

*“AWK is a language for processing files of text. A file is treated as a sequence of records, and by default each line is a record. Each line is broken up into a sequence of fields, so we can think of the first word in a line as the first field, the second word as the second field, and so on. An AWK program is of a sequence of pattern-action statements. AWK reads the input a line at a time. A line is scanned for each pattern in the program, and for each pattern that matches, the associated action is executed.” – Alfred V. Aho*

The design of `nawk` was influenced by shell scripting languages that we already used. In turn `nawk` influenced the development of more advanced and complex languages such as `perl`. It shares much syntax with `C`, but `nawk` is not a compiled language to be used on large data files. It is, however, extremely powerful on short or text-based files for advanced searches or data extraction. It can be executed from a command-line or from within a shell script as

```
% nawk 'pattern {action}' infile >outfile
```

The “pattern” is basically a statement, which, if true, will enable the “action.” A set of one-line programmes, variations of which I use a lot, are:

```
% nawk 'END{print NR}' infile          #NR is a line number, printed after all of infile is read
% nawk '$0 ~ "string" {print $0}' infile # if a line matches "string", print that line
% nawk '{print NR, NF}' infile          # for each line, print line number and number of column
% nawk '$5 == 4 {print sqrt($5)}' infile # if column 5 equals 4, print the square root of 4
```

The program content ‘pattern {action}’ can also be placed within a file which can be executed either from a command line or from within a shell script as

```
% nawk -f file.awk par=set_pa infile >outfile
```

where the passing of a variable called `par` with a value `set_pa` is optional as is the direction of the output to `outfile`. Check the Wiki page above (and/or google `awk`) for more details.

You are here tasked to write an `awk`-file called `RadianScales.awk` (`ReflecScales.awk`) to extract the scaling factors that convert the MODIS science integer values to radiance (reflectance) values. For this purpose you will make extensive use of the file `HeaderInfo.ascii`:

14.	Extract relevant information on calibration, file size, scan numbers, etc, from a header dump, e.g. <code>hdfdump -h \$name.L1B_LAC &gt;HeaderInfo.ascii</code>
15.	How many lines does the <code>HeaderInfo.ascii</code> file contain? Are the number of column per line constant?
16.	Search the <code>HeaderInfo.ascii</code> file from the command line for any occurrence of a text string “radiance”: <code>nawk '\$0 ~ "string" {print \$0}' HeaderInfo.ascii</code>
17.	You will get too many lines, so your string “pattern” was too general, modify it to, say, “radiance_scales”
18.	The output may still be too cryptic, so search also for the text string “band_names” and see if you can make sense of the different radiance scales. How many bands are there? What’s special about band-26?
19.	Extract the <code>radiance_scales</code> for the bands 1 and 2 by designing a unique “pattern” that returns only one line.
20.	Note that the scales contain the letter “f” at the end, print a sub-string without it; the “action” <code>print substr(\$x,StartCharacter,NumberOfCharacters)</code> will be useful, here “x”, “StartCharacter”, “NumberOfCharacters” are integers you must specify.
21.	Make file <code>RadianScales.awk</code> that contains the above and add the following line to your <code>proc2.csh</code> shell script: <code>nawk -f RadianScales.awk HeaderInfo.ascii &gt;log.dat</code> where you replace the ‘pattern {action}’ construct with <code>{ if (pattern);{action};endif }</code>