

[Credit: [John Swinbank & Dave Russell](#)]

These instructions have been tested on Debian 5.0.4 (lenny) on the AMD64 platform. This system was freshly installed from the [netinst CD image](#) . As part of the installation, I chose the "Standard system" and "Desktop environment" software collections. After installation, I applied all the outstanding system updates.

This instructions are intended to be read in conjunction with the documentation supplied with the various software packages!

Note that \$ indicates commands which should be typed at a terminal prompt as a normal user. # indicates commands which are typed at a root prompt (ie, with elevated privileges).

Preparation

Install the required packages from the Debian repository:

```
$ su
# aptitude update
# aptitude install make libc6-dev g++ gfortran x11proto-core-dev libx11-dev
libxt-dev libncurses5-dev libpng12-dev zlib1g-dev
```

Unfortunately, the version of slang supplied by the current Debian release is outdated & not compatible with the latest versions of Isis. However, it is straightforward to build and install an updated version using standard Debian tools:

```
# echo "deb-src http://mirrors.nl.kernel.org/debian/ testing main contrib non-free"
>> /etc/apt/sources.list
# aptitude update
# aptitude build-dep libslang2-dev
# apt-get -b source -t testing libslang2-dev
# dpkg -i libslang2_2.2.2-4_amd64.deb libslang2-dev_2.2.2-4_amd64.deb
```

```
libslang2-modules_2.2.2-4_amd64.deb slsh_2.2.2-4_amd64.deb
```

Note that the precise version numbers of the packages to install might vary between the time of writing and the summer school: you can always use `ls` to see exactly what has been built. After the above installation, you are free to remove the files that have been created in your working area.

You can now press `ctrl+d` to leave the root level shell and proceed with the installation.

SLIRP for S-Lang

Available directly from Debian:

```
$ su
# aptitude install slang-slirp
```

S-Lang Stats Module

For S-lang we require the [stats module](#) : download it from [here](#) (using your browser) or `wget`, then `untar` and `uncompress` it:

```
$ wget
http://space.mit.edu/cxc/software/slang/modules/stats/src
/slstats-1.0.2.tar.gz
$ tar zxvf slstats-1.0.2.tar.gz
```

Then configure and build it:

```
$ cd slstats-1.0.2/
$ ./configure
$ make
```

```
$ sudo make install
```

You will need to type your password.

This will install the module into `/usr/local`

. You now need to instruct S-Lang to look for scripts in that location.

Add the following lines to

```
$HOME/.slshrc
```

(creating the file if necessary):

```
append_to_slang_load_path  
("/usr/local/share/slsh/local-packages");  
set_import_module_path
```

```
("usr/local/lib/slang/v2/modules:"$+  
get_import_module_path ());
```

To use the module in a S-lang script (more info [here](#)), it must be loaded into the interpreter:

```
$ slsh  
slsh> require ("stats");
```

To quit slsh:

```
slsh:> quit
```

Heasoft 6.9

HEASoft includes FTOOLS (FITS

manipulation tools) and XANADU (X-ray data analysis).

Any problems with your Heasoft installation? First check for clues [h](#)
[ere](#)

then

[here](#)

Detailed installation guides are also available [here](#) and [here](#)

You'll need approximately up to 900 Mb free disk space

Download the source (300 Mb)

either using

[this direct link](#)

or by visiting

[the Heasoft website](#)

- Untar and uncompress it
in the directory you
want to unpack it in
:

```
$ tar zxvf  
heasoft6.9src.tar.gz
```

Then configure it.
You need to specify
the installation
location as part of
this process: we'll
use your home
directory in this
example.

NB if you don't use
your home directory,

you'll need to change
the other instructions
to match!

```
$ cd
```

```
heasoft-6.9/BUILD_  
DIR
```

```
$ ./configure
```

```
--prefix=$HOME/hea  
soft > config.out
```

```
2>&1
```

Wait until the configuration has finished, and build:

```
$ make > build.log
```

```
2>&1
```

Then install the

software:

\$ make install

At the start of your session, the heasoft environment must be initialised. We can make that happen automatically:

```
$ echo "export  
HEADAS=$HOME/  
heasoft/x86_64-unk  
nown-linux-gnu-libc  
2.7/" >>  
$HOME/.bashrc  
$ echo ".  
$HEADAS/headas-i  
nit.sh" >>
```

`$HOME/.bashrc`

Note that the platform string above (`x86_64-unknown-linux-gnu-libc2.7`) might vary if

Debian release
updated libc
packages. Check
what has been
created in
\$HOME/heasoft
to confirm the
correct value for

your system.

