## Update on UrQMD Production and Reading

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## Overview

Source: ui:/ustcfs/HICUser/fsi/urqmd\_250421.tar.gz

- Location: ui:/ustcfs/HICUser/fsi/urqmd
- Subdirectories
  - urqmd (original UrQMD production code)
  - curqmd (this production code)
  - run (script for production)
  - UrqmdDst (classes for storage)
  - UrqmdUtilities (new since current version)
  - PdgData (new since current version)

• C++98 is supported (PdgData requires C++11)

## Introduction: \$Location/urqmd

- Taken from https://itp.uni-frankfurt.de/~bleicher/index.html?content=urqmd
- Change: pythia6409.f: -> pytahi6428.f (final PYTHIA6, https://pythia.org/)
- Change: mk/Linux.mk: new gfortran option -std=legacy
  Otherwise unable to be compiled with high versions of gfortran
- Change: GNUmakefile: pythia6409 -> pytahi6428

## Introduction: \$Location/curqmd

• For UrQMD production and its storage in ROOT Tree

• Contents:

- main.{cxx,h}: C++ -> (Fortran <-> C++)
- c\_%.f: modified from \$Location/urqmd/%.f

• GNUmakefile

- Calls \$Location/urqmd/GNUmakefile
- Calls \$Location/UrqmdDst/GNUmakefile
- Compiles c\_\*.f and main.cxx
- Links \$Location/urqmd/\*.o (only those different from \$Location/curqmd/c\_\*.o), \$Location/curqmd/{c\_\*.o,main.o} to executable main

### Introduction: \$Location/run

#### • Contents:

- in.urqmd: input configuration of UrQMD production
- random.C: ROOT code for random number generation to random.list
  - Input: n (random number counter, required), name (file name for existing numbers, default="")
  - Output: file random.list containing n different random numbers (beginning with existing ones)
    - 1<= random number <= maximum of 4-byte integer
  - Used for UrQMD input seed for random number generator
  - Able to read existing numbers since current version
- run.bash: script for calling UrQMD production
  - Input: index (serial number for the job)
  - Sets the seed as the index-th number in random.list and saves the updated in.urqmd to in/\${index}.urqmd
  - Output: out/\${index}.root (event IDs begin with \${index}\*(#events per job))

### Introduction: \$Location/run

#### • Contents:

- run.condor: condor configuration template
- run\_ui.condor: condor configuration template with specific settings for ui
- submit.bash: script for submitting condor jobs using run.condor
  - Input: number (number of jobs, default=100)
  - Saves the updated run.condor to script/run\_\${number}.condor
  - Log files (error, log, output) saved in log/job\_\$(Process).{err,log,out}
- resubmit.bash: script for resubmitting condor jobs
  - Input: condor configuration file (e.g. script/run\_\${number}.condor)
  - Resubmits jobs with errors or incomplete jobs

## Introduction: \$Location/UrqmdDst

- Classes for storage of UrQMD data
- Contents:
  - UrqmdDst.{cxx,h}: saves pointers to sub-structures and statuses of them
  - UrqmdEvent.{cxx,h}: saves event information
  - UrqmdColllision.{cxx,h}: saves collision information
  - UrqmdTrack.{cxx,h}: saves track information
  - UrqmdTrackEx.{cxx,h}: saves track extended information
  - GNUmakefile
    - Compiles code in this directory to (lib)UrqmdDst.so

#### Introduction: class UrqmdDst

- Useful functions:
  - UrqmdEvent\* event()
  - UrqmdTimestep\* timestep(UInt\_t i)
  - UrqmdCollision\* collision(UInt\_t i)
  - UrqmdTrack\* track(UInt\_t i)
  - UrqmdTrackEx\* trackEx(UInt\_t i)
  - o UInt\_t numberOfTimesteps()
  - o UInt\_t numberOfCollisions()
  - o UInt\_t numberOfTracks()
  - TChain\* initInputChain(TString inputFileName, const TString inputTreeName = "urqmd")
    Reads UrqmdDst tree from file (\*.root or \*.list)
  - TTree\* initOutputTree(const TString outputTreeName = "urqmd")
    - Creates UrqmdDst tree for writing

#### Introduction: class UrqmdEvent

- Useful functions:
  - o UInt\_t id()
    - Serial number of event
  - Double\_t b(): impact parameter

