Introduction to MCPL

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Practical: P. Willendrup





Acknowledgements:

1: DTU 2: ESS 3: CSNS

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Key MCPL features

MCPL: Monte Carlo Particle Lists

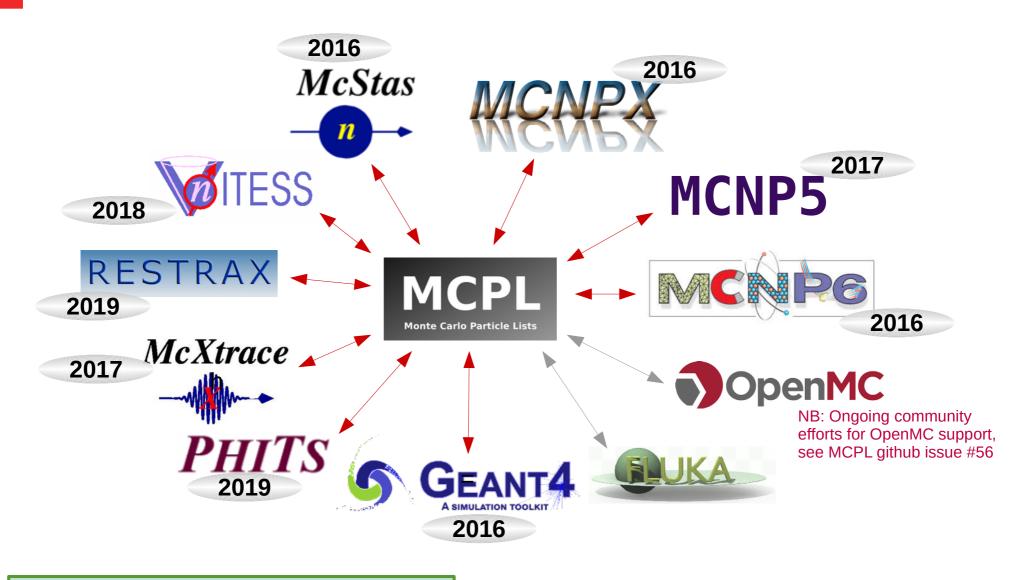
- It is a **simple** binary file-format. Each file contains a list of MC particles with enough info to seed simulations.
- MCPL files can contain *meta-data*. This makes it possible to tell what data is in a file, where it came from, how it should be interpreted.
- The format is **flexible**: can contain a lot of information if needed, or can contain only minimal information if small filesize is important. Can be gzip'ed.
- It is **easy** to make code dealing with MCPL, so it is easy to make plugins & converters for the various Monte Carlo frameworks. → End-users will simply use those converters.
- MCPL comes with **tools and APIs**, such as for inspecting or editing contents.
- **Well-defined** versioned format, focus on backwards compatibility.

MCPL background/philosophy

MCPL: Monte Carlo Particle Lists

- In principle simple: just a "bag of Monte Carlo particles", with properties such as particle type, energy, position, direction, weights, ...
- Goal of the MCPL project: Make this a new **standard particle**exchange format.
 - Original motivation: we needed to chain MCNP→McStas→Geant4
- To achieve this, we tried to make it **attractive to use**:
 - Have custom hooks for most major Monte Carlo particle codes.
 - Have cmdline tools and easy to use C/C++/Python API
- .. and we tried hard to avoid:
 - Annoying dependencies.
 - "MCPL is too bloated/slow for my usecase, I'll roll my own custom solution"
- So it must be **flexible in what is actually on-disk** (e.g. don't need "type" field in McStas output since it is always neutrons)
- But should **always look the same when opening the files** to make reading files trivial.

