

AMPL

A Modeling Language for Mathematical Programming

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Introduction

- ▶ Algebraic **modeling language** for linear and nonlinear optimization problems, in discrete or continuous variables.
- ▶ Allow the **development of models and algorithms**.
- ▶ **Linked** to the most widely used **solvers** for LP/MILP/MINLP programming.
- ▶ **Download:**
`https://www.lix.polytechnique.fr/~dambrosio/teaching/`
- ▶ **AMPL book:**
`www.ampl.com/BOOK/download.html`
- ▶ **Quick-start guide:**
`https://www.lix.polytechnique.fr/~dambrosio/teaching/ampl-quick-start-guide_dambrosio.pdf`

AMPL files

- ▶ Each problem instance is coded in AMPL using three files:
 - ▶ a **model file** (extension .mod): contains the mathematical formulation of the problem.
 - ▶ a **data file** (extension .dat): contains the numerical values of the problem parameters.
 - ▶ a **run file** (extension .run): specifies the solution algorithm (external and/or coded by the user in the AMPL language itself).

File .mod

- ▶ parameters, lines starting with the keyword
`param`
- ▶ sets, lines starting with the keyword
`set`
- ▶ decision variables, lines starting with the keyword
`var`
- ▶ objective function(s), lines starting with the keyword
`minimize` or `maximize`
- ▶ constraints, lines starting with the keyword
`subject to`

File .mod

```
param n > 0;
```

```
param w{1..n} > 0;
```

```
param n > 0;  
param m > 0;  
param a{1..n, 1..m};
```

File .mod

```
set N := 1..n;
```

```
param w{N} > 0;
```

```
param w{N} > 0;  
param p{j in N} <= 10*w[j];
```

File .mod

Decision variables:

```
var x{j in 1..n} >= 0, <= 1, binary;
```

Objective function:

```
maximize total_profit:  
    sum{j in N} p[j]*x[j];
```

File .mod

```
subject to capacity_constraint:  
    sum{j in N} w[j]*x[j] <= c;
```

```
subject to random_constraint{j in 2..n}:  
    w[j]*x[j] - w[j-1]*x[j-1] <= 1;
```


Non linear knapsack problem: File .mod

```
param N > 0;           # number of objects
set VARS ordered := {1..N};
param U {j in VARS} > 0, default 100;
param a {j in VARS} > 0;
param b {j in VARS} > 0;
param c {j in VARS} > 0;
param d {j in VARS} < 0;
param C > 0; # knapsack capacity

var x {j in VARS} >= 0, <= U[j]; # variables

maximize Total_Profit: # objective function
  sum {j in VARS} (a[j]+b[j]*x[j]+c[j]*x[j]**2+d
    [j]*x[j]**3);

subject to KP_constraint: # constraint
  sum{j in VARS} x[j] <= C;
```

Non linear knapsack problem: File .dat

```
param N := 10;  
param C := 546.000000;
```

```
param: a :=  
1      0.172274  
2      0.134944  
3      0.101030  
4      0.163588  
5      0.152350  
6      0.196601  
7      0.181208  
8      0.126588  
9      0.184087  
10     0.187434  
;
```

```
param: b :=  
1      78.770199  
2      77.468892  
3      93.324757  
4      96.180080  
5      55.137398  
6      40.101851  
7      36.007819  
8      5.317250  
9      9.964929  
10     60.265707  
;
```

```
param: c :=  
1      3.062328  
2      43.280130  
3      52.983122  
4      62.101010  
5      58.531125  
6      47.574366  
7      53.101406  
8      6.902601  
9      16.985577  
10     62.576610  
;
```

```
param: d :=  
1      -81.876165  
2      -56.455229  
3      -56.428945  
4      -43.813029  
5      -28.246895  
6      -81.108142  
7      -37.839956  
8      -1.138258  
9      -80.968161  
10     -11.536015  
;
```

```
param: U :=  
1      100.000000  
2      100.000000  
3      100.000000  
4      100.000000  
5      100.000000  
6      100.000000  
7      100.000000  
8      100.000000  
9      100.000000  
10     100.000000  
;
```

Non linear knapsack problem: File .run

```
# Author: Claudia D'Ambrosio
# Date: 20190121
# nlkp.run

reset;
reset data;

model nlkp.mod;

data "/mypath/nlkp.dat";

option solver "/usr/local/bin/baron";

option baron_options 'prfreq=100 outlev=1';

solve > nlkp.out;
```