### Introduction: class UrqmdTimestep

- Useful functions:
  - Double\_t time(): time of this timestep
  - o UInt\_t nPart(): not defined by UrQMD
    - Calculated by (total # projectile and target nucleons) (# tracks with parentCollisionType%100==0)
      parentCollisionType%100==0 means particle did not go through inelastic scattering
  - UInt\_t nColl(): nElasticColl+nInelasticColl+nBlockedColl
  - o UInt\_t nElasticColl(), UInt\_t nInelasticColl(), UInt\_t nBlockedColl()
  - o UInt\_t nDecays(), UInt\_t nHardRes(), UInt\_t nSoftRes(), UInt\_t nDecayRes()
  - UInt\_t nTracks(): number of tracks at this timestep
  - UInt\_t startTrack(): "nTracks" tracks starting from track index "startTrack" belong to this timestep
- If CTOption(4)==1, original UrQMD outputs initial configuration (t==0)
  - This code saves timestep at t==0

## Introduction: class UrqmdTimestep

- Useful functions:
  - Double\_t time(): time of this timestep
  - o UInt\_t nPart(): not defined by UrQMD
    - Calculated by (total # projectile and target nucleons) (# tracks with parentCollisionType%100==0)
      parentCollisionType%100==0 means particle did not go through inelastic scattering
  - UInt\_t nColl(): nElasticColl+nInelasticColl+nBlockedColl
  - UInt\_t nElasticColl(), UInt\_t nInelasticColl(), UInt\_t nBlockedColl()
  - UInt\_t nDecays(), UInt\_t nHardRes(), UInt\_t nSoftRes(), UInt\_t nDecayRes()
  - UInt\_t nTracks(): number of tracks at this timestep
  - UInt\_t startTrack(): "nTracks" tracks starting from track timestep
- If CTOption(4)==1, original UrQMD outputs initial
  - This code saves timestep at t==0

colum	n#	contents
0		"E" (only in016)
1		# of collisions
2		# of elastic collisions
3	$\checkmark$	# of inelastic collisions
4		# of Pauli-blocked collisions
5		# of decays
6		# of produced <i>hard</i> baryon resonances
7	$\checkmark$	# of produced <i>soft</i> baryon resonances
8		# of baryon resonances produced via a decay of another resonance

if ( (itypt(1).eq.ityp(inew(i))).and.
 (iso3t(1).eq.iso3(inew(i))) ) then
 origin(inew(i))=origint(1)+100
 uid(inew(i))=uidt(1)
elseif ( (itypt(2).eq.ityp(inew(i))).and.
 (iso3t(2).eq.iso3(inew(i))) ) then
 origin(inew(i))=origint(2)+100
 uid(inew(i))=uidt(2)

## Introduction: class UrqmdCollision

- (Serial number (starting from 1) not saved, equal to index; 0th collision always empty)
- Useful functions:
  - Double\_t time(): time of this collision
  - Double\_t sqrtS():  $\sqrt{s}$  of this collision
  - Double\_t sigmaTotal(): total cross section (mbarn)
  - Double\_t sigmaPartial(): partial cross section (mbarn) of actual subprocess
  - Double\_t rhoB(): baryon density (in units of  $\rho_0$ ) at collision point
  - UInt\_t type(): type index of process defined by UrQMD
  - UInt\_t nIn(): number of incident particles in this collision
  - UInt\_t nOut(): number of outgoing particles in this collision
  - o UInt\_t nTracks(): nIn+nOut
  - UInt\_t startTrack(): "nTracks" tracks starting from track index "startTrack" belong to this collision; first "nIn" tracks are incident, last "nOut" tracks are outgoing

co	olumn#	format	contents
1		(i8)	number of ingoing particles Nin
2	$\checkmark$	(i8)	number of outgoing particles Nout
3	$\checkmark$	(i4)	process ID (see Table 11)
4		(i7)	collision/entry counter
5	$\checkmark$	(f8.3)	collision time $t_{\rm coll}$ in fm/c
6	$\checkmark$	(e12.4)	center of mass energy of the collision $\sqrt{s}$ in GeV
7	$\checkmark$	(e12.4)	total cross-section of the collision $\sigma_{\rm tot}$ in mbarn
8		(e12.4)	partial cross-section of the actual sub-process $\sigma_i$ in mbarn
9		(e12.4)	Baryon density at collision point $\rho_{\rm B}$ in units of $\rho_0$

- Useful functions:
  - UInt\_t id(): serial number (not particle type ID)
  - o UInt\_t timestep()
    - If timestep==0 and collision!=0, this track belongs to that collision
  - UInt\_t collision(): collision==0 should be ignored
    - If collision==0 (even if timestep==0), this track belongs to that timestep
  - Int\_t iType(): type index according to isospin defined by UrQMD
  - Int\_t i3(): double 3rd component of isospin
  - o Int\_t q(): double charge
  - Int\_t s(): double spin (not saved by original UrQMD, may be ineffective)
  - Double\_t m(): mass (in UrQMD, may not be equal to that in PDG)
  - o Double\_t x(), Double\_t y(), Double\_t z()
  - o Double\_t px(), Double\_t py(), Double\_t pz()

