. set ranseed < 0
 - 2. if < 0, set ranseed by time
 - if different from old
 - 3. set iseed = -ranseed
 - 4. modify iseed (extra if < 0)
 - initialize random number generator
 - call auto-seed generator only for first event and if no seed was fixed if(.not.firstseed.and.(.not.fixedseed)) then
 - ranseed=-(1*abs(ranseed)) call sseed(ranseed)

else

firstseed=.false.

subroutine sseed(ranseed)

reset the random number generato

implicit none

real*8 dummy, ran2 integer iseed, ranseed, oldseed, time, timeseed common /seed/iseed,oldseed

```
if (ranseed.le.0) then
    timeseed = abs(time())
    if (timeseed.eq.oldseed) return
    ranseed = timeseed
endif
oldseed = ranseed
iseed = -ranseed
dummy = ran2(iseed)
```

return end

function ran2(idum)

Long period (>2E18) random number generator of L'Ecuyer with Bays-Durham shuffle and added safeguards. Returns a uniform random deviate between 0.0 and 1.0 (exclusive of the endpoint values). Call with idum a negative integer to initialize; thereafter, do not alter idum between successive deviates in a sequence. RNMX should approximate the largest floating value that is less than 1.

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implicit none

integer idum,IM1,IM2,IMM1,IA1,IA2,IQ1,IQ2,IR1,IR2,NTAB,NDIV
real*8 ran2,AM,EPS,RNMX

parameter (IM1=2147483563,IM2=2147483399,AM=1.0D0/IM1,IMM1=IM1-1
\$ IA1=40014,IA2=40692,IQ1=53668,IQ2=52774,IR1=12211,IR2=3791,
\$ NTAB=32,NDIV=1+IMM1/NTAB,EPS=1.0D0-16,RNMX=1.0D0-EPS)
integer idum2, j, k, iv(NTAB), iy
save iv, iy, idum2
data idum2/123456789/, iv/NTAB*0/, iy/0/

```
if (idum.le.0) then
    idum=max(-idum,1)
    idum2=idum
    do 11 j=NTAB+8,1,-1
        k=idum/IQ1
        idum=IA1*(idum-k*IQ1)-k*IR1
        if (idum.lt.0) idum=idum+IM1
        if (j.le.NTAB) iv(j)=idum
        continue
        iy=iv(1)
```

iy=iv(i) endif k=idum/IQ1 idum=IA1*(idum-k*IQ1)-k*IR1 if (idum.lt.0) idum=idum+IM1 k=idum2/IQ2 idum2=IA2*(idum2-k*IQ2)-k*IR2 if (idum2.lt.0) idum2=idum2+IM2 j=1+iy/NDIV iy=iv(j)-idum2 iv(j)=idum if (iy.lt.1) iy=iy+IMM1 ran2=min(AM*iy,RNMX)

return end

8

endif

How to set seed for random generator

Set seed for gRandom by random_device
True random number if supported by compiler
Set c_use_external_seed = true

- Set ranseed uniformly in [-2147483648, 2147483647]
 - Not finalized yet
 - Set ranseed only positive?
 - Add effect of time?
- Call ran2 to modify iseed

```
std::random_device rd;
gRandom->SetSeed(rd());
```

```
c_sets_l_.c_use_external_seed = kTRUE;
```

```
void c_set_seed_() {
```

```
sys_.ranseed = (gRandom->Rndm()-0.5)*4294967296;
seed_.oldseed = sys_.ranseed;
seed_.iseed = -sys_.ranseed;
ran2 (&seed .iseed);
```

```
initialize random number generator
call auto-seed generator only for first event and if no seed was fixed
if (c use external seed) then
 if(firstseed)firstseed=.false.
 if (fixedseed) then
   write (6, *) 'c use external seed is neglected due to fixedseed'
   c_use_external seed=.false.
 else
    call c set seed
 endif
elseif(.not.firstseed.and.(.not.fixedseed)) then
   ranseed=-(1*abs(ranseed))
  call sseed(ranseed)
else
   firstseed=.false.
endif
```

Other developments

- Parent particle Information in collision output
 - In temporary quantities named t*
 - No spin, or Nin, rho_b values
 - Saved to my defined global quantity
 - \circ Extraction method is the same as other t*
- Now time step tree filled after final particle decays rather than after final time step
 - Unstable particles' decays after all time steps & before the end of event
 - The same as origin output f14
- If CTOption(4)==1, initial information (t==0) is stored in tree
 - The same as origin output f14

