

# Multiple Nested Structures: the Curse (or Blessing?) of Applied Mathematics

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- 1 Structure: top-down
- 2 Structure: bottom-up
- 3 The Core Elements of SMS++
- 4 The quasi-Core Elements of SMS++
- 5 Example: SMS++ for the Unit Commitment
- 6 Conclusions and (a Lot of) Future Work

- General optimization problem  $(P) \min \{ c(x) : x \in X \}$ :  
clearly unsolvable if  $c(\cdot)$  and  $X$  are “any” function/set
- To do anything one needs assumptions on the structure of  $c(\cdot)/X$
- Many different cases, most of them hard
- Let's take it easy: strong structure  $\equiv$  easy problem:

$$\text{Linear Program } (P) \min \{ cx : Ax = b, 0 \leq x \leq u \}$$

$$A \in \mathbb{R}^{n \times m} \text{ (sparse)}, b \in \mathbb{R}^n, c \in \mathbb{R}^m, u \in \mathbb{R}^m, m > n$$

- Structure  $\implies$  useful properties: dual problem to  $(P)$

$$(D) \max \{ yb - wu : yA + z - w = c, z \geq 0, w \geq 0 \}$$

fundamental tool for solving  $(P)$

- **Karush-Kuhn-Tucker** optimality conditions ( $\text{diag}(V) = v$ ,  $e = \text{all } 1\text{s}$ )

$$(KKT) \quad \begin{cases} Ax = b, \quad x + s = u, \quad yA + z - w = c & (\text{linear}) \\ XZe = 0, \quad SWe = 0 & (\text{nonlinear}) \\ [x, s, z, w] \geq 0 & (\text{inequalities}) \end{cases}$$

- **Interior Point** methods for LP: “slacken + linearize”:

i)  $\mu > 0$ ,  $(KKT_\mu) \equiv (KKT)$  except  $XZe = \mu e$ ,  $SWe = \mu e$

$\implies (2m\mu)$ -optimal solution

ii) **feasible**  $[\bar{x}, \bar{s}, \bar{z}, \bar{w}] > 0$ ,  $v = \bar{v} + \Delta v$  (stepsize ensures  $v > 0$ )  $\implies$

$$\begin{cases} A\Delta x = 0, \quad \Delta x + \Delta s = 0, \quad \Delta yA + \Delta z - \Delta w = 0 \\ \bar{X}\bar{Z}e + \bar{X}\Delta z + \bar{Z}\Delta x = \mu e, \quad \bar{S}\bar{W}e + \bar{S}\Delta w + \bar{W}\Delta s = \mu e \end{cases}$$

ignore second-order terms  $\equiv$  **Newton's method** for nonlinear equations

- $[\bar{x}, \bar{s}, \bar{z}, \bar{w}]$  satisfies  $(KKT_\mu)$ : one iteration,  $\mu \searrow$  (fast), repeat
- Many improvements (infeasible method, predictor corrector, ...)

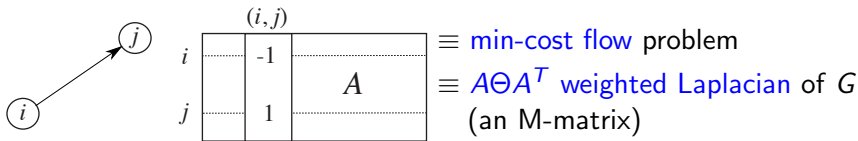
- Boils down to **Reduced KKT** or **Normal equations** ( $\Theta > 0$  diagonal)

$$\begin{bmatrix} -\Theta & A^T \\ A & 0 \end{bmatrix} \quad A\Theta A^T$$

$m + n \times m + n$ , sparse       $n \times n$ , a lot less sparse

- Special case of **Saddle-Point system**, lots of applications (physics, engineering, economy, computer science, ...), very active research<sup>1</sup>
- Specific twists in the LP case:
  - large size**:  $m \approx 10^{6+}$ ,  $n \approx 10^{5+}$  ...
  - must be solved **many times**, but **rather inexactly** (at the first iterations)
  - fixed nonzero structure** ( $A$ ) and **variable data** ( $\Theta$ )
  - special **evolution of data** over the iterations
  - no discretization, no underlying smooth operator
- Ultimate performances** require **assumptions** on (structure of)  $A$

