HERACLES and DJANGO6: Updates for version 4.6.8 - 4.6.10

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Abstract

Modifications introduced into DJANGOH-4.6.8 to allow calculations for nuclear targets and for version 4.6.10: nucleon polarization.

1 File lhaglue-copy.f

- Parton distribution functions from LHAPDF required a modification in lhaglue.f. The modified version, lhaglue-copy.f is part of the DJANGOH source code now.
 - Removed reference to common block /ludat1/ because of type mismatch with JETSET 7.4; removed the corresponding assignment to unused mstu5(11).
 - To allow use of nuclear corrections from [3, 4, 5] via calls to structm from PYTHIA, structm in lhaglue-copy.f has been renamed to structf and structa into structm. The argument list of structa is shortened: the nuclear number a is transferred via the common block /HSNUCL/ from DJANGOH.

2 File djangoh_h.f:

• To allow linkage to the LHAPDF library [1], initialization of the path to the PDFsets folder has been introduced in subroutine HSWPDF and the path name has to be provided in the input file with code word LHAPATH:

LHAPATHI: Path name to the PDFsets folder of LHAPDF, character*80, no default defined.

The code was tested with LHAPDF version 5.8.6.

• The new code word NUCL-MOD with input INUMOD (integer, default = 0) has been introduced to select a model for nuclear shadowing or nuclear parton distribution functions.

The variable INUMOD extends the common block /HSNUCL/.

- INUMOD = 0: No correction. Parton distribution functions are used as selected by ILQMOD, ILIB, ICODE from code word STRUCTFUNC.
 This is the appropriate setting if nuclear parton distribution functions like those from [2] are used which incorporate nuclear effects.
- INUMOD = 1000: Structure functions are calculated from the isoscalar average of unmodified parton distribution functions and a Q^2 -independent correction factor. The same correction factor is applied to both F_2 and F_L .
- INUMOD = 2000: Apply correction factors to free parton distribution functions from EKS98 [3].
- INUMOD = 3000: Apply correction factors to free parton distribution functions from EPS08 [4].

INUMOD > 4000: Apply correction factors to free parton distribution functions from EPS09 [5].

LO or NLO fits (iorder = 1 or 2) and the error set ipset = $1, \ldots, 23$, are selected according to INUMOD = $100 \times iorder + ipset$.

The EPS09 analysis is based on CTEQ6.1M partons, but any PDF set can be combined with the nuclear corrections in this scheme.

Code in subroutines HSSTRF and HSSTR1 has been reorganized to allow for this option.

• Since version 4.6.10, the proton can be polarized [6]. The prescription for the calculation of the cross section is based on the parton model and distribution functions for polarized partons are used (see section 4, File polpdf.f). The following new input options are available:

HPOLAR: The degree of the proton polarization. Input is expected as a second double real in one line after the proton energy, i.e. with the code word PR-BEAM. If not between -1 and +1, HPOLAR is set to zero. This variable extends the common block HSPARM.

POLPDF: The new code word to read input for the selection of a parametrisation of polarized PDFs. Input for the integer variable IDPVR (on common block HSSTRP) is expected, where

IDPVR = 100*ISET+MODE with the following meaning:

ISET = 1: DSSV [7]; ISET = 2: DNS [8]; ISET = 3: DS [9]; ISET = 4: GSLO [10]; ISET = 5: GSNLO [10]; ISET = 6: BB [11]; ISET = 7: AAC [12]; ISET = 8: LSS [13]; ISET = 9: GRSV [14];

The new subroutine HSDPVR provides an interface to the necessary code which is collected in file polpdf.f (see below). A separate routine (HSCPVR) is called after integration and after sampling to print out the number of counts for which x and Q^2 were outside the range of validity of the parametrisation for polarized PDFs.

• The new code word OUTFILENAM allows to give a name for output files. The input is of type character*80 and has to start with a non-blank character. The output files are

- OUTFILENAM_out.dat is used for standard output. Redirection to the file starts after code word START appears in the input file;
- OUTFILENAM_smp.dat is used for sampling information needed for iterated calls;
- OUTFILENAM_rnd.dat takes information about the state of the random number generator;
- OUTFILENAM_his.paw is for histograms generated in the user routine contained in djangoh_u.f.

If no input to code word OUTFILENAM is given, standard output continues to appear at the console and the default name djangoh-default-output is used for the other output files.

Note that logical unit numbers 6, 8, 9, and 31 are used in the program (changed since version 4.6.10 to avoid a conflict with routines for polarized PDFs).

- If a negative value for input variable ISDINP (with code word RNDM-SEEDS) is given, the Cernlib random number generator ranlux is used with date and time as seed.
- A number of modifications was introduced to comply with gnu-fortran standards. The argument list of the subroutines DX1FCF and D01AJF was shortened and includes now used variables only

3 File djangoh_1.f:

• The target nucleon, proton or neutron, is chosen event by event according to its cross sections and using input from HNA and HNZ.

4 File polpdf.f:

• A collection of third-party code for polarized parton distribution functions. The most recent one (DSSV, [7]) is based on the most complete set of data and on theoretical calculations performed fully at next-to-leading order. The other programs were obtained from http://hepdata.cedar.ac.uk/pdfs.

The necessary grid-files are searched in a sub-directory called polpdf-gridfiles.

5 File gmc_random.f:

• Random number generator ranlux from Cernlib. The function PYR has been replaced by an interface to ranlux.

6 File djangoh_u.f:

• Filling of histograms is done with HFILL after a preceding call to HBARX. The present version in more accurate statistics with large event numbers. Output is written to file OUTFILENAM_evt.dat.

7 Usage

- Input to code word NUCLEUS is EPRO, the energy per nucleon, and HNA (nucleon number) and HNZ (proton number).
- Parton distribution functions are used from LHAPDF [1] if ILQMOD=0 and ILIB=2. Important values for the PDF set are

ICODE = 10150: CTEQ61M. Note that this set was used as a reference in the EKS and EPS fits of nuclear correction factors to the PDFs.

ICODE = 10550: CTEQ66.

ICODE = 29061: MRST98dis.

ICODE = 100151: HKNnlo for proton; in general: 100150+n with n=2 (deuteron), n=3 (Helium), n=6 (C), n=17 (Au), n=18 (Pb), etc.

• For version 4.6.10, a test run was performed with input file erhic-nc-test.in and erhic-cc-test.in. The corresponding output files are named as described above with OUTFILENAM = erhic-nc-test and OUTFILENAM = erhic-cc-test. Generation of 1,000,000 events (including F_L and fully fragmented) took about 14 minutes on eic0001.rcf.bnl.gov $(7.5 \times 10^{-4} \text{ sec per event plus initialization})$.

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