



Bugs: mkarf

Caveats

1. *The energy grid of the ARF and RMF files must be the same for use in XSpec. (25 Aug 2006)*

Sherpa allows you to use different energy grids for your ARF and RMF files, but *XSpec* does not. Note that *XSpec* will still run if the grids do not match, but it issues a warning and sets all values in the ARF to unity (1).

1. Create the RMF first

Since `mkacisrmf` can change the requested grid to match the calibration data, create the RMF first and then use it to define the energy grid when creating the ARF. This will work for both `mkarf` and `mkwarf`:

```
unix% pset mkarf \  
      engrid="grid(sources_ciao.wrmf[cols ENERG_LO,ENERG_HI])"  
  
or  
  
unix% pset mkwarf \  
      egridspec="grid(sources_ciao.wrmf[cols ENERG_LO,ENERG_HI])"
```

2. Match an existing ARF

If the `specextract`, `psextract` or `acisspec` scripts were used, you already have an ARF file for the data. Rather than remake both the RMF and ARF, get the grid information from the history in the ARF file:

```
unix% dmhistory acis_srcl.warf tool=all  
# dmhistory (CIAO 3.4): WARNING: Found "pixlib" library parameters  
  
# dmhistory (CIAO 3.4): WARNING: Found "ardlib" library parameters  
  
mkwarf infile="acis_srcl.[WMAP]" outfile="acis_srcl.warf" weightfile="acis_srcl.wfe  
spectrumfile="" egridspec="0.3:9.5:0.01" threshold="0" feffile="CALDB"  
mskfile="" mirror="HRMA" detsubsysmod="" ardlibpar="ardlib" geompar="geom"  
clobber="no" verbose="2"
```

Your file may have been created with `mkarf` instead of `mkwarf`; the `dmhistory tool=all` will show the tool used in either case.

Use the `egridspec` value (or `engrid` in the `mkarf` case) as input for the energy parameter in `mkacisrmf`:

```
unix% pset mkacisrmf energy="0.3:9.5:0.01"
```

The [mkacisrmf analysis thread](#) has information on creating the RMF file.

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URL:
<http://cxc.harvard.edu/ciao3.4/bugs/mkarf.html>
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