# INF580 - Advanced Mathematical Programming <br> TD3 - Complexity and MP 

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## Simple AMPL codes

Write AMPL code for the following problems:

- Max Clique
- SUBSET-SUM
- KNAPSACK
- Hamiltonian Cycle and test them with the feasible and infeasible (whenever applicable) instances given in the course slides


## Random instance generators

- Coding up instances by hand is boring
- Let's use AMPL to generate random instances!
- Each problem needs its own generator
- In general, for a problem called prob:

1. copy the index sets / parameters from prob.mod to prob-instgen.run
2. set sizes by hand (e.g. "let $\mathrm{n}:=5$;")
3. use AMPL imperative sublanguage to randomly fill set/param values
4. print to file

## Random instance generators

## Example: generating random graphs, Erdős-Renyi model

```
option randseed 0; # pseudornd gen starts from rnd seed
## start from sizes/index sets/params of original problem
param n integer, > 0;
set V := 1..n;
set E within {V,V};
## randomly generate missing index sets/params
let n := 50; # initialize number of vertices
param p := 0.5; # probability of creating edge
let E := {}; # initialize the edge set to empty
for {i in V, j in V : i < j} { # no loops or antiparallel arcs
    if Uniform(0,1) < p then {
        let E := E union {(i,j)}; # create the edge
    }
}
## print out a .dat file (MIND YOU DON'T OVERWRITE OLD .dat FILES!)
print "# file generated by clique-instgen.run" > rndcliq.dat;
printf "param n := %d;\n", n >> rndcliq.dat;
printf "set E :=" >> rndcliq.dat;
for {(i,j) in E} {
    printf " (%d,%d)", i,j >> rndcliq.dat;
}
printf ";\n" >> rndcliq.dat;
```


## Random instance generators

## Example: uniformly distributed knapsack instances

```
option randseed 0; # pseudornd gen starts from rnd seed
param n integer, > 0;
set N := 1..n;
param c{N} integer;
param w{N} integer;
param K integer, >= 0;
## randomly generate missing index sets/params
let n := 20; # initialize number of objects
param cL := 1; param cU := 10; # bounds for object volume
param wL := 1; param wU := 10; # bounds for object value
let {i in N} c[i] := round(Uniform(cL,cU));
let {i in N} w[i] := round(Uniform(wL,wU));
let K := round((sum{i in N} c[i])/2); # generate capacity
## print out a .dat file (MIND YOU DON'T OVERWRITE OLD .dat FILES!)
print "# file generated by knapsack-instgen.run" > rndknap.dat;
printf "param n := %d;\n", n >> rndknap.dat;
printf "param K := %d;\n", K >> rndknap.dat;
printf "param : c w :=\n" >> rndknap.dat;
for {i in N} {
    printf " %i %d %d\n", i, c[i], w[i] >> rndknap.dat;
}
printf ";\n" >> rndknap.dat;
```


## Random instance generators

```
Example: normally distributed knapsack instances Change
```

```
param cL := 1; param cU := 10; # bounds for object volume
```

param cL := 1; param cU := 10; \# bounds for object volume
param wL := 1; param wU := 10; \# bounds for object value
let {i in N} c[i] := round(Uniform(cL,cU));
let {i in N} w[i] := round(Uniform(wL,wU));
to
param cavg := 4.5; param cstdev := 1.9; \# object volume
param wavg := 5.8; param wstdev := 3.1; \# object value
let {i in N} c[i] := round(Normal(cavg,cstdev));
let {i in N} w[i] := round(Normal(wavg,wstdev));

```

\section*{Playing with instances}
- How many vertices/edges does the largest Max Clique instance have, that CPLEX can solve in 30s on your laptop?
- Generate 9 random graphs, each with 160 vertices, and with edge generation probability \(p \in P=\{0.1,0.2, \ldots, 0.9\}\)
- Find max cliques on all these graphs
- use CPLEX as a solver
- use "option cplex_options "mipdisplay=2";" after "option solver cplex;" (shows CPLEX progress)
- record size \(\omega\left(G_{p}\right)\) of max clique of each graph \(G_{p}\) and CPU time \(\gamma_{p}\) for \(p \in P\)
```

with bash: name random instances rndcliq-0.1.dat, ..., then type (1 line):

```
for i in 0.10 .20 .30 .40 .50 .60 .70 .80 .9 ; do
cp rndcliq-\$i.dat rndcliq.dat ; ampl clique.run ;
done > clique.log 2>\&1 \&
- plot \(\omega\left(G_{p}\right)\) versus \(p\) and \(\gamma_{p}\) versus \(p\) with bash: grep OUT: clique.log | cut -d ':' -f 2

\section*{The Motzkin-Straus formulation}

Write AMPL code to implement the Motzkin-Straus formulation for solving Max Clique
- make sure this formulation can read the same . dat files as those you already worked on
- test this formulation on the instance given in the course slides
- use a global optimization solver (e.g. baron) and also a local optimization one (e.g. snopt): what results do you obtain?
- can this formulation be solved using cplex?
- what is the maximum instance size you can solve to global optimality with this formulation? What about local optimality?

\section*{Structured formulation for SAT}
- Propose a numerical encoding for SAT instances
- Based on this, write a structured MP formulation for SAT
- Implement it in AMPL and test it using an appropriate solver```