- $A =$  *node-arc incidence matrix* of directed graph  $G$



- Can tell a lot on the system by looking at the graph<sup>2,3</sup>
- Can do a lot about the system by working on the graph:
  - preconditioners are (chordal) sub-graphs, can be obtained by efficient graph algorithms (Kruskal<sup>4</sup>, Prim<sup>5</sup>, ...)
  - projection in algebraic multigrid is merging nodes<sup>6</sup>
  - projection and preconditioning is a unique graph-based process<sup>7</sup>

<sup>2</sup> Cvetković, Doob, Sachs "Spectra of graphs", 1980 — Brouwer, Haemers "Spectra of Graphs", 2012

<sup>3</sup> F., Serra Capizzano "Spectral Analysis of (Sequences of) Graph Matrices" *SIMAX*, 2001

<sup>4</sup> F., Gentile "New Preconditioners for KKT Systems of Network Flow Problems" *SIOPT*, 2004

<sup>5</sup> F., Gentile "Prim-based Support-Graph preconditioners for Min-Cost Flow Problems" *CO&A*, 2006

<sup>6</sup> Dell'Acqua, F., Serra Capizzano "Computational Evaluation of Multi-Iterative Approaches [...]" *CALCOLO*, 2015

<sup>7</sup> Dell'Acqua, F., Serra Capizzano "Accelerated Multigrid for Graph Laplacian Operators" *Appl. Math. & Comp.*, 2015

- These systems have been approached by many different angles:
  - graph theory
  - computer science
  - numerical linear algebra
  - optimization
  - physics ...
- Lots of ingenuity, theoretical results, implementations
- Applied mathematics at its best: **focus on one structure with relevant applications, drill it down until it cries**

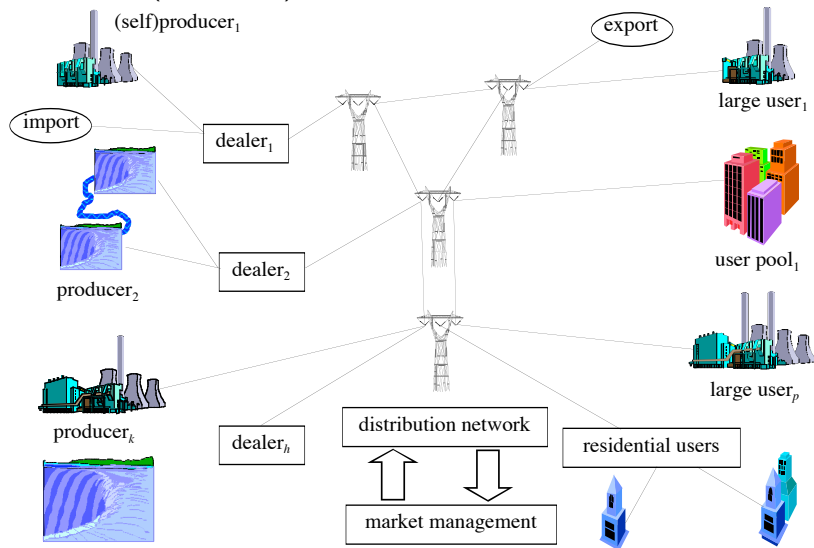
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- Applied mathematics at its best: **focus on one structure with relevant applications, drill it down until it cries**
- **Is this always enough?**



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# Motivation: The Unit Commitment (UC) problem

- Schedule a set of **generating units**  $\mathcal{U}$  over a **set of time instants**  $\mathcal{T}$  to satisfy the (forecasted) **demand**  $d_t$  at each  $t \in \mathcal{T}$