Codes with MCPL support







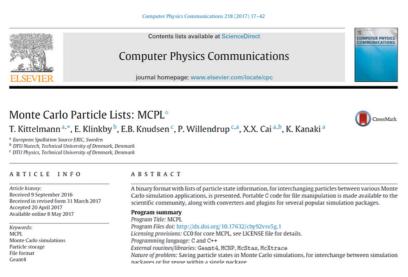


Download, follow, and report issues @GitHub

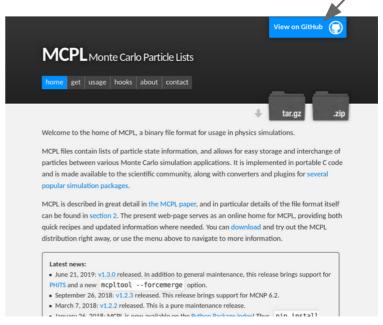
focus on availability:

- Extremely liberal license (CC0) encourage bundling.
- API for C/C++/Python code (all versions).
- "fat" single-file versions of all C code (even embedding zlib)
- Can "pip install" Python API+pymcpltool.

... and documentation:



Detailed paper for release 1.1.0: (DOI 10.1016/j.cpc.2017.04.012)



 Online docs with recipes (https://mctools.github.io/mcpl/)

What form does MCPL support take?

- Built-in support in instrument simulation codes:
 - McStas, McXtrace, VITESS, RESTRAX/SIMRES
 - Batteries included → great for users!

Most work done by developers of these applications!

- C++ helper classes for particle capture or event seeding available for Geant4 (in line with how most Geant4 users work) T. Kittelmann
- MCNP support relies on inbuilt ability to dump particles to/seed from "SSW" files.
 - We provide ssw2mcpl and mcpl2ssw tools.
- T. Kittelmann+E. Klinkby
- Somewhat high maintenance burden due to plethora of MCNP flavours + closed nature of programme.
- Complication is that particles need "surface ID". Can be provided as MCPL userflags or via global setting.
- mcpl2ssw must be provided with sample SSW files from target setup.
- PHITS support: Like MCNP, but simpler. More details later.

Data in MCPL files

All generic parameters always Available to reading code, no matter source of MCPL file.

Flexibility in how this is actually stored!

_			
	Particle state information		
	Field	Description	
	PDG code	32 bit integer indicating particle type.	
	Position	Vector, values in centimetres.	
	Direction	Unit vector along the particle momentum.	
	Kinetic energy	Value in MeV.	
	Time	Value in milliseconds.	
	Weight	Weight or intensity.	
	Polarisation	Vector.	
	User-flags	32 bit integer with custom info.	

FD can be single (22bit)

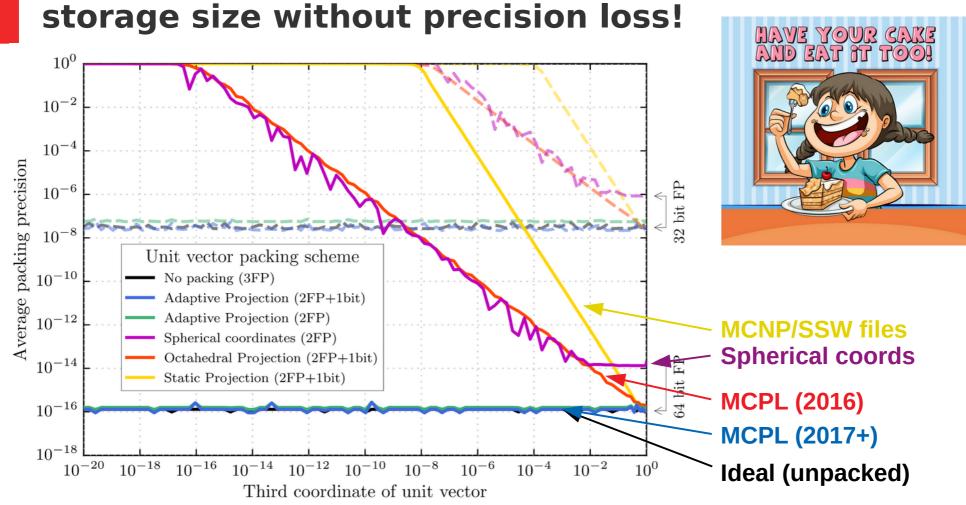
Detailed layout of the data associated with each particle in an MCPL file.