Other developments

- Pdgid (int) -> mass (float), iType (short), iso3 (char), charge (char)
 - Strange masses of protons
 - Different masses between UrQMD and PDG
 - Some particles in UrQMD has no pdgid in ityp2pdg.f
 - Such as iType = 12, 13, 15, 16

c Neutron		
•	1, -1, 2112,	
c Proton		
	1, 1, 2212,	
c N*		
	2, -1, 12112,	2, 1, 12212,
	3, -1, 1214,	3, 1, 2124,
	4, -1, 22112,	4, 1, 22212,
	5, -1, 32112,	5, 1, 32212,
	6, -1, 2116,	6, 1, 2216,
	7, -1, 12116,	7, 1, 12216,
	8, -1, 21214,	8, 1, 22124,
	9, -1, 42112,	9, 1, 42212,
	10, -1, 31214,	10, 1, 32124,
	14, -1, 1218,	14, 1, 2128,
c Delta		
•	17, -3, 1114,	17, -1, 2114, 17, 1, 2214, 17, 3, 2224,
	18, -3, 31114,	18, -1, 32114, 18, 1, 32214, 18, 3, 32224,
	19, -3, 1112,	19, -1, 1212, 19, 1, 2122, 19, 3, 2222,
	20, -3, 11114,	20, -1, 12114, 20, 1, 12214, 20, 3, 12224,
	21, -3, 11112,	21, -1, 11212, 21, 1, 12122, 21, 3, 12222,
	22, -3, 1116,	22, -1, 1216, 22, 1, 2126, 22, 3, 2226,
•	23, -3, 21112,	23, -1, 21212, 23, 1, 22122, 23, 3, 22222,
	24, -3, 21114,	24, -1, 22114, 24, 1, 22214, 24, 3, 22224,
	25, -3, 11116,	25, -1, 11216, 25, 1, 12126, 25, 3, 12226,
•	26, -3, 1118,	26, -1, 2118, 26, 1, 2218, 26, 3, 2228,

data	baryon_names/
•	'Nukleon',
•	'N(1440)',
•	'N(1520)',
•	'N(1535)',
•	'N(1650)',
	'N(1675)',
	'N(1680)',
•	'N(1700)',
•	'N(1710)',
•	'N(1720)',
•	'N(1900)',
•	'N(1990)',
•	'N(2080)',
	'N(2190)',
•	'N(2220)',
•	'N(2250)',

ityp	nucleon	ityp	delta	ityp	lambda	ityp	sigma	ityp	xi	ityp	omega
1	N_{938}	17	Δ_{1232}	27	Λ_{1116}	40	Σ_{1192}	49	Ξ_{1317}	55	Ω_{1672}
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3	N_{1520}	19	Δ_{1620}	29	Λ_{1520}	42	Σ_{1660}	51	Ξ_{1690}		
4	N_{1535}	20	Δ_{1700}	30	Λ_{1600}	43	Σ_{1670}	52	Ξ_{1820}		
5	N_{1650}	21	Δ_{1900}	31	Λ_{1670}	44	Σ_{1775}	53	Ξ_{1950}		
6	N_{1675}	22	Δ_{1905}	32	Λ_{1690}	45	Σ_{1790}	54	Ξ_{2025}		
7	N_{1680}	23	Δ_{1910}	33	Λ_{1800}	46	Σ_{1915}				
8	N_{1700}	24	Δ_{1920}	34	Λ_{1810}	47	Σ_{1940}				
9	N1710	25	Δ_{1930}	35	Λ_{1820}	48	Σ_{2030}				
10	N_{1720}	26	Δ_{1950}	36	Λ_{1830}						
11	N_{1900}			37	Λ_{1890}						
12	N_{1990}			38	Λ_{2100}						
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16	N_{2250}										

ityp	0^{-+}	ityp	1	ityp	0^{++}	ityp	1++	ityp	charmed
101	π	104	ρ	111	<i>a</i> 0	114	a_1	133	D
106	K	108	K^*	110	K_0^*	113	K_1^*	134	D^*
102	η	103	ω	105	f_0	115	f_1	135	J/Ψ
107	η'	109	ϕ	112	f_0^*	116	f'_1	136	χ_c
ityp	1+-	ityp	2^{++}	ityp	(1)*	ityp	(1)**	137	Ψ'
122	b_1	118	a_2	126	ρ_{1450}	130	ρ_{1700}	138	D_s
121	K_1	117	K_2^*	125	K_{1410}^{*}	129	K_{1680}^{*}	139	D_s^*
	h_1	119	f_2	127	ω_{1420}	131	ω_{1662}		
123	n_1		12		1420		1002		

Wrong name?

What to be determined

- Npart: Aproject + Atarget #(pptype%100==0)
 - ppType = origin%100
 - # scatters = origin/100%10

• Some particles only explore scatters, contribute to Npart?