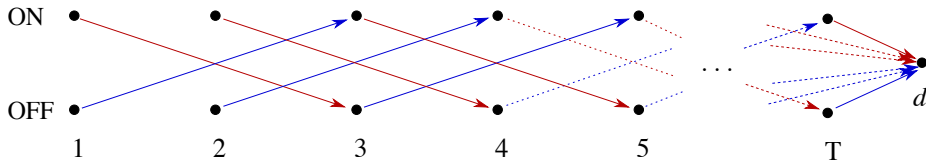


- Gazzillions €€€ / \$\$\$, enormous amount of research<sup>8,9</sup>
- **What has it to do with networks?** More than it would seem
- Different types of production units, different constraints:
  - Thermal (comprised nuclear): min/max production, min up/down time, ramp rates on production increase/decrease, start-up cost depending on previous downtime, others (modulation, ...)
  - Hydro (valleys): min/max production, min/max reservoir volume, time delay to get to the downstream reservoir, others (pumping, ...)
  - **Non programmable** (ROR hydro) **intermittent** units (solar/wind, ...)
  - Fancy things (small-scale storage, demand response, smart grids, ...)
- Plus the **interconnection network** (AC/DC, transmission/distribution)

<sup>8</sup> van Ackooij, Danti Lopez, F., Lacalandra, Tahanan "Large-scale Unit Commitment Under Uncertainty [...]" AOR, 2018

<sup>9</sup> The plan4res project: <https://www.plan4res.eu/>

- Again, what did this have to do with graphs, please?
- Specialized DP algorithms for thermal single-Unit Commitment<sup>10</sup>  
 $\equiv$  shortest path on appropriate acyclic graph

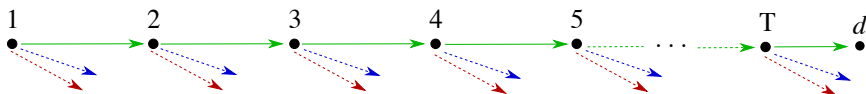


- Not too many nodes  $2(T = |\mathcal{T}|)$ , but rather dense:  $O(T^2)$  arcs
- $((t, \text{ON}), (\tau, \text{OFF})) \equiv$  startup at  $t$  and shutdown at  $\tau > t \dots$
- Costs require another nested DP per arc,  $O(T^3)$  overall
- Hence, (strong but large) formulation as a flow problem<sup>11</sup>

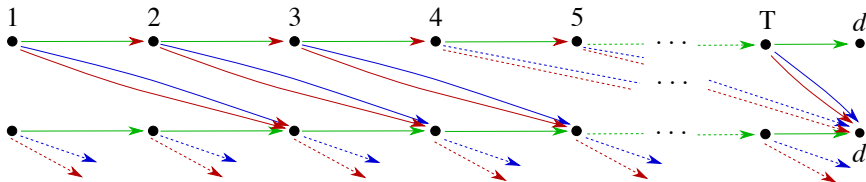
<sup>10</sup> F., Gentile "Solving Nonlinear Single-Unit Commitment Problems with Ramping Constraints" OR, 2006

<sup>11</sup> F., Gentile "New MIP Formulations for the Single-Unit Commitment Problems with Ramping Constraints" IASI 15-06, 2015

- Water flowing over time is a flow problem (surprise!)



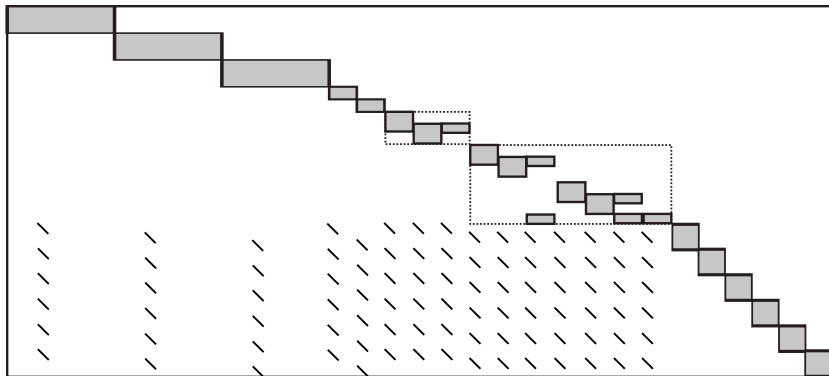
- Quite skinny graph,  $O(T)$  nodes and arcs, too
- Turbining/spilling arcs produce/not energy (max reservoir capacity)
- However, **hydro units are often whole interconnected hydro valleys**