<ul> <li>Useful functions:</li> </ul>	<b>0</b> 013	<b>O</b> O14	0 <sup>015</sup>	<b>O</b> O16
• UInt t id()· serial number (not particle type ID)	1 2	1 2	2 3	1 2
	3	3 4	4 5	3 √ 4 √
• UInt_t timestep()	5	5	6	5
• If timestep==0 and collision!=0, this track belongs to that coll	6 7	6 7	7 8	6 √ 7 √
• UInt t collision(): collision==0 should be ignored	8 9	8 9	9 10	8 √ 9 √
	10	10	11	10 🗸
• If collision==0 (even if timestep==0), this track belongs to the	11	11	12	11 V 12 V
• Int_t iType(): type index according to isospin defined by 1	13 14	13 14	14 15	13 14
• Int_t i3(): double 3rd component of isospin	15	15	16	15
• Int_t q(): double charge	16		17	
Int t s(): double spin (not sayed by original UrOMD may	18			
• mt_t s(). double spin (not saved by original OrQMD, may	19 20			
• Double_t m(): mass (in UrQMD, may not be equal to that	21			
• Double $t x()$ , Double $t y()$ , Double $t z()$	22	1/*		
$- \sqrt{2} - \sqrt{2} - \sqrt{2}$		10* 17*		
• Double_t px(), Double_t py(), Double_t pz()		18*		1
		19*		N

O <sup>013</sup>	<b>O</b> 014	<b>O</b> 015	016	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 <sub>coll</sub> 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				Efr : freeze-out energy of particle					
21				$p_x^{\rm fr}$ : freeze-out momentum x component					
22				$p_y^{\rm fr}$ : freeze-out momentum y component					
23				$p_z^{\rm fr}$ : freeze-out momentum z component					
	16*			$\tau_{dec}$ decay time of particle					
	17*			$\tau_{form}$ formation time of particle					
	18*			$R_{\sigma}$ cross section reduction factor					
	19*		√	unique particle number (not ID!)					
			16*	ityp <sub>1</sub> <sup>old</sup> : particle-ID of parent particle # 1					
			17*	ityp <sub>2</sub> <sup>old</sup> : particle-ID of parent particle # 2					

#### • Useful functions:

- o UInt\_t nCollisions()
- UInt\_t lastCollision(): index of last collision experienced by this particle
- UInt\_t parentCollision(): index of parent collision creating this particle
  - $\circ$  Should <= lastCollision, because elastic scattering (A+B -> A+B) not recorded as parent collision
  - Not saved in UrQMD
  - To encode this and lastCollision, a variable "lstColl" in UrQMD is changed
- o UInt\_t parentCollisionType()
  - $\circ$  = nElasticScatterings\*100 + collisionType (>=0, <100) according to UrQMD
  - If only collision type required, calculate parentCollisionType%100
- o Int\_t iTypeParent1(), Int\_t i3Parent1()
- o Int\_t iTypeParent2(), Int\_t i3Parent2()
  - Parent i3 not saved in original UrQMD
  - To encode these four, a variable "origin" in UrQMD is changed