Particle data layout		FP can be single (32bit)	
Presence	Count & type	Description or double precision (64bit)	
OPTIONAL	3 × FP	Polarisation vector (if enabled in file).	
ALWAYS	$3 \times FP$	Position vector	
ALWAYS	$3 \times FP$	Packed direction vector and kinetic energy.	
ALWAYS	$1 \times FP$	Time.	
OPTIONAL	$1 \times FP$	Weight (if file does not have universal weight).	
OPTIONAL	$1 \times INT32$	PDG code (if file does not have universal PDG code).	
OPTIONAL	1 × UINT32	User-flags (if enabled in file).	

This implies from 28 to 96 bytes/particle. Already good, but most files are gzip'ed (by MCPL or user) and consume less. (NB: MCPL code can read .mcpl.gz files directly)



Novel packing of direction vectors: Optimal 2xFP



Breakdown of the Adaptive Projection Packing method, in which a unit vector, (u_x, u_y, u_z) is stored into two floating point numbers, FP1 and FP2, and one extra bit of information.

Adaptive Projection Packing						
Scenario	FP1	FP2	+1 bit	Packed signature		
$ u_x $ largest	$1/u_z$	u_y	$sign(u_x)$	FP1 > 1, FP2 < 1		
$ u_y $ largest	u_x	$1/u_z$	$sign(u_y)$	FP1 < 1, FP2 > 1		
$ u_z $ largest	u_x	u_y	$sign(u_z)$	FP1 < 1, FP2 < 1		



Example file Inspected with (py)mcpltool

```
Opened MCPL file recordfwd.mcpl.gz:
    Basic info
                         : MCPL-3
      Format
      No. of particles
                         : 542199
      Header storage
                         : 826 bytes
      Data storage
                         : 17350368 bytes
    Custom meta data
      Source
                                                                                 Custom meta-data
      Number of comments: 8

    This file is from ESS-DG Geant4

            -> comment 0 : "Created with the Geant4 MCPLWriter in the ESS/dgco
            -> comment 1 : "MPCLWriter volumes considered : ['RecordFwd']"

    Comments reminding us of setup

            -> comment 2 : "MPCLWriter steps considered : <at-volume-exit>"
                                                                                  used to create file
            -> comment 3 : "MPCLWriter write filter : <unfiltered>"
            -> comment 4 : "MPCLWriter user flags : <disabled>"

    Binary "blobs" keep more complete

            -> comment 5 : "MPCLWriter track kill strategy : <none>"
                                                                                  configuration details, here ESS-DG
            -> comment 6 : "ESS/dgcode geometry module : G4StdGeometries/GeoSl
            -> comment 7 : "ESS/dgcode generator module : G4StdGenerators/Simp
                                                                                  geo/gen parameters. Could be
      Number of blobs
                                                                                  McStas instrument file, input deck
            -> 74 bytes of data with key "ESS/dgcode_geopars"
                                                                                   from MCNP/PHITS, etc.
            -> 231 bytes of data with key "ESS/dgcode_genpars"
    Particle data format
      User flags
      Polarisation info
                        : no
      Fixed part. type
                         : no
                                                                        Columns of particle data
      Fixed part. weight : yes (weight 1)
                                            NB: compresses to
      FP precision
                           single
                                                                        In this file: No userflags or polarisation
                                             19.2bytes/particle
      Endianness
                         : little
      Storage
                         : 32 bytes/particle
                                                  y[cm]
  index
            pdgcode
                      ekin[MeV]
                                      x[cm]
                                                              z[cm]
                                                                              ux
                                                                                                      uz
                                                                                                            time[ms]
                                                 3.5344
                                                                        -0.43426
      0
                 22
                         1.2238
                                    -13.327
                                                                                   -0.036564
                                                                                                 0.90005
                                                                                                             0.14113
                 22
                                    -15.976
                                                 14.788
                                                                        -0.63971
                                                                                   0.082934
                                                                                                 0.76413
                        0.12059
                                                                                                             0.14113
                 22
                        0.10212
                                    -22.452
                                                -7.1864
                                                                 40
                                                                        -0.58735
                                                                                    -0.35527
                                                                                                 0.72718
                                                                                                             0.14113
                 22
                                     12.547
                                                 36.899
                                                                        0.19775
                                                                                                 0.85987
                          7.695
                                                                 40
                                                                                     0.47066
                                                                                                             0.20354
               2112
                                                                 40
                                                                                                              0.1829
                        2.5e-08
                                                                                                       1
                                                                         n 21854
                                                                                     0.33885
                                                                                                 0.46387
                                                                                                           0.0047377
PDG codes: 2112 = neutron, 22 = gamma
                                                                             666
                                                                                     0.38747
                                                                                                 0.91761
                                                                                                           0.0047367
                                                                             866
                                                                                   -0.075343
                                                                                                 0.97717
                                                                                                             0.12339
More at http://pdg.lbl.gov/2015/reviews/rpp2015-rev-monte-carlo-numbering.pdf
                                                                              0
                                                                                                              0.1829
                                                                                                              0.1829
```