	2 2 17	233 1.554 0	.5056E+01 0.3978	8E+02 0.1166E+02	2 0.8930E+00								
362	0.15541405E+01 -	-0.22213285E+01	0.35829759E+01 -	-0.71742345E+00	0.17039231E+01	-0.17381395E+00	0.13815719E+01 -0.29078794E+00	0.93800002E+00	1 -1	0 2	13	10	100
233	0.15541405E+01 -	-0.27680167E+01	0.33953480E+01 -	-0.60064611E+00	0.85553163E+01	0.20642659E+00	0.13534976E+00 -0.85059975E+01	0.88346555E+00	1 1		0	0 0	0
362	0.15541405E+01 -	-0.22213285E+01	0.35829759E+01 ·	-0.71742345E+00	0.84217759E+01	0.32168377E+00	0.25335456E-01 -0.83631539E+01	0.93800002E+00	1 -1	0 2	33	20	200
233	0.15541405E+01 -	-0.27680167E+01	0.33953480E+01 -	-0.60064611E+00	0.18374634E+01	-0.28907113E+00	0.14915862E+01 -0.43363154E+00	0.93800002E+00	1 1	1 2	33	1 0	100

Check	=====> EVENT:0 b	= 13.9877
73 0.74690319E-01 0.16514065E+01 0.84431142E-01 0.11221819E-01 0.10077745E+03 0.21420381E+01 -0.72458558E+00 -0.10074771E+03 0.93800002E+00 1 1 1 1 1 1 0 100	рх	= 1 = 0.0746903 = 199.043 = 52.0758 = 10.3074 = 14.489 = 17 = 2 = 2 = 4 = 73, 352, 73, 352 = 1, 1, 1, 1 = 1, -1, 1, -1 = 1, 0, 1, 0 = 1, -1, -1, 1 = 1, 65141, 0.385643, 1.65141, 0.385643 = 0.0844311, 0.0137725, 0.0844311, 0.0137725 = 0.0112218, 0.0112218, 0.0112218 = 0.014267, 0.0926628, 2.14204, -2.03525 = 0.148645, -0.105794, -0.724586, 0.767437 = 98.2774, -100.773, -100.748, 98.2526 = 0.917558, 0.919356, 0.938, 0.938 = 0, 0, 1, 1 = 0, 0, 100, 100 = 0, 0, 0, 0 = 0, 0, 0, 0

Check		root [3] UrgmdCollisionDst->Show(3); =====> EVENT:3
2 3 28 4 0.085 0.6134E+02 0.6747E+01 0.5664E+01 0.9465E+01 0.28688709E-01 -0.10823742E+01 -0.4772494E+00 6.2210789E-01 0.5237234E+02 -0.22741829E+00 0.10696294E+00 -0.12958522E+02 0.12080407E+ 200 0.38888709E-01 -0.0803534E+00 -0.70689933E+00 0.22102056E-01 0.10598629E+02 -0.3930945E+00 -0.120921140E+00 0.100336E+02 0.2122020E+ 200 0.38888709E-01 -0.082934E+00 -0.70689933E+00 0.2192056E-01 0.7373737E+02 -0.21838169E+00 0.100350E+00 0.1295532E+02 0.120804002E+ 200 0.38888709E-01 -0.8903534E+00 -0.70689933E+00 0.2192056E-01 0.10732337E+02 -0.11539479E+00 0.19100350E+00 0.10658570E+01 0.38864038E+ 200 0.38880709E-01 -0.8903534E+00 -0.70689933E+00 0.2192056E+01 0.10732337E+02 -0.11539479E+00 0.19100350E+00 0.10658570E+02 0.08564038E+ 200 0.38880709E-01 -0.8903534E+00 -0.70689933E+00 0.2192056E+01 0.10732337E+02 -0.11539479E+00 0.19100350E+00 0.10658570E+02 0.08564038E+	1 113 -1 0 2 1 -1 99331 1 17 -1 0 4 2 0 169902 0 101 -2 -1 4 2 0 169902	$\begin{array}{rrrrrrrrrrrrrrrrrrrrrrrrrrrrrrrrrrrr$