Other commands

```
reset;  
reset data;
```

```
option solver gurobi;  
option gurobi_options "outlev 1";  
solve;
```

Other commands

```
model myMILPmodel.mod;
data myInstance.dat;
option solver cplex;
option relax_integrality 1; # relaxing the
    integrality requirements on all the decision
    variables
solve;
option relax_integrality 0; # restoring the
    integrality requirements on all the decision
    variables
solve;
```

Other commands

```
display n, c;  
display N;  
display w, p;
```

```
n = 7  
c = 19  
  
set N := 1 2 3 4 5 6 7;
```

```
:      w      p      :=  
1      11     10  
2       6      3  
3       6      4  
4       5      5  
5       5      6  
6       4      7  
7       1      2  
:
```

Other commands

```
display x;  
display cost;
```

```
x [*] :=  
1  0.363636  
2  0  
3  0  
4  1  
5  1  
6  1  
7  1  
;  
  
cost = 23.6364
```

Other commands

```
display capacity_constraint;
```

```
capacity_constraint = 0.909091
```


Other commands

```
expand capacity_constraint;
```

```
subject to capacity_constraint:  
    11*x[1] + 6*x[2] + 6*x[3] + 5*x[4] + 5*x[5] + 4*x[6] + x[7] <= 19;
```

Other commands

```
printf "param n := %d;\n", n;
printf "\n";

# param c
printf "param c :";
for {j in N} {
  printf "\t%d", j;
}
printf "\t:=\n";
for {i in N} {
  printf "\t%d", i;
  for {j in N} {
    printf "\t%d", c[i,j];
  }
  printf "\n";
}
printf ";\n\n";
```

Other commands

```
for {j in 1..n} {  
  ...  
}
```

```
repeat {  
  . . .  
}  
until x[n] > 0;
```

How to call AMPL from the command line:

```
Claudias-MacBook-Pro-8: ampl myrunfile.run
```

or

```
Claudias-MacBook-Pro-8: ampl myrunfile.run > myoutputfile.out
```

How to call AMPL from the AMPL environment/IDE:

```
AMPL: include "../myfolder/myrunfile.run";
```

Non linear knapsack problem: File nlkp.out obtained

```
BARON 18.11.12 (2018.11.12): prfreq=100 outlev=1
=====
Preprocessing found feasible solution with value 1.60010400000
Doing local search
Preprocessing found feasible solution with value 1093.96317285
Solving bounding LP
Starting multi-start local search
Done with local search
=====
Iteration      Open nodes      Time (s)      Lower bound      Upper bound
-----
1              1                0.15         1093.96         11340.2
100            16               1.62         1093.96         1121.50
200            30               2.04         1093.96         1094.05
300            35               2.19         1093.96         1094.00
400            31               2.38         1093.96         1093.97
500            28               2.55         1093.96         1093.97
600            13               2.67         1093.96         1093.97
631            0                2.70         1093.96         1093.96
Cleaning up
*** Normal completion ***
Wall clock time:          3.00
Total CPU time used:     2.70

Total no. of BaR iterations:    631
Best solution found at node:    -1
Max. no. of nodes in memory:   37

All done
=====
BARON 18.11.12 (2018.11.12): 631 iterations, optimal within tolerances.
Objective 1093.963173
Profit = 1093.96
```

To install AMPL

- ▶ Download one of the following .zip files:
 - ▶ http://www.lix.polytechnique.fr/~dambrosio/teaching/ampl_linux-intel64.tgz
 - ▶ http://www.lix.polytechnique.fr/~dambrosio/teaching/ampl_macos64.tgz
 - ▶ http://www.lix.polytechnique.fr/~dambrosio/teaching/ampl_mswin64.zip
- ▶ Follow installation instructions:
<https://ampl.com/ampl-course-install/>
- ▶ Available solvers: baron, conopt, cplex, gurobi, ilogcp, knitro, lgo, loqo, minos, snopt, xpress

References

- ▶ Modeling languages like **ampl**: ampl.com
or **gams**: www.gams.com or **jump**
<https://jump.dev/JuMP.jl/>
- ▶ Open source solvers like **scip**: scip.zib.de
- ▶ **NEOS Server**, State-of-the-Art Solvers for Numerical Optimization: www.neos-server.org/neos/