C API

- Stable C API for reading/creating/editing MCPL
- Use to create most application-specific hooks
- Some users use it to analyse or tailor MCPL files

```
#include "mcpl.h"
void read example()
  mcpl file t f = mcpl open file("myfile.mcpl");
  const mcpl particle t* prtcl;
  while ( ( prtcl = mcpl read(f) ) ) {
    //<Access here: prtcl->ekin, prtcl->time, ...>
  mcpl close file(f);
```

C not C++ to support more apps (C is "lingua franca" of SW)

Despite being C, interface is "object oriented" and hopefully easy.

```
#include "mcpl.h"
void create example()
  mcpl outfile t f = mcpl create outfile("myfile.mcpl");
  mcpl hdr set srcname(f, "Custom C code");
  mcpl hdr add comment(f, "Just an example.");
  mcpl enable doubleprec(f);
  mcpl particle t * prtcl = mcpl get empty particle(f);
  for (i = 0; i < 1000; ++i)
    //<Set here: prtcl->ekin, prtcl->time, ...>
    mcpl add particle(f,prtcl);
  mcpl close outfile(f);
```

Custom filtering via C API

Filtering files with custom code in very few lines:

mcpl_transfer_metadata does all the hard work of configuring output file

```
#include "mcpl.h"
void filter example()
  mcpl file t fi = mcpl open file("all.mcpl");
  mcpl_outfile_t fo = mcpl_create_outfile("lowEneutrons.mcpl");
  mcpl_transfer_metadata(fi, fo);
  mcpl hdr add comment(fo, "Only neutrons, ekin<0.1MeV");</pre>
  const mcpl particle t* prtcl;
  while ( ( prtcl = mcpl read(fi) ) ) {
    if ( prtcl->pdgcode == 2112 && prtcl->ekin < 0.1 )</pre>
      mcpl transfer last read particle(fi,fo);
  mcpl close outfile(fo);
  mcpl close file(fi);
```

mcpl_transfer_last_read_particle from MCPL v1.3.0 prevents lossy unpacking+repacking of data. If need to edit particles fields, replace with: mcpl_add_particle(fo,prtcl);

Python API (readonly)

To enable MCPL Python module, download mcpl.py or do python -mpip install mcpl (this incidently also installs the pymcpltool...)