Check

1 2 20 11 3.864 0.1321E+01 0.0000E+00 0.0000E+00 0.9497E-02

462 0.38641318E+01 -0.44654879E+00 -0.12298002E+01 -0.27913244E+01 0.29055463E+00 0.16513024E+00 -0.16992283E+00 -0.96105291E-01 0.13800000E+00

	root [4] UrqmdCo1 =====> EVENT:10	lisionDst->Show(10);
	b	= 13.9877
	number	= 11
	time	= 3.86413
	sqrts	= 1.32149
	sigma_total	= 0
	sigma_partial	= 0
	rho_B	= 0.00949667
	pType Nam	= 20
	Nin Nout	= 1 = 2
	Ntracks	= 3
	uid	= 447,
		462. 463
	iType	= 118,
		101. 104
	iso3	= 2,
	abamga	2, 0
	charge	2, 0 = 1, 1, 0
	spin	= -2,
	-*	0, 0
	х	= -0. 446549, -0. 446549, -0. 446549
		-0.446549, -0.446549
028 020	У	= -1.2298, -1.2298, -1.2298
020 020	-	-1.2298, -1.2298 = -2.79132, -2.79132, -2.79132 = 0.242734
020	Z	2.19132, -2.79132 -2.79132
	рх	= 0. 242734,
	¥	0.16513, 0.0776042
	ру	= -0.285845,
		-0.169923, -0.115922
	pz	= -1.54135,
	m	-0.0961053, -1.44524 = 1.32149,
	111	
	Nc	= 1,
		0.138, 1.0194 = 1, _2, 1
	pcNumber	= h.
	Т	11, 11
	ррТуре	= 28, 20, 20
	iType_old1	= 101,
	11, 10, 10, 10, 10, 10, 10, 10, 10, 10,	118, 118
	iso3_o1d1	= 0, 2, 2
		2, 2
	iType_old2	= 1,
	iso3_o1d2	0, 0 = -1,
	1803_0142	1, 0, 0
		0, 0

1	
-	\sim

- (b_{max})_{max} = nucrad(Ap)+nucrad(At)+2*CTParam(30)
 CTP(30): radius offset for initialization (1.5 default)
 - 197 Au+ 197 Au: 15.82 = 6.41+6.41+1.5*2

```
function nucrad(AA)
     implicit none
     real*8 nucrad, r 0
     integer A,AA
     include 'coms.f'
     include 'options.f'
     A=abs(AA)/CTParam(67)
c root mean square radius of nucleus of mass A
c r 0 corresponding to rho0
     if (CTOption(24).eq.1) then
c root mean square radius of nucleus of mass A (Mayer-Kuckuck)
        nucrad = 1.128 + a^{*}(1./3.) - 0.89 + a^{*}(-(1./3.))
     else
        r 0 = (0.75/pi/rho0) ** (1./3.)
c subtract gaussian tails, for distributing centroids correctly
        nucrad = r 0*(0.5*(a + (a**(1./3.)-1.)**3.))**(1./3.)
     endif
      return
     end
```