	<b>~</b> 013	<b>~</b> 014	<b>_015</b>	<b>_016</b>	contents
• Useful functions:		Ŭ	1		ind: index of particle (see CTOption (56))
Should be last colligion number	1	1	2	1	t: time of particle
• UInt t nCollisions()	2	2	3	2	$r_x$ : x coordinate
(prevents reaction between 2 particles just reacted)	3	3	4	3 V	$r_y$ : y coordinate
• UInt_t lastCollision(): index of last collision experienced	5	4 5	6	4 <b>∨</b> 5	$r_z$ : 2 coordinate E: energy of particle
	6	6	7	6 🗸	$p_x$ : x momentum component
• Ulnt t parentCollision(): index of parent collision creating	7	7	8	7 🗸	$p_y$ : y momentum component
	8	8	9	8 🗸	$p_z$ : z momentum component
• Should $\leq$ =lastCollision, because elastic scattering (A+B $\rightarrow$ A)	9	9	10	9 1	m: mass of particle
	10	10	11	10 V	ityp: particle-ID
• Not saved in UrQMD	11	11	12	12	$1805: 2 \cdot I_3$ (see Section [1.2])
	13	12	14	13 1	barent collision number (see Table 10)
• To encode this and lastCollision, a variable "IstColl" in UrQN	14	14	15	14 🗸	N <sub>coll</sub> number of collisions
			16		S: strangeness
• Unt_t parentCollision Type()	15	15		15 🗸	parent process type (see Table III)
$\Gamma_{1}$ $\Gamma_{2}$ $\Gamma_{1}$ $\Gamma_{2}$ $\Gamma_{1}$ $\Gamma_{2}$ $\Gamma_{2$	- 16		17		history information (debugging only)
$\circ = nElasticScatterings*100 + collisionType (>=0, <100) accord$	16				$t^{\rm fr}$ : freeze-out time of particle
o If only colligion type required coloulate nerent Colligion Type	18				$r_x$ : freeze-out x coordinate
• If only consion type required, calculate parentConsion type?	19				$r_{*}^{\text{fr}}$ : freeze-out z coordinate
Int tiTypoPoront1() Int ti2Poront1()	20				$\tilde{E^{\text{fr}}}$ : freeze-out energy of particle
<sup>o</sup> Int_t Hyper atent (), Int_t ISr atent ()	21				$p_x^{\rm fr}$ : freeze-out momentum x component
Int tiTupoDoront?() Int ti2Doront?()	22				$p_y^{\rm fr}$ : freeze-out momentum y component
<sup>o</sup> Int_t TypeParent2(), Int_t TSParent2()	23	1/*			$p_z^{\rm Ir}$ : freeze-out momentum z component
· Derent is not caual in original UrOMD		10° 17*			$\tau_{dec}$ decay time of particle
• Falent 15 not saved in original OrQIVID		18*			$R_{-}$ cross section reduction factor
• To encode these four, a variable "origin" in UrOMD is change		19*		√ √	unique particle number (not ID!)
To cheode these tour, a variable origin in origin to change				16* 🗸	ityp <sub>1</sub> <sup>old</sup> : particle-ID of parent particle # 1
				17* 🗸	ityp20ld : particle-ID of parent particle # 2

- Index always equal to UrqmdTrack
- In origin UrQMD, this information not saved for tracks belong to collisions
  Can be saved via this code
- Useful functions:
  - Double\_t tFreezeOut(),
  - Double\_t xFreezeOut(), Double\_t yFreezeOut(), Double\_t zFreezeOut()
  - o Double\_t pxFreezeOut(), Double\_t pyFreezeOut(), Double\_t pzFreezeOut()
  - Double\_t tauDecay(): decay time
  - Double\_t tauForm(): formation time
  - Double\_t rSigma(): cross section reduction factor

	<b>_013</b>	0014	<b>_015</b>	016	contents
• Index always equal to Urqmd Irack			1	$\checkmark$	ind: index of particle (see CTOption (56))
	1	1	2	1	t: time of particle
	2	2	3	$2 \sqrt{2}$	$r_x$ : x coordinate
	3	3	4		$r_y$ : y coordinate
• In origin UrOMD this information not saved for track	5	5	6	5	$F_z$ : energy of particle
In origin or givib; und information not baved for track	6	6	7	6 🗸	$p_x$ : x momentum component
• Can be saved via this code	7	7	8	7 🗸	$p_y$ : y momentum component
Can be saved via this code	8	8	9	8 🗸	$p_z$ : z momentum component
	9	9	10	9 🗸	m: mass of particle
	10	10	11	10 1	ityp: particle-ID
I Jacful functional	12	12	12	$11$ $\sqrt{12}$	$1505: 2 \cdot I_3$ (see Section [1.2])
• Useful functions:	13	13	14	13 1	parent collision number (see Table 10)
$\Delta D_{avel} = \frac{1}{2} \left( \frac{1}{2} + \frac{1}{2} \frac{1}{2} - \frac{1}{2} \frac{1}{2} \right)$	14	14	15	14 🗸	N <sub>coll</sub> number of collisions
<sup>o</sup> Double_t tFreezeOut(),			16		S : strangeness
Double tu Encare Out() Double tu Encare Out() Double	15	15		15 🗸	parent process type (see Table 11)
<sup>o</sup> Double_t xFreezeOut(), Double_t yFreezeOut(), Double_	16		17		history information (debugging only)
$\mathbf{D}$	17				$r^{\text{fr}}$ : freeze-out x coordinate
<sup>o</sup> Double_t pxfreezeOut(), Double_t pyfreezeOut(), Doubl	18				$r_{y}^{\text{fr}}$ : freeze-out y coordinate
Develate the Decent (), decent times	19			$\checkmark$	$r_z^{\rm fr}$ : freeze-out z coordinate
<sup>o</sup> Double_t tauDecay(): decay time	20				$E^{\rm fr}$ : freeze-out energy of particle
$\mathbf{D}$	21				$p_x^{\text{fr}}$ : freeze-out momentum x component
• Double_t tauForm(): formation time	22				$p_y^{\text{ir}}$ : freeze-out momentum y component
$\mathbf{D}_{\mathbf{r}} = 1 1_{\mathbf{r}} + \mathbf{C}_{\mathbf{r}} + \mathbf{C}_{r$	23	16*		N	$p_z$ : freeze-out momentum z component
• Double_t rSigma(): cross section reduction factor		17*		√	$\tau_{dec}$ decay time of particle $\tau_{form}$ formation time of particle
		18*			$R_{\sigma}$ cross section reduction factor
		19*		$\checkmark$	unique particle number (not ID!)
				16* 🔨	ityp <sup>old</sup> : particle-ID of parent particle # 1
				17* 🗸	ityp <sup>ord</sup> : particle-ID of parent particle # 2