You will need to
open a new
terminal for this
to take effect.

Please do that,
and close the

old one, before
proceeding to
the next
section.

Now you are
ready to run

the various
Heasoft
software
packages:
Introduction,
tutorials (where
available) or

manuals: FTO

OLS

XANADU:

XSPEC

XRONOS

XIMAGE

Isis 1.6.0

ISIS

interactively
fits X-ray
spectra.
Any
problems

with your
ISIS
installation?
First check
for clues he

re

then (more
generally)

here

A detailed

installation
guide is
also
available [he](#)
[re](#)

Download
the source.
You can do
this directly
with the

command:

```
$ wget
```

```
ftp://space.
```

```
mit.edu/pu
```

b/cxc/isis/is

is-1.6.0-12.

tar.gz

or by

visiting the

ISIS

website

Untar and

uncompress
it in the
directory
you want
to unpack

it in :

\$ tar

zxvf

isis-1.6.0-

12.tar.gz

\$ cd

isis-1.6.0-

12/

Configure
and build:

\$

./configur

e

--with-hea

das=\$HE

ADAS

\$ make

There are
some
build in
diagnosti

c tests.

Before

proceedin

g, run

them and
check
that
everything

g is ok:

\$ make

check

If all the
tests
pass, you
can install

the
software
by
running:

\$ su

make

install

Press
ctrl+d to
leave the
root

environm
ent.

Now you

are ready

to run

ISIS: Tut

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Manual

Fermi Science Tools (v9r15p2

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in the
build

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for the
Fermi
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Tools
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platform.

However

, it is

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to

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and

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pre-com
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from the

Fermi
Science
Support
Centre

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the
version
compiled

for

Scientific

Linux 5

libc2.5,

and 32
or 64 bit
dependin
g on your

platform.

If you're
not sure,
check by

running:

\$

uname

-m

**A 64 bit
machine**

will

return x8

6_64

; a 32 bit

machine

will

return

i686



After you
have

identified

the

correct

package,

download
d with
your
browser

or with

wget:

\$ wget

<http://fermi.gsfc.nasa.gov/ssc/data/>

analysis/
software/
tar/Scien
ceTools-

v9r15p2-
fssc-200
90808-x8
6_64-unk

nOWN-lin

ux-gnu-li

bc2.5.tar.

gz

Note that
the instru
ctions
provided

by FSSC

for

installing

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binary
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on also
contain

some
errors. In
particular
, you

should

skip

steps 3 &

4.

Installati
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proceeds
as

follows:

\$ tar

zxvf

ScienceT
ools-v9r1
5p2-fssc-
2009080

8-x86_64
-unknow
n-linux-g
nu-libc2.

5.tar.gz

You may
wish to

move the
resulting
files to a
more

conveniently
memorable

location:

\$ mv

ScienceT

ools-v9r1

5p2-fssc-

2009080

8-x86_64

-unknow

n-linux-g

nu-libc2.

5

\$HOME/f

ermi_sci

ence_too

ls

Then set
the \$FE
RMI_DIR
environm

ent
variable
to point
to the

newly-ins

talled

tools:

\$ export

FERMI_

DIR=\$H

OME/fer

mi_scien
ce_tools/
x86_64-u
nknown-l

linux-gnu- libc2.5

You may

wish to
add this
to \$HOME
E/.bashr

c , so
that it
happens
automati

cally

when

you log

in:

\$ echo

"export

FERMI_

DIR=\$H

OME/fer
mi_scien
ce_tools/
x86_64-

unknown

-linux-gn

u-libc2.5

”

>>

\$HOME/

.bashrc

In order
to
access
the

science
tools,
you can
now run:

\$ sourc

e

\$ FERMI

_
DIR/fer

mi-init.s

h

Note

that

rather

than

integrati

ng with
the rest
of the
system,

the

Fermi

Science

Tools

include

an

additional

al (and

often

outdated

) copy of

a

number
of
standard
system

tools
(such as
Python,
tcl/tk,

etc). For
the
duration
of your

session
after you
source
the

fermi-init

.sh

script,

you will

automati
cally use
the
Fermi-s

applied

tools

rather

than

your
system
defaults.
Therefore

e, it is
strongly
suggested
to you

only
source
fermi-init
.sh

when

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lly need

to work
with
Fermi
tools,

rather
than as
part of
your

normal

login

process!

SAOIm

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DS9

6.1.2

DS9 is

a tool to

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and

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images.
It is
availabl

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downloa
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either

directly

using

your
browser
or using
wgget on

the
command
line.

Grab

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"Linux6

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version

if you
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e, or
"Linux"
for 32
bit. If

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check

using:

\$

uname

-m

"x86_64

"

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s 64 bit;
"i686"
32.

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ressed
and run
directly

without
any
further
configur

ation:

\$ wget

<http://he>

a-www.

harvard.

edu/sao

rd/down

load/ds

9/linux6

4/ds9.li

nux64.6

.1.tar.gz

\$ tar

zxvf

ds9.linu

x64.6.1.

tar.gz

\$./ds9

&

You
may
wish to
move it

to a
location
on your
path so

you can

access

it

without

specifying
the
full
location:

`$ sudo`

`mv ds9`

`/usr/loc`

`al/bin`

\$ ds9

&

AIPS

31 DEC

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AIPS

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CVS

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file (a Perl script)

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AIPS,
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\$

mkdir

aips

\$ cd

aips

\$

wget

ftp://ftp

.aoc.nr

ao.edu

**/pub/software/
aips/31**

DEC10

/install.

pl

\$

chmod

+X

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pl

\$ perl

install.

pl -n

Read &
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Refer

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d notes

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accept
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default

choice)

Screen

4b --

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enter

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Screen

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choose

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Screen

5a --

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"y"

Screen

5b --

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(LINUX6
4 on a

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Screen

6 -- hit

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LOCAL

HOST Screen 7 ..

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9 --

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"A" to

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setting

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Screen

9b --

press

"A" to

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11 --

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That's
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should

now
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latest
AIPS!
Enjoy.

**Edit
your \$
HOME/**

`.bashrc`
to
ensure

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AIPS is

initialis

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when

you log

in:

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~/.bash

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Now

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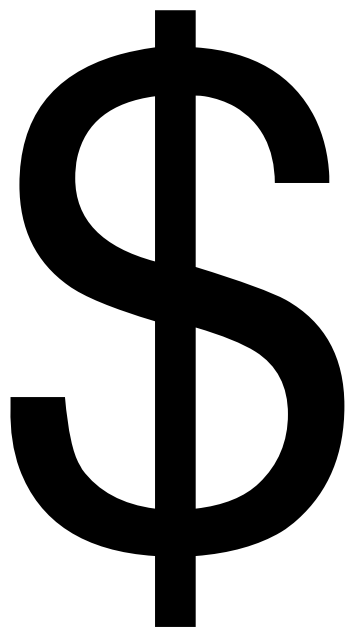
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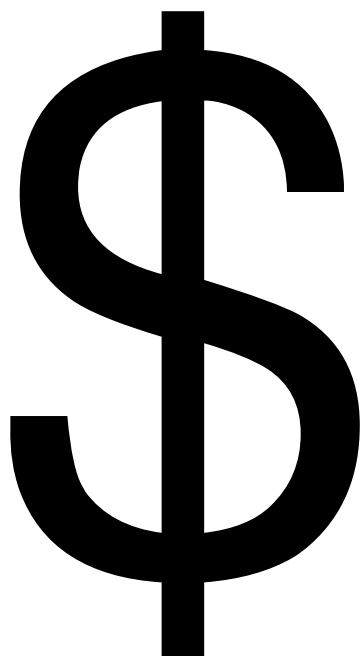
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