• pro: 197 79

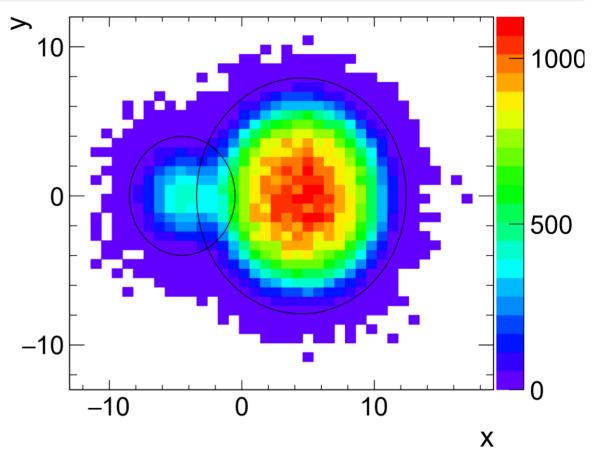
• tar: 16 8

 $\circ \sqrt{s_{\rm NN}} = 200 {\rm ~GeV}$

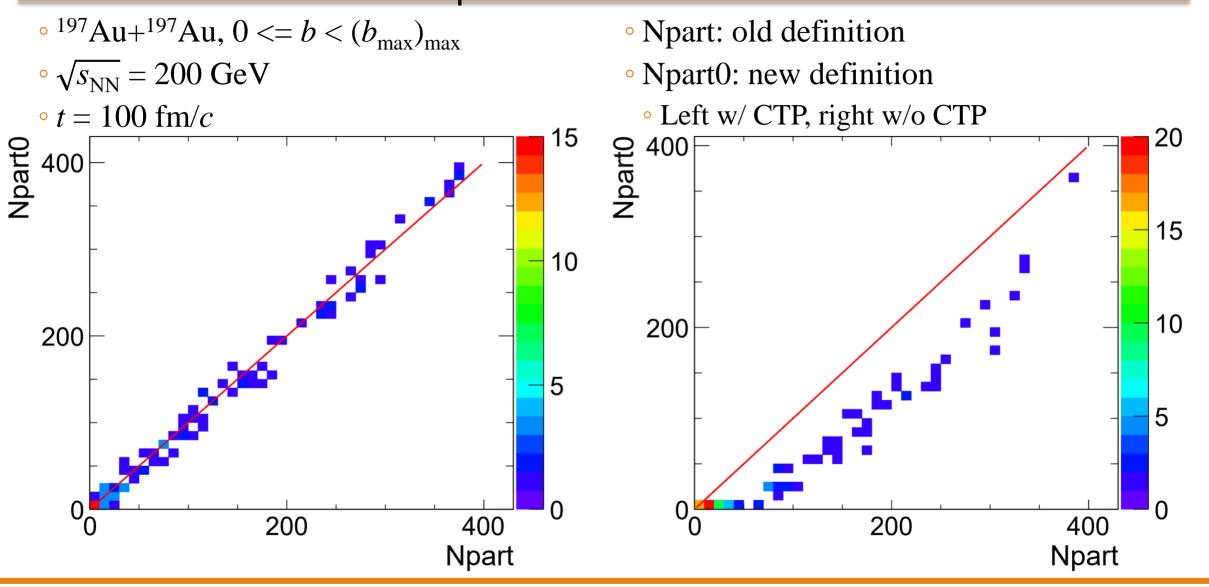
• *b* = 9 fm

• t = 0

- Center: pro (b/2, 0), tar (-b/2, 0)
- Radius: nucrad(A)+CTParam(30)
 - But CTP(30) only contributes to $(b_{\max})_{\max}$ rather than nuclei initialization

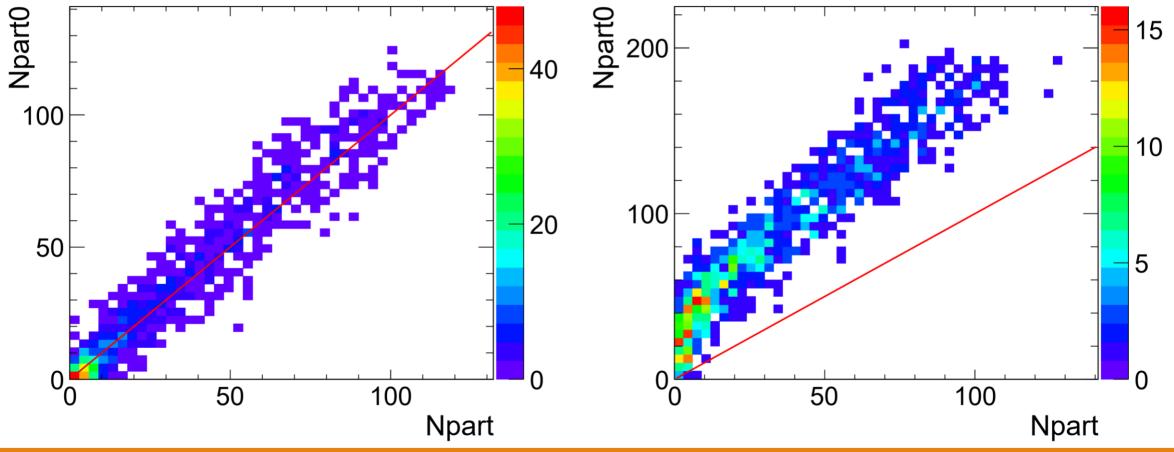


• $N_{\text{part}} = #((x-b/2)^2 + y^2 < R(\text{pro}) \&\& (x+b/2) + y^2 < R(\text{tar}))$ at t = 0



- ¹⁹⁷Au+¹⁶O, $0 \le b \le (b_{\text{max}})_{\text{max}}$
- $\circ \sqrt{s_{\rm NN}} = 200 {
 m GeV}$
- t = 100 fm/c

Left: CTP(30) = 1.5 (default)
Right: CTP(30) = 4



subroutine boostnuc(i1,i2,pin,b,dst)