# **Configuration File**

#### • Before xxx

- The same configuration as original UrQMD
- f14 -> does not save f14 output file (#f14 -> saves f14 output file)

#### • After xxx

- Specific configuration for this code
- #ske -> skips empty events (Ncoll + Ndecays == 0)
  - For non-empty events, Npart (only counts inelastic collisions) may be 0 (can be manually skipped when reading)
- #qut -> quiet standard output
- col -> does not save collision information
- tke -> does not save track extended information
- Removed since this version: esd -> does not use external seed for random



## How to save parent particle information

• 4-byte integer: "origin" (original definition)

- = parentCollisionType + 100\*(# elastic scatterings)
  - + 1000\*(|iType\_parent1| + 1000\*|iType\_parent2|)
- 0 <= parentCollisionType < 100, 0 <= |iType\_ parent| < 1000
- No "i3\_parent" or sign of "iType\_parent"



parentCollisionType = origin%100
# elastic scatterings = (origin/100)%10

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

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

ityn	0-+	ityn	1	ityn	0++	ityn	1++	ityn	charmed
ng p	0	101	-		•	ng p	-	100	D
101	$\pi$	104	$\rho$	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	$J/\Psi$
107	$\eta'$	109	$\phi$	112	$f_0^*$	116	$f'_1$	136	$\chi_c$
ityp	1+-	ityp	$2^{++}$	ityp	(1)*	ityp	(1)**	137	$\Psi'$
122	$b_1$	118	$a_2$	126	$\rho_{1450}$	130	$\rho_{1700}$	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	$\omega_{1420}$	131	$\omega_{1662}$		
124	$h_1'$	120	$f'_2$	128	$\phi_{1680}$	132	$\phi_{1900}$		

## How to save parent particle information

• 4-byte integer: "origin" (new definition)

- = parentCollisionType + 100\*(# elastic scatterings)
  - $+ 1000*((i3_parent1+3)+7*(t_iType_parent1+100))$
  - + 1400000\*((i3\_parent2+3)+7\*(t\_iType\_parent2+100))
- $\circ$  0 <= parentCollisionType < 100, 0 <= i3\_parent+3 <= 6
- t\_iType\_parent=iType\_parent; if(>=100) -=40; if(<=-100) +=40
  - $0 < t_iType_parent + 100 < 200$

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

- !!!Also affects \*.txt output
- parentCollisionType = origin%100
  # elastic scatterings = (origin/100)%10

ityp	nucleon	ityp	delta	ityp	lambda	ityp	sigma	ityp	xi	ityp	omega
1	$N_{938}$	17	$\Delta_{1232}$	27	$\Lambda_{1116}$	40	$\Sigma_{1192}$	49	$\Xi_{1317}$	55	$\Omega_{1672}$
2	$N_{1440}$	18	$\Delta_{1600}$	28	$\Lambda_{1405}$	41	$\Sigma_{1385}$	50	$\Xi_{1530}$		
3	$N_{1520}$	19	$\Delta_{1620}$	29	$\Lambda_{1520}$	42	$\Sigma_{1660}$	51	$\Xi_{1690}$		
4	$N_{1535}$	20	$\Delta_{1700}$	30	$\Lambda_{1600}$	43	$\Sigma_{1670}$	52	$\Xi_{1820}$		
5	$N_{1650}$	21	$\Delta_{1900}$	31	$\Lambda_{1670}$	44	$\Sigma_{1775}$	53	$\Xi_{1950}$		
6	$N_{1675}$	22	$\Delta_{1905}$	32	$\Lambda_{1690}$	45	$\Sigma_{1790}$	54	$\Xi_{2025}$		
7	$N_{1680}$	23	$\Delta_{1910}$	33	$\Lambda_{1800}$	46	$\Sigma_{1915}$				
8	$N_{1700}$	24	$\Delta_{1920}$	34	$\Lambda_{1810}$	47	$\Sigma_{1940}$				
9	$N_{1710}$	25	$\Delta_{1930}$	35	$\Lambda_{1820}$	48	$\Sigma_{2030}$				
10	$N_{1720}$	26	$\Delta_{1950}$	36	$\Lambda_{1830}$						
11	$N_{1900}$			37	$\Lambda_{1890}$						
12	$N_{1990}$			38	$\Lambda_{2100}$						
13	$N_{2080}$			39	$\Lambda_{2110}$						
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	$\pi$	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	$J/\Psi$
107	$\eta'$	109	$\phi$	112	$f_0^*$	116	$f'_1$	136	$\chi_c$
ityp	1+-	ityp	$2^{++}$	ityp	(1)*	ityp	(1)**	137	$\Psi'$
122	$b_1$	118	$a_2$	126	$\rho_{1450}$	130	$\rho_{1700}$	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	$\omega_{1420}$	131	$\omega_{1662}$		
124	$h'_1$	120	$f_2'$	128	$\phi_{1680}$	132	$\phi_{1900}$		