- All in all (**without pumping**) still a flow problem, on a **structured graph** (composition of lines with a reverse tree)

- The transmission/distribution network is a graph (surprise!)  
nodes are zones/buses, arcs are links (bi-directed)
- Kirchhoff's current law:  $Af = n$  ( $f$  = flows,  $n$  = net injection)
- Kirchhoff's voltage law + Ohm's law for AC current  $\implies f = \gamma^T A^T \theta$   
( $\theta$  = voltage angles,  $\gamma$  = arc susceptances =  $1/\text{impedance}$ )
- AC  $\implies$  currents and voltages are periodic  $\equiv$  complex numbers
- **DC approximation:**  $|\theta_i - \theta_j| \ll 1$  ( $(i, j) \in A$  (small phase differences between neighbours)  $\implies$  can linearize the trigonometric functions
- $A\Gamma A^T \theta = n$  (Laplacian!) +  $\underline{f} \leq \gamma^T A^T \theta \leq \bar{f}$  (capacity)
- Fixing one reference voltage  $A\Gamma A^T$  nonsingular:  
$$\underline{f} \leq \gamma^T A^T (A\Gamma A^T)^{-1} n \leq \bar{f}$$
- True AC version nonlinear nonconvex, rather hard ...

- Not a single flow but a multicommodity flow (of sorts)



- Many blocks, either  $A$  or  $A\Gamma A^T$ , but of rather different shape and size
- Nontrivial linking constraints

- Of course we can, in fact with several different approaches:
  - Lagrangian decomposition<sup>12</sup> and related methods<sup>13</sup>, even in parallel<sup>14</sup>
  - Structured Interior-Point methods<sup>15</sup>
  - Structured active-set (simplex) methods<sup>16</sup>
  - Structured Dantzig-Wolfe decomposition<sup>17,18</sup>
  - ...
- Most can exploit the “inner” graph structure of (the many)  $A(s)$
- Significantly more complex: two-level approaches ( $\equiv$  more fun)

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<sup>12</sup> F., Gallo “A Bundle type dual-ascent approach to linear multicommodity min cost flow problems” *INFORMS JOC*, 1999  
<sup>13</sup> Grigoriadis, Khachiyan “An exponential function reduction method for block angular convex programs” *Networks*, 1995  
<sup>14</sup> Cappanera, F. “Symmetric and asymmetric parallelization of a cost-decomposition algorithm [...]” *INFORMS JOC*, 2003  
<sup>15</sup> Castro “Solving difficult multicommodity problems through a specialized interior-point algorithm” *Ann. OR*, 2003  
<sup>16</sup> McBride “Progress made in solving the multicommodity flow problem” *SIOPT*, 1998  
<sup>17</sup> F., Gendron “A stabilized structured dantzig-wolfe decomposition method” *Math. Prog.*, 2013  
<sup>18</sup> Mamer, McBride “A decomposition-based pricing procedure for large-scale linear programs [...]” *Man. Sci.*, 2000