implicit none

continue return end

```
include 'coms.f'
                                                                                      include 'options.f'
    dstp=dstp+CTParam(30)
    dstt=dstt+CTParam(30)
                                                                                      integer i1,i2,i
       if(bdist.gt, (nucrad(Ap)+nucrad(At)+2*CTParam(30)))
                                                                                      real*8 b,<mark>dst</mark>,ei,ti
            bdist=nucrad(Ap)+nucrad(At)+2*CTParam(30)
&
                                                                                      real*8 pin,beta,gamma
       if(bdist.gt.(nucrad(Ap)+nucrad(At)+2*CTParam(30)))
            bdist=nucrad(Ap)+nucrad(At)+2*CTParam(30)
&
                                                                                      do 1 i=i1,i2
 CTParam(30)=1.5
                                                                                      beta = pin/sqrt(pin**2+fmass(i)**2)
real*8 nucrad, dstt, dstp, pcm, eb, embeam, emtarget
                                                                                      gamma = 1.d0/sgrt(1.d0-beta**2)
       dstp = nucrad(Ap)+CTParam(41)
       dstp = nucrad(Ap)*ratio**(2.0/3.0)+CTParam(41)
                                                                                   Gallilei-Trafo in x-direction (impact parameter)
    dstp=dstp+CTParam(30)
     dst=0.5d0*(dstt+dstp)
                                                                                  projectile hits at POSITIVE x
       dstp=0d0
                                                                                         rx(i) = rx(i) + b
                      pboost, 0.5*bimp, -dstp)
                                                                                  distance between nuclei: projectile at NEGATIVE z for dst < 0
&
               pboost, 0.5*bimp, -dstp+CTParam(20))
&
                                                                                         if(CTOption(23).eq.0)then
              call boostnuc(npart, npart, -pbeam, 0.5*bimp, -dstp)
                                                                                           ti = r0(i)
              call boostnuc(npart, npart, 0. d0, 0. 5*bimp, -dstp)
                                                                                           rz(i) = rz(i)/gamma+dst/gamma
              call boostnuc(npart, npart, -ppeg, 0.5*bimp, -dstp)
                                                                                         else
real*8 nucrad, dstt, dstp, pcm, eb, embeam, emtarget
                                                                                          rz(i) = (rz(i) + dst)
                                                                                         end if
                                                                                         Ei = p0(i)
                                                                                         p0(i) = gamma*(p0(i) - beta*pz(i))
                                                                                         pz(i) = gamma*(pz(i) - beta*Ei)
```

24	1	initialization mode
	0	hard sphere (used for EOS \neq 0)
	1	Woods-Saxon (used for CASCADE mode)
	2	Fast Woods-Saxon (used for CASCADE mode)

if (CTOption(24).eq.1) then
R2 = nucrad(A) + 10.0
else
R2 = nucrad(A)
endif

subroutine nucfast(IA,JJ)

include 'inputs.f'
real*8 nucrad

am=0.545 rad=nucrad(ia)/am

- UrqmdDst
 - TClonesArray*[4] containing the following classes
 - initOutputTree: returns TTree for writing
 - initInputChain: reads input file (*.lis
 - chain->GetEntry(i);
 - urqmdDst-> ...

- UrqmdEvent
- UrqmdTimestep
- UrqmdCollision
- UrqmdTrack

	6
st. *	root), and returns TChain for analyzing
	s UrqmdDst{
pu	blic:
	enum {Event = 0, Timestep, Collision, Track, NUrqmdArrays};
	<pre>static const TString UrqmdArrayNames[NUrqmdArrays];</pre>
	<pre>static const TString UrqmdArrayTypes[NUrqmdArrays];</pre>
	static const UInt_t UrqmdArraySizes[NUrqmdArrays];
	UrgmdDst();
	virtual ~UrqmdDst();
	UrqmdEvent* event(){return (UrqmdEvent*)mUrqmdArrays[Event]->UncheckedAt(0);}
	<pre>UrqmdTimestep* timestep(UInt_t i){return (UrqmdTimestep*)mUrqmdArrays[Timestep]->UncheckedAt(i);}</pre>
	<pre>UrqmdCollision* collision(UInt_t i){return (UrqmdCollision*)mUrqmdArrays[Collision]->UncheckedAt(i);}</pre>
	UrqmdTrack* track(UInt_t i){return (UrqmdTrack*)mUrqmdArrays[Track]->UncheckedAt(i);}
	UInt_t numberOfTimesteps(){return mUrqmdArrays[Timestep]->GetEntriesFast();}
	<pre>UInt_t numberOfCollisions() {return mUrqmdArrays[Collision]->GetEntriesFast();}</pre>
	<pre>UInt_t numberOfTracks() {return mUrqmdArrays[Track]->GetEntriesFast();}</pre>
	TTree* initOutputTree(const TString outputTreeName = "urgmd");
	TChain* initInputChain(const TString inputFileName, const TString inputTreeName = "urgmd");
pr	otected:
	TClonesArray* mUrqmdArrays[NUrqmdArrays];
ļ	
1 C	efROOT
1	ClassDef(UrqmdDst, 1)
#end	if
};	

• UrqmdEvent

- id: event id
 - $\circ 0 \sim 4294967295$
- b: impact parameter
 0 ~ 26.214

class UrqmdEvent : public TObject{

enum {BScale = 2500};

UrqmdEvent() : mId(0), mB(0){}
UrqmdEvent(const UInt_t id, const Double_t b) : mId(id), mB(TMath::Nint(b*Double_t(BScale)))
virtual ~UrqmdEvent(){}