## How to store parent collision information

- Original UrQMD saves serial number of last collision
  - "lstColl" in code; parent collision number in users guide (better to say last collision number)
  - To prevent reaction between 2 particles from the same collision
  - If a particle goes through elastic collision (A+B -> A+B), "lstColl" is recorded, and its parent collision (in which collision the particle is produced) is not known
- In this code, both last collision and parent collision are saved
  - parentCollision <=lastCollision, because elastic scattering not recorded as parent collision
  - "lstColl" changed to encode both quantities (decoded when required in UrQMD simulation)
    - lastCollision = integer((sqrt(1+8.0d0\*lstcoll)-1)/2)
    - For odd "lastCollision", parentCollision = lstColl-(lastCollision+1)/2\*lastCollision
    - For even "lastCollision", parentCollision = lstColl-lastCollision/2\*(lastCollision+1)

#### • !!!Also affects \*.txt output

#### Check

root [46] u->co1	lision(2634)->p	rint()									
5.0000e+01	9.2480e-01	0.0000e+00	0.0000e <sup>.</sup>	+00	1.0897e <sup>.</sup>	-02	20	1	2	3	10190
root [47] u->tra	ack(10190)->prin	t()									
5543 0	2634 104	0 0	-2	9.2490e-	-01	6.2051e-	+00	4.9854e	+00	-4.7102	e+01 3?
.0487e-01	3.3380e-01	-2.9844e+00	1	2575	2575	20	-117	1	0	0	
root [48] u->tra	ack(10191)->prin	t()									
5666 0	2634 101	2 1	0 _	1.3800e-	-01	6.2051e	+00	4.9854e	+00	-4, 7102	e+01 1?
.4145e-01	4.8468e-01	-2.7690e+00	2	2634	2634	20	104	0	0	0	
root [49] u->tra	ack(10192)->prin	t()									
5667 0	2634 101	-2 -1	0	1.3800e-	-01	6.2051e-	+00	4.9854e	+00	-4,7102	e+01 1
.6339e-01	-1.5085e-01	-2.1552e-01	1	2634	2634	20	104	0	0	0	

	1 2 20	2634 50.000 (	).9249E+00 0.0000E+00 0.0000E+0	0 0.1090E-01							
2458	0.50000000E+02	0.62052729E+01	0.49856325E+01 -0.47099683E+02	0.31569740E+01	0.30484766E+00	0.33381796E+00 -0.29844185E+01	0.92485896E+00	104 0 0	2575	1 0	117020
2458	0.5000000E+02	0.62052729E+01	0.49856325E+01 -0.47099683E+02	0.28179373E+01	0.14146304E+00	0.48467783E+00 -0.27688991E+01	0.13800000E+00	101 2 1	2634	2 0	104020
2523	0.50000000E+02	0.62052729E+01	0.49856325E+01 -0.47099683E+02	0.33903666E+00	0.16338462E+00	-0.15085987E+00 -0.21551943E+00	0.13800000E+00	101 -2 -1	2634	1 0	104020