Technical details:

- Pure Python, does not use mcpl.c
- Usage of Numpy for efficiency.
- Works with both Python 2 and 3.
- Readonly access for now.

```
import mcpl
myfile = mcpl.MCPLFile("myfile.mcpl")
for p in myfile.particles:
                                               Accessing particles is
   print( p.x, p.y, p.z, p.ekin )
                                               straight-forward
 Can also process blocks of
                              for p in myfile.particle blocks:
 N particles at a time, for
                                  print( p.x, p.y, p.z, p.ekin )
 increased efficency.
                                              Numpy arrays of length N
print( myfile.sourcename,
        myfile.nparticles,
         myfile.opt singleprec
                                               Can of course access
for cmt in myfile.comments:
                                               meta data as well.
     print( 'Comment: "%s"' % cmt )
                                                                   12 / 18
```

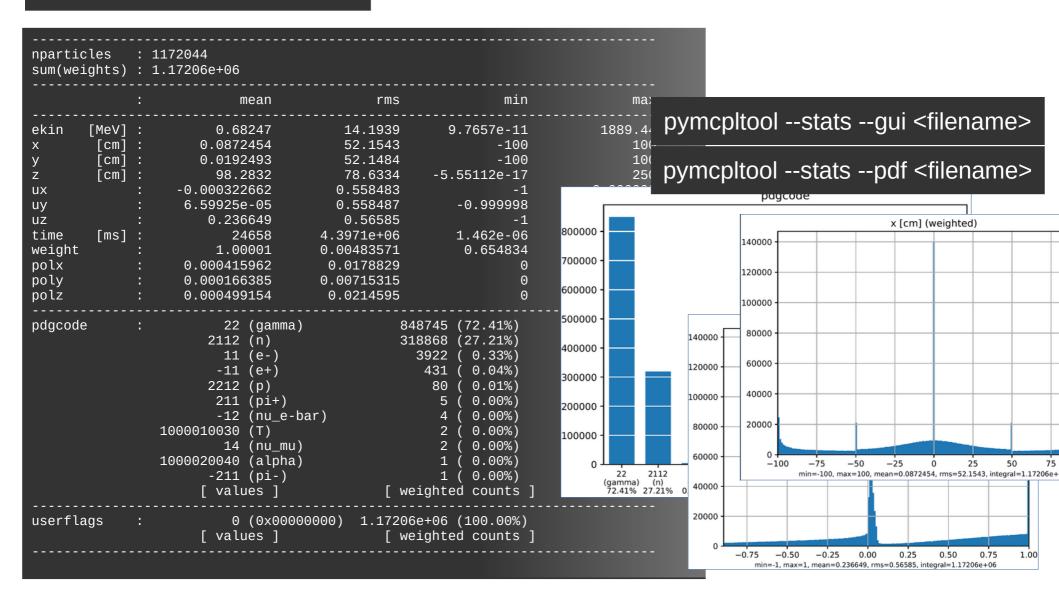
Command-line tools

- **mcpltool** and **pymcpltool**, both can:
 - **Inspect** files, extract binary blobs metadata to stdout
 - Convert MCPL to (inefficient) **ASCII** files for interoperability with software lacking MCPL support.
 - Show all options with --help
- The **mcpltool**:
 - Compiled executable with C compiler (from "fat" or proper linked code)
 - Can edit files:
 - **Merge** files
 - **Extract** subset of particles to smaller file (select by type or file idx)
 - **Repair** files leftover by crashed jobs
- The **pymcpltool**:
 - Built upon Python API (fast because of Numpy)
 - Download 1 file + run, or "pip install mcpl"
 - Can provide **statistics** (see next slide)



File statistics with pymcpltool

pymcpltool --stats <filename>



Merging files

- Ability to merge files is crucial for collecting output of concurrent simulations.
 - But other use-cases exists for combining files.
- Done via "mcpltool --merge" or "mcpl merge files(..)" in the C API.
- As a quality concern, MCPL is conservative about not producing files with misleading meta-data.
- All meta-data must be identical and will be transferred to the newly created file.
- On some occasions this restriction has caused problems...