UInt_t id() const{return mId;}
Double_t b() const{return mB/Double_t(BScale);}

void print() const; void set(const UInt_t id, const Double_t b);

protected: UInt_t mId; UShort_t mB;

```
#ifdef __ROOT___
ClassDef(UrqmdEvent, 1)
#endif
```

- UrqmdTimestep
 - time: Float16_t (24 bit, truncated mantissa)
 - nPart: old definition
 - nColl = Elastic + Inelastic + Bloacked
 - nElasticColl
 - nInelasticColl
 - nBloackedColl
 - nDecays
 - nHardRes
 - nSoftRes
 - nDecayRes
 - nTracks

• startTrack: array index of 1st track at this timestep

```
class UrgmdTimestep : public TObject{
  UrgmdTimestep() : mTime(0.), mNPart(0), mNElasticColl(0), mNInelasticColl(0),
   UrgmdTimestep(const Double t time, const UInt t nPart, const UInt t nElasticCo
   virtual ~UrgmdTimestep(){}
   Double t time() const{return mTime;}
   UInt t nPart() const{return mNPart;}
   UInt t nColl() const{return nElasticColl()+nInelasticColl()+nBloackedColl();
   UInt t nElasticColl() const{return mNElasticColl;}
  UInt t nInelasticColl() const{return mNInelasticColl;}
   UInt t nBloackedColl() const{return mNBloackedColl;}
   UInt t nDecays() const{return mNDecays;}
   UInt t nHardRes() const{return mNHardRes;}
  UInt t nSoftRes() const{return mNSoftRes;}
   UInt t nDecayRes() const{return mNDecayRes;}
  UInt t nTracks() const{return mNTracks;}
   void print() const;
   void set(const Double t time, const UInt t nPart, const UInt t nElasticColl,
  Float16 t mTime;
  UShort t mNPart;
   UShort t mNElasticColl;
   UShort t mNInelasticColl;
   UShort t mNBloackedColl;
   UShort t mNHardRes;
   UShort t mNSoftRes;
   UShort t mNDecayRes;
  UInt t mStartTrack;
```

ifdef __ROOT

ClassDef(UrqmdTimestep, 1) eendif 1.

• UrqmdCollision: -	~ 20000 for 200 GeV, $b = 0$, $t = 50$ fm/c	<pre>class UrqmdCollision : public TObject{ public: UrqmdCollision() : mTime(0.), mSqrts(0.), mSigmaTotal</pre>
• time		<pre>UrqmdCollision(const Double_t time, const Double_t sq virtual ~UrqmdCollision() {}</pre>
• Sqrts		<pre>Double_t time() const{return mTime;} Double_t sqrts() const{return mSqrts;}</pre>
 sigmaTotal 		<pre>Double_t sigmaTotal() const{return mSigmaTotal;} Double_t sigmaPartial() const{return mSigmaPartial;} Double t rhoB() const{return mRhoB;}</pre>
 sigmaPartial 	1.1 wall interaction (infinite matter calculations)	<pre>UInt_t type() const{return mType;} UInt_t nIn() const{return mNIn;}</pre>
• rhoB: $-1 \sim 30+$	1, $N > 1$ decay 2,1 annihilation	<pre>UInt_t nOut() const{return mNOut;} UInt_t nTracks() const{return nIn()+nOut();} UInt t startTrack() const{return mStartTrack;}</pre>
• Type: < 100	 2,2 scattering 2,N > 2 string excitation and decay 1,0 Pauli-blocked decay 	<pre>void print() const;</pre>
• nIn: < 2?	2,0 Pauli-blocked scattering<i>N</i>,0 Fluidization: All particles at the beginning of the fluid-dynamical part (Process-ID "91"). Please	<pre>void set(const Double_t time, const Double_t sqrts, c protected:</pre>
• nOut	Note: this also includes spectators and corona-particles that are not transferred to the fluid! 0 ; N Particlization: All particles at the end of the fluid-dynamical part (Process-ID "96"). Here, t_{coll}	// UInt_t mId;
• $nTracks = nIn + n$		<pre>Float16_t mSqrts; Float16_t mSigmaTotal; Float16 t mSigmaPartial;</pre>
 startTrack 		Float16_t mRhoB; UChar_t mType;
		UChar_t mNIn; UChar_t mNOut; UInt_t mStartTrack;
• An empty collisio		#ifdefROOT ClassDef(UrqmdCollision, 1) #endif

};