#### Check

root [51] u-1	>collision(1759)-	>print()										
2.1062e+01	2.7930e+00	8.8867e+00	2. 122	6e+00	2.7710	e+00	26	2	2	4	7333	
root [52] u-1	>track(7333)->pri:	nt()										
4063 0	1759 101	2 1	0	1.3800	le-01	-2.170	)4e+00	2.720	2e+00	-2.08	98e+01	-
2.0889e-01	-1.1183e-01	-7.8552e-01	1	1742	1742	15	17	-1	16	1		
root [53] u-1	>track(7334)->pri:	nt()										
3512 0	1759 17	1 1	-1	1.2320	le+00	-2.284	17e+00	2.916	0e+00	-2.09	26e+01	6
.0555e-02	2.7295e-01	-6.6250e+01	7	1366	1366	15	17	1	17	-1		
root [54] u-1	>track(7335)->pri:	nt()										
3512 0	1759 17	1 1	-3	1.2320	le+00	-2.284	17e+00	2.916	0e+00	-2.09	26e+01	- 7
.0203e-01	2.7917e-01	-5.7695e+01	8	1759	1366	115	17	1	17	-1		
root [55] u-1	>track(7336)->pri:	nt()										
4063 0	1759 101	2 1	0	1.3800	le-01	-2.170	)4e+00	2.720	2e+00	-2. 08	98e+01	-
8.5046e-01	-1.1807e-01	-9.3301e+00	2	1759	1742	115	17	-1	16	1		

	2 2 26	1759 21.062 0	).2793E+01 0.888	7E+01 0.2123E+0	1 0.2771E+01									
1869	0.21061567E+02	-0.21704844E+01	0.27203925E+01	-0.20897817E+02	0.83200094E+00	-0.20889727E+00	-0.11183377E+00	-0.78551684E+00	0.13800000E+00	101	2 1	1742	1 0	1601701
351	0.21061567E+02	-0.22844273E+01	0.29161623E+01	-0.20926401E+02	0.66255078E+02	0.60558097E-01	0.27293928E+00	-0.66243033E+02	0.12320000E+01	17		1366	70	1701701
351	0.21061567E+02	-0.22844273E+01	0.29161623E+01	-0.20926401E+02	0.57715714E+02	0.70206737E+00	0.27917483E+00	-0.57697617E+02	0.12320000E+01	17		1759	80	1701711
1869	0.21061567E+02	-0.21704844E+01	0.27203925E+01	-0.20897817E+02	0.93713649E+01	-0.85040655E+00	-0.11806933E+00	-0.93309327E+01	0.13800000E+00	101	2 1	1759	2 0	1601711
	2 4 15	1742 20 573	0 4078E+01 0 400	)OE+O2 0 7113E+O	1 0 3180E+01									
200		2,2205500171,01	0.17000101 0.400	0.004045077.00	0 175255277.00	0 702420217.00	0 140014000 01	0 174755000.00	0 101067038.01	17	1 0	1740	F 0	110111
289	0.205/3349E+02	-0.23855281E+01	0.1/656513E+01	-0.20424527E+02	0.1/53552/E+02	-0.78342021E+00	0.14621490E-01	-0.1/4/5502E+02	0.12196703E+01	±/-	-1 0	1/40	5 0	110111
231	0.20573349E+02	-0.20479036E+01	0.27860165E+01	-0.20436877E+02	0.47860026E+02	0.36748606E-01	-0.68083006E+00	-0.47785087E+02	0.25889575E+01	16	1 1	1462	7 0	112322
289	0.20573349E+02	-0.23855281E+01	0.17656513E+01	-0.20424527E+02	0.21800109E+02	-0.89090823E+00	0.53665146E-02	-0.21747027E+02	0.12320000E+01	17 -	10	1742	60	1601701
231	0.20573349E+02	-0.20479036E+01	0.27860165E+01	-0.20436877E+02	0.36313948E+02	0.52319997E+00	-0.39157334E+00	-0.36284463E+02	0.13090051E+01	17 -	·1 0	1742	8 0	1601701
1871	0.20573349E+02	-0.20479036E+01	0.27860165E+01	-0.20436877E+02	0.64494956E+01	-0.17006607E+00	-0.16816797E+00	-0.64435818E+01	0.13800000E+00	101	0 0	1742	1 0	1601701
1872	0.20573349E+02	-0.20479036E+01	0.27860165E+01	-0.20436877E+02	0.83200094E+00	-0.20889727E+00	-0.11183377E+00	-0.78551684E+00	0.13800000E+00	101	2 1	1742	1 0	1601701

## Update on writting

• Add forward compatibility for ROOT version < 6.30

- Error when ROOT files written by ROOT >= 6.30 are read by ROOT < 6.30
  - Solution 1: file->SetBit(TFile::k630forwardCompatibility);
  - Solution 2: put TFile.v630forwardCompatibility: true in ~/.rootrc
    - Fully valid for ROOT 6.32 (>= 6.32.10) and ROOT 6.34 (>=6.34.04)
    - ROOT on ui is 6.32.02