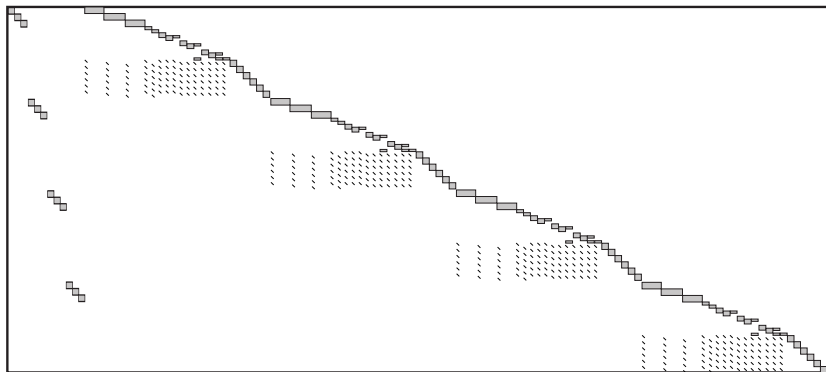


- Maybe if this were the end, but it is **just the beginning**
- **Data is uncertain**: demand, wind/solar production, units/network state ... which **cannot be ignored** (increased RES penetration ...)
- Unit commitment is decided in advance (here-and-now) but actual dispatch can be changed in real time (recourse)
- Many methods to represent uncertainty: Stochastic Optimization<sup>19</sup>, Robust Optimization, Chance-Constrained Optimization, hybrid<sup>20</sup>
- Simplest approach **scenario-based**: each  $\approx$  a full UC  
 $\implies$  **yet another two-level structure**
- Cons: **size increases of a factor  $\#$  scenarios** (which should be large)

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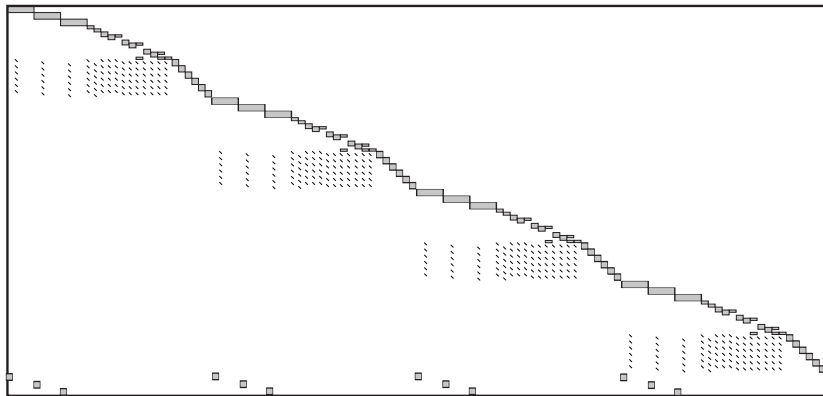
<sup>19</sup> Scuzziato, Finardi, F. "Comparing Spatial and Scenario Decomposition for Stochastic [...]" *IEEE Trans. Sust. En.*, 2018

<sup>20</sup> van Ackooij, F., de Oliveira "Inexact Stabilized Benders' Decomposition Approaches, with Application [...]" *CO&A*, 2016



- Perfect structure for Benders' decomposition
- Benders' decomposition with Lagrangian decomposition inside<sup>21</sup>
- ... with (different) graph structure(s) inside

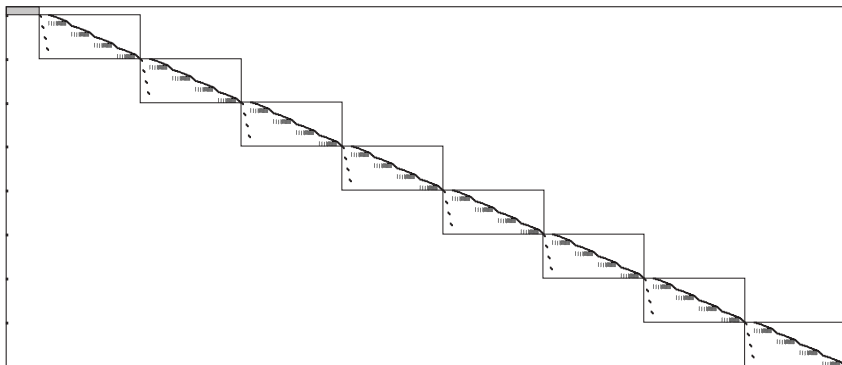
<sup>21</sup> van Ackooij, Malick "Decomposition algorithm for large-scale two-stage unit-commitment" *Ann. OR*, 2016