• UrqmdTrack: ~86000 for 200 GeV, b = 0, t = 50 fm/c (collision+1timestep) • mType = (shifted_itype+100)+200((iso3+3)+7((charge+2)+5(spin+3))) enum {RSigmaScale = 250}; UInt t mId; • $|itype| \ge 140$, shifted_itype = 0 (PdgId+1000 used for some c/b hadrons) UChar t mTimestep; UShort t mCollision; • $|itype| \ge 100$, shifted_itype = itype-40*|itype|/itype UShort t mType; JInt t id() const{return mId;} • iso $3 = 2I_3, -3 \sim 3$ JInt t timestep() const{return mTimestep;} JInt t collision() const{return mCollision;} • charge, $-2 \sim 2$ Float16 t mX; [nt t iso3() const{return mType/200%7-3;} Int t charge() const{return mType/1400%5-2;} • spin = $2s_3?, -3 \sim 3$ Float16 t mZ; Int t spin() const{return mType/7000-3;} Double t x() const{return mX;} Float16 t mPx; Double t y() const{return mY;} Float16 t mPy; \circ m: 0 - 6.5535 Double t z() const{return mZ;} Float16 t mPz; UShort t mM; Double t py() const{return mPy;} UChar t mNCollisions; • mOrigin)ouble t pz() const{return mPz;} UShort t mParentCollision; Double t m() const{return mM/Double t(MScale);} JInt t nCollisions() const{return mNCollisions;} parentCollisionType, itypeOld1, JInt t parentCollision() const{return mParentCollision;} JInt t parentCollisionType() const{return mOrigin%1000;} iso3Old1, itypeOld2, iso3Old2 Int t itypeOld1() const; Int t iso30ld1() const{if(mOrigin<1000) return 0; return mOrigin/200000%7-3;}</pre> Int t itypeOld2() const; • rSigma: 0 − 1 Int t iso30ld2() const{if(mOrigin<1400000) return 0; return mOrigin/280000000-3;}</pre> Float16 t mXFreezeOut; Double t xFreezeOut() const{return mXFreezeOut;} Float16 t mYFreezeOut; Double t yFreezeOut() const{return mYFreezeOut;} • If mTimestep=mCollision=0 Float16 t mZFreezeOut; Double t zFreezeOut() const{return mZFreezeOut;} Float16 t mPxFreezeOut; Double t pxFreezeOut() const{return mPxFreezeOut;} Float16 t mPyFreezeOut; Double t pyFreezeOut() const{return mPyFreezeOut;} 1 0 Float16 t mPzFreezeOut; ouble t pzFreezeOut() const{return mPzFreezeOut;} Double t tauDecay() const{return mTauDecay;} Float16 t mTauDecay; Double t tauForm() const{return mTauForm;} Float16 t mTauForm;

Double t rSigma() const{return mRSigma/Double t(RSigmaScale);}

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UChar t mRSigma;

Other updates

- Random seed (Int_t) =
 - $(Int_t((gRandom->Rndm()-0.5)*0x10000000))^(Int_t(ULong64_t(gSystem->Now()))%0x10000)$
 - gRandom->Rndm(): $-1 \sim 1$, ULong64_t(gSystem->Now()): current time in milliseconds
- All input settings summarized together
 - New input parameter: event ID offset
- Extended track information extracted for collision history
- GNUmakefile updated, rootcint compiler called
 - ROOT IO functions required by UrqmdDst
- New Npart definition in UrqmdEvent?
- Define 1 or 2 switches for some infrequently used track information?
 - Split UrqmdTrack to 2 or 3 classes

Result

- 10 events, 200 GeV, b = 0, t = 50 fm/c (collision+1timestep)
 32 MB
 - ~20000 collisions, ~86000 tracks
 - Generally, 3~4 tracks per collision (decay 3, collision 4)
 - If storeCollision is turned off, required space is much smaller

gSystem->Load("UrgmdDst/UrgmdDst.so"); root [1] UrgmdDst* u=new UrgmdDst(); root [2] TChain* c=u->initInputChain("out.root") Read in root file out.root . Total 1 file has been read in. root [3] c->GetEntries() (1ong 1ong) 10 root [4] c->GetEntry(0); <u>oot [5] u->event()->b()</u> (double) 0.0000000 root [6] u->numberOfTimesteps() (unsigned int) 1 root [7] u->numberOfCollisions() (unsigned int) 20629 root [8] u->numberOfTracks() (unsigned int) 86390 root [9] u->timestep(0)->startTrack() (unsigned int) 76058 root [10] u->track(76057)->collision() (unsigned int) 20628

root [0] gSystem->Load("UrqmdDst/UrqmdDst.so"); root [1] UrqmdDst* u=new UrqmdDst(); root [2] TChain* c=u->initInputChain("test.list") Read in root file out.root . Total 1 file has been read in.