### **Known Issues**

- Negligible particles (such as some c-baryons) do not have UrQMD iType indices
  - In Original UrQMD, PDG ID saved in iType (4-byte integer)
  - $\circ$  In this code, only values between  $\pm\,100$  are reserved for iType
    - iType == 0 is saved for these particles (can be identified using mass)
- Some UrQMD iType indices may not correspond to PDG ID
  Current data structure (iType, i3) not affected
- Particles of the same species in UrQMD may not have the same masses
  For proton, UrQMD mass = 0.938, some protons may < 0.938</li>

### **Known Issues**

- Different simulation results from CentOS 7 gfortran 4.8.5 and Ubuntu gfortran 11+
   Unknow reason
- UrQMD simulation using some input seeds may be broken (exit with error)
  - Both original UrQMD and this code
  - Solution: input a new seed and resubmit the job

## Introduction: \$Location/UrqmdUtility

- Contents:
  - UrqmdUtility.{cxx,h}
  - GNUmakefile
    - Compiles Fortran \$Location/urqmd/{blockres.f,error.f,ityp2pdg.f,upmerge.f} and C++ code in this directory to (lib)UrqmdUtility.so
- Useful functions:
  - static Int\_t getPdgId(Int\_t iType, Int\_t i3)
  - o static Int\_t getPdgId(std::pair<Int\_t, Int\_t>id)
  - static TString getParticleName(Int\_t iType)
  - static void getUrqmdId(Int\_t& iType, Int\_t& i3, Int\_t pdgId)
  - o static std::pair<Int\_t, Int\_t> getUrqmdId(Int\_t pdgId)
  - o static Int\_t getUrqmdIType(Int\_t pdgId)
  - o static Int\_t getUrqmdI3(Int\_t pdgId)

## Introduction: \$Location/PdgData

- Contents:
  - ParticleDataEntry.{cxx,h}
  - PdgData.{cxx,h}
  - GNUmakefile
    - Compiles code in this directory to (lib)PdgData.so
- Useful functions:
  - PdgData: static const ParticleDataEntry\* get(const Int\_t pid)
  - ParticleDataEntry: Int\_t pid(): PDG particle ID
  - ParticleDataEntry: Int\_t spin(): double spin
  - ParticleDataEntry: Int\_t charge(): triple electric charge
  - ParticleDataEntry: Int\_t color()
  - o ParticleDataEntry: Double\_t mass()

# Usage

- By ROOT: gSystem->Load(\$Location/UrqmdDst/libUrqmdDst.so");
- By GCC: -L\$Location/UrqmdDst -lUrqmdDst -Wl,-rpath,\$Location/UrqmdDst
- How to read tree

0

- o UrqmdDst\* u = new UrqmdDst();
- TChain\* c = u->initInputChain("PATH\_OF\_FILE");//\*.root or \*.list
- o for(Long64\_t iEntry=0; iEntry<c->GetEntries(); iEntry++){
- c->GetEntry(iEntry);

. . .

- o for(UInt\_t iTimestep=0; iTimestep<u->numberOfTimesteps(); iTimestep++){
- const UrqmdTimestep\* timestep = urqmdDst->timestep(iTimestep);
- o for(UInt\_t iTrack=0; iTrack<timestep->nTracks(); iTrack++){
- const UrqmdTrack\* track = u->track(timestep->startTrack()+iTrack);

# Usage

- By ROOT: gSystem->Load(\$Location/UrqmdUtility/libUrqmdUtility.so");
- By GCC: -L\$Location/UrqmdUtility -lUrqmdUtility -Wl,-rpath,\$Location/UrqmdUtility
- By ROOT: gSystem->Load(\$Location/PdgData/libPdgData.so");
- By GCC: -L\$Location/PdgData -lPdgData -Wl,-rpath,\$Location/PdgData

root [2] PdgData::get(UrqmdUtility::getPdgId(101,2))->mass()
(double) 0.13957000
root [3] UrgmdUtility::getUrgmdId(UrgmdUtility::getPdgId(101,2))
(std::pair <int_t, int_t="">) { 101, 2 }</int_t,>
root [4] UrgmdUtility::getParticleName(UrgmdUtility::getUrgmdIType(2212))
(TString) "Nukleon"[7]
root [5] UramdUtility::getParticleName(UramdUtility::getUramdIType(-2212)
(TString) "*Nukleon"[8]

# Summary

- A few updates on UrQMD production and reading
- Almost all information can be accessed with as small disk space required as possible
- New version compatible with older ones
- New codes UrqmdUtility and PdgData help with reading UrQMD data

- Existing UrQMD samples in ui:/ustcfs/HICUser/fsi/MODEL/UrQMD/
   Naming conventions: ProjectileTarget\_energy(\_configuration)
  - E.g., UU\_2p1, AuAu\_3\_eos1