Introduced "mcpltool --forcemerge" in release 1.3.0

- Can always merge, but will **throw away all meta-data**.
 - Should be considered as a last resort only!
- Particle data format options adapted to accommodate particles from all input files.
 - Options concerning FP prec., polarisation, fixed pdg/weight set as needed.
 - Discards userflags by default, since these are normally documented in metadata [override with --keepuserflags]
- Loss-less particle data transfer whenever possible.

```
Opened MCPL file forceme ged.mcpl:
  Basic info
                           CPL-3
    Format
    No. of particles
                            170823
    Header storage
                             bytes
    Data storage
                         : 79615964 bytes
  Custom meta data
    Source
                          "mcpl_forcemerge_f
                                              les (from MCPL v1.3.0)"
    Number of comments
     lumber of blobs
  Particle Lata format
    Priaris tion info
     ixed part. type
     ixed part. weight : no
     precision
                        : double
                         : little
    En lanness
                        : 68 byt s/particle
    Stora
                     ekin[MeV]
index
          pdgcode
                                      x[cm]
                                                  y[cm]
                                                                z[cm]
                      0.040287
                                     59.118
                                                  67.828
                                                                  250
                                                               197.92
                      0.048627
                                     19.774
                                                 -98.025
                                    -81.242
                      0.044083
                                                  58.308
                                                               71.342
                                     70.895
                      0.042855
                                                 -70.526
                                                               8.9938
               22
                                                               160.88
                          0.05
                                    -68.413
                                                 -72.936
               22
                      0.049592
                                    -95.998
                                                 -28.005
                                                               223.32
               22
                      0.042521
                                      84.72
                                                 -7.3153
                                                                  250
               22
                       0.04898
                                    -52.851
                                                 -84.892
                                                                26.98
               22
                      0.045358
                                     66.239
                                                 -74.916
                                                               127.78
                       0.04368
                                    -98.073
                                                  19.537
                                                               219.24
```

How to use MCPL in McStas

There's not much to it! Ships with components for **input**:

```
COMPONENT vin = MCPL input( filename="myfile.mcpl" )
AT(0,0,0) RELATIVE Origin
```

- Ignores non-neutrons.
- See mcdoc MCPL input for options controlling max energy of neutrons, or smearing of input particle properties.
- And the **output** component:

```
COMPONENT mcplout = MCPL output( filename="myoutput.mcpl" )
AT(0,0,0) RELATIVE PREVIOUS
```

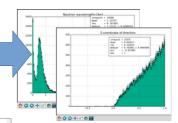
- See mcdoc MCPL output for options controlling e.g. floating point precision.
- Also possible for advanced users to set MCPL userflags.
- Good to be aware of a few design decisions by Peter & Erik in how McStas deals with MCPL files:
 - Will always use the full MCPL file (so ignores -n flag to mcrun). Use the repeat count option of MCPL_input allows replaying the file multiple times.
 - Concurrent modes (MPI/GPU) will replay the full MCPL file in each 7 / 18 process, possibly with smearing.

Outlook / wishful thinking

- Github issue 6: Mergeable statistics? E.g. "NEvtsSimulated" which would be added when files are merged. Would allow easier book-keeping.
- Github issue 44: In ESS Detector Group we have internal C++-based enhanced tools for working with MCPL files, based on our ExpressionParser and histogram classes:

mcplfilterfile in.mcpl.gz out.mcpl.gz "time<2ms and is neutron and neutron wl>2.2Aa"

mcplbrowse in.mcpl.gz where "pdgcode!=11 and ekin<10keV"





```
gen.input_file = "myfile.mcpl.gz"
gen.input_filter = "ekin>1keV && sqrt(x^2+y^2) < 10 cm"</pre>
```

- It would be great to export these tools to the greater community, but needs significant work to disentangle and prepare.
- IMHO if the Python API would not be read-only, we could easily build and easily distribute a lot of great new tools (e.g. GUI for editing). It would also be easy for people to compose/filter their own MCPL files from cmdline or code.
- Nice-to-have: An actual modern C++ interface with all the safety ant 18 convenience guarantees that entails.