

<https://www.r-ccs.riken.jp/labs/cbrt/>

1.6.0

- Intel Parallel Studio XE 2018 Update 4
- CUDA 11.1 Update 1

- genesis-1.6.0.tar.bz2
- tests-1.6.0.tar.bz2
- fortdep.py.patch

```
--- fortdep.py.org    2020-12-28 14:05:33.000000000 +0900
+++ fortdep.py       2020-12-28 14:39:43.000000000 +0900
@@ -24,6 +24,7 @@
     re_fcomment = re.compile( "^(^!)*!(.*$)" )
     re_module   = re.compile( "^(.*/|s*)s*module\s*{([^\s,]*)s*", re.I )
     re_use      = re.compile( "^(.*/|s*)s*use\s*{([^\s,]*)s*", re.I )
+   re_include  = re.compile( "^(s*)#include\s*{([^\s,]*)s*" )
     mod_ext    = ".mod"

     def __init__( self, fname = "", ext = ".o" ):
@@ -31,6 +32,7 @@
         self.filename = fname
         self.modules  = []
         self.depmods  = []
+   self.includes = []

     def setFilename( self, fname, ext = ".o" ):
         if len(fname) == 0:
@@ -54,6 +56,14 @@
         if ( ma ):
             modname = ma.group(2).lower()
             self.appendDependMods( modname )
+   mi = FortranFile.re_include.search( line )
+   if ( mi ):
+       srcname = mi.group(2)
+       if re.search( r".h$", srcname, re.I ):
+           #self.includes.append( srcname )
+       pass
+   else:
+       self.includes.append( re.sub( r"\.[a-zA-Z0-9]+$", ".o", srcname ) )
         myf.close()
         self.makeUniqList()

@@ -64,8 +74,9 @@
         self.depmods.append( modname + FortranFile.mod_ext )

     def makeUniqList( self ):
-   self.modules = list(set(self.modules))
-   self.depmods = list(set(self.depmods))
+   self.modules = list(set(self.modules))
+   self.depmods = list(set(self.depmods))
+   self.includes = list(set(self.includes))

     def getMyModuleFileNames( self ):
```

```

return self.modules
@@ -87,7 +98,8 @@
    if m.lower() in mods_avail or not mods_avail:
        depmods.append(m)
    ret = ""
-   ret += self.objname + ": " + self.filename + " " + ".join(depmods) + " " + static_deps
+   ret += self.objname + ": " + self.filename + " " + ".join(depmods) + \
+       " " + ".join(self.includes) + " " + static_deps
    if len(self.modules) > 0:
        ret += "\n"
    ret += " ".join(self.modules) + ": " + self.filename + " " + self.objname

```

ビルド手順

```

#!/bin/sh

VERSION=1.6.0
BASEDIR=/home/users/${USER}/Software/GENESIS/${VERSION}
SRC_TARBALL=${BASEDIR}/genesis-${VERSION}.tar.bz2
TESTS_TARBALL=${BASEDIR}/tests-${VERSION}.tar.bz2
FORTDEP_PATCH=${BASEDIR}/fortdep.py.patch

INSTALLDIR=/local/apl/lx/genesis160-CUDA

WORKDIR=/work/users/${USER}
BUILDDIR=${WORKDIR}/genesis-${VERSION}
TESTSDIR=${WORKDIR}/tests-${VERSION}

PARALLEL=12
PARALLEL_TESTS=8

# -----
umask 0022

module purge
module load intel_parallelstudio/2018update4
module load cuda/11.1

export LANG=C
export LC_ALL=C
export OMP_NUM_THREADS=1

cd ${WORKDIR}
if [ -d genesis-${VERSION} ]; then
    mv genesis-${VERSION} genesis-erase
    rm -rf genesis-erase &
fi

if [ -d tests-${VERSION} ]; then
    mv tests-${VERSION} tests-erase
    rm -rf tests-erase &
fi

tar xjf ${SRC_TARBALL}
tar xjf ${TESTS_TARBALL}

cd ${BUILDDIR}
patch -p0 < ${FORTDEP_PATCH}
FC=mpiifort CC=mpiicc ./configure --prefix=${INSTALLDIR} \
    --enable-gpu \
    --enable-single \
    --with-cuda=/local/apl/lx/cuda-11.1
make depend

make -j ${PARALLEL} && make install

```

```
SPDYN=${INSTALLDIR}/bin/spdyn
```

```
cd ${TESTSDIR}/regression_test
```

```
# spdyn tests
```

```
./test.py      "mpirun -np ${PARALLEL_TESTS} $SPDYN" gpu
```

```
./test_remd.py "mpirun -np ${PARALLEL_TESTS} $SPDYN" gpu
```

```
./test_rpath.py "mpirun -np ${PARALLEL_TESTS} $SPDYN" gpu
```

```
./test_gamd.py "mpirun -np ${PARALLEL_TESTS} $SPDYN" gpu
```

```
./test_fep.py  "mpirun -np ${PARALLEL_TESTS} $SPDYN" gpu
```

メモ

- テストは全て通過(仕様上動作しない test_remd の 9 個のテストを除く)
- バイナリ(spdyn)は P100, V100 のどちらでも動作します(ccgpup, ccgpuv で確認)。ただし、ネイティブサポートにはなっていない可能性があります。