

Source file: example

```
#####
#
# A simple example of a Maple 'source' file. Such a file can
# be easily created and maintained using your favorite text
# editor and can contain arbitrary Maple commands.
# I find this mechanism particularly useful for developing
# and maintaining Maple procedures.
#
# The file can most easily be read into a Maple session by
# first 'cd'-ing to the directory which contains this file:
#
# % cd /Public/Members/matt/Src/maple/examples
#
# starting up maple or xmaple
#
# % maple
#
# or
#
# % xmaple &
#
# then typing
#
# > read example;
#
# If maple (or xmaple) isn't running in the directory
# containing this file, then you must use an absolute
# pathname and be sure to enclose the name in backquotes
# or double-quotes. (This also applies to filenames
# containing a '.', which is why I tend to use simple
# names (no extensions) for files containing Maple source.)
#
# > read '/Public/Members/matt/Src/maple/examples/example';
#
# Recall that use of the colon (:) as terminator rather
# than semi-colon (;) inhibits echoing of results.
#
#####
aa := 23 / 155;

myadd := proc(x::numeric, y::numeric)
    x + y;
end;
```

Source file: ladd

```
#####
# ladd: Adds all elements of a list.
#####
ladd := proc(l::list)

#-----
# Define local variables.
#-----
local lsum, i;
#-----
# Check for valid argument, exit with error message
# if not valid.
#-----
if nops(l) = 0 then
    ERROR('argument is the NULL list');
fi;
#-----
# Initialize sum to first element of list.
#-----
lsum := l[1];
#-----
# Loop over rest of elements accumulating the sum.
#-----
for i from 2 to nops(l) do
    lsum := lsum + l[i];
od;
#-----
# Return the sum.
#-----
lsum;
end:
```

```
#####
# ladd: Alternative, more compact implementation using
# 'add' procedure. Not possible before Maple V.4.
#####
laddnew := proc(l::list)
    local i;
    add( l[i], i=1..nops(l) );
end:
```

Source file: tladd

```
#####
# Tests
#      ladd
#      laddnew
#####
read ladd;

l1 := [1,2,3,4];

#####
# Note the use of the 'printf' procedure, and the '%a'
# format specification.
#####
printf("ladd(...) is %a \n", ladd(l1));
printf("laddnew(...) is %a \n", laddnew(l1));
```

Source file: tladd-trace

Script started on Mon Sep 15 14:27:25 2003

```
lnx1% cat tladd
#####
# Tests
#      ladd
#      laddnew
#####
read ladd;

l1 := [1,2,3,4];

#####
# Note the use of the 'printf' procedure, and the '%a'
# format specification.
#####
printf("ladd(...) is %a \n", ladd(l1));
printf("laddnew(...) is %a \n", laddnew(l1));
```

```
lnx1% maple
| \^/|      Maple 6 (IBM INTEL LINUX)
._\|_|_ /|/_ . Copyright (c) 2000 by Waterloo Maple Inc.
\ MAPLE / All rights reserved. Maple is a registered trademark of
<---- ----> Waterloo Maple Inc.
|      Type ? for help.
> read tladd;
               l1 := [1, 2, 3, 4]

ladd(...) is 10
laddnew(...) is 10
> quit
bytes used=112428, alloc=131048, time=0.01
lnx1% exit
exit
```

Script done on Mon Sep 15 14:27:46 2003

Source file: polyinterp

```
#####
#
# polyinterp: Constructs Lagrange Interpolating Polynomial
#
# Given n distinct "data points" (x_i,f_i) , i = 1 ... n, and a name,
# this procedure returns the unique polynomial (in name) of degree
# n - 1 which passes through (interpolates) all the points.
#
# Input parameters:
#
#   ldata:    list of lists, which defines (x_i,f_i)
#   var:      name, returned interpolating polynomial is
#             a polynomial in 'var'
#
# Usage example:
#
#   > polyinterp([ [0,1], [1,6], [2,4], [3,0] ], 'x' );
#
#               3      2
#               5/6 x  - 6 x  + 61/6 x + 1
#
# Implementation notes:
#
#   This routine converts the list of input pairs (each pair
#   itself a two-element list) to separate *sequences* of
#   the x_i and f_i. You could also build up separate *lists*
#   but it is syntactically easier to build sequences in Maple.
#
#####
polyinterp := proc(ldata::list(list),var::name)
#-----
# Local variables:
#
#   n:      number of data points
#   i, j:   loop variables used in evaluation of Lagrange formula
#   sx, sf: for building up sequences of x_i, f_i respectively
#   num, den: for building up the numerators and denominators of the
#             characteristic polynomials.
#   p:      for building up the interpolating polynomial itself
#
#-----
local n, i, j, sx, sf, num, den, p;
#
#-----
# Determine number of data points
#
#-----  
n := nops(ldata);
```

```

#-----
#  Initialize polynomial and x_i and f_i sequences
#-----
p := 0;
sx := NULL;
sf := NULL;
#-----
# Convert input list-of-lists into separate sequences of x_i and f_i
#-----
for i from 1 to n do;
    sx := sx , ldata[i][1];
    sf := sf , ldata[i][2];
od;

#-----
# For each of the x_i ...
#-----
for i from 1 to n do;
#-----
#     ... build up the numerators and denominators of the ith
#     characteristic polynomial. First initialize the numerator
#     and denominator ...
#-----
num := 1;
den := 1;
#-----
#     ... and then build them up using the Lagrange formula. Note that
#     both the numerator and denominator are products of n - 1
#     terms, one term for each j = 1..n such that j <> i.
#-----
for j from 1 to n do;
    if j <> i then
        num := num * (var - sx[j]);
        den := den * (sx[i] - sx[j]);
    fi
od;
#-----
#     Update the polynomial
#-----
p := p + sf[i] * (num / den);
od;
#-----
# Return the polynomial in expanded form
#-----
expand(p);

end:

```