- Or was it the **perfect structure for Lagrangian decomposition**?
- Lagrangian decomposition with Lagrangian decomposition inside ...
- Which is better? Very hard to say beforehand<sup>19</sup>

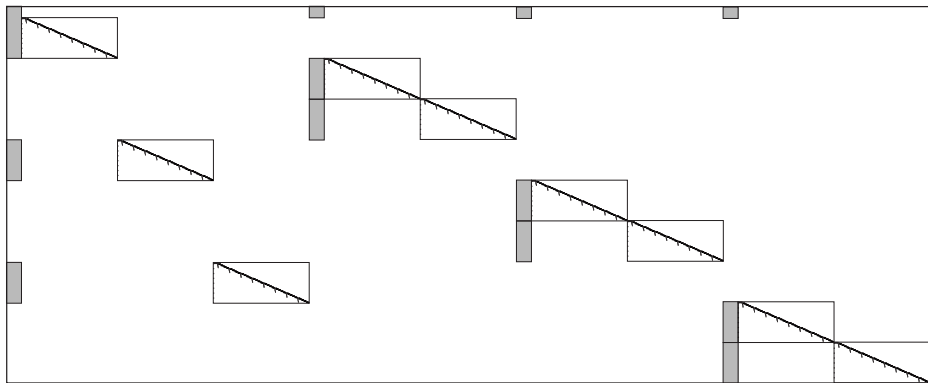
- Unit-Commitment is a short-term problem, lacks long-term strategies
- Issue: **cost of water** (none) / minimum reservoir volume (very low)  
 $\implies$  lot of water used  $\implies$  no water most of the year
- Hydro production most useful for **peak shaving** every day
- Computing **value of water left in the reservoirs at  $T$**   
 $\equiv$  **solving a parametric (uncertain) UC problem**  
**for each (significant) day of the year**
- Can approximate it by **dual variables**/Lagrangian multipliers of minimum reservoir volume constraints
- Better a **piecewise linear representation** (cutting-plane model)
- Then, stochastic dual **dynamic programming**<sup>22</sup> (another graph)

<sup>22</sup> Pereira, Pinto "Multi-stage stochastic optimization applied to energy planning" *Math. Prog.*, 1991



- This is **not really** how you'd do that (integer variables)
- Still OK for Benders-like decomposition
- Benders + Benders + Lagrange + Graph or  
Benders + Lagrange + Lagrange + Graph or  
Lagrange + Benders + Lagrange + Graph or ...

- The energy system changes all the time, but modifications **slow**, **extremely costly**, with **huge inertia**
- **Demand and production subject to very significant uncertainties:**  
climate = RES production + demand, shifts in consumption patterns (EV, cryptocurrencies, ...), new technologies (shale, LED, ...), geo-political factors (energy security), economical factors (boost or boom), regulatory factors (EU energy market, ...), political factors (CO<sub>2</sub> emission treaties, nuclear power, ...), ...
- **Planning long-term evolution very hard**, yet **necessary**
- 20/30 years, 2/5 years steps (**multi-level** recourse), **many scenarios**



- Huge size, multiple nested structure
- Still OK for either Benders or Lagrange
- Benders + Lagrange + Benders + Lagrange + Graph or ...

- **Modeling system:** easily construct a huge, flat = unstructured matrix to be passed to a general-purpose, flat solver
- Some solvers offer one-level decomposition (Benders, CG = DW)
- Attempts at automatically recovering structure from a matrix<sup>23</sup>, but only one level and anyway conceptually awkward
- Only one tool (that I know of) for multiple nested structure<sup>24,25</sup>, but only solves continuous problems by Interior Point methods
- Nothing for multilevel, heterogeneous approaches (such as, but not only, decomposition), e.g., allowing specialized solvers for each block
- So far

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<sup>23</sup> Gamrath, Lübbecke "Experiments with a Generic Dantzig-Wolfe Decomposition for Integer Programs" *LNCS*, 2010

<sup>24</sup> Gondzio, Grothey "Exploiting Structure in Parallel Implementation of Interior Point Methods [...]" *Comput. Man. Sci.*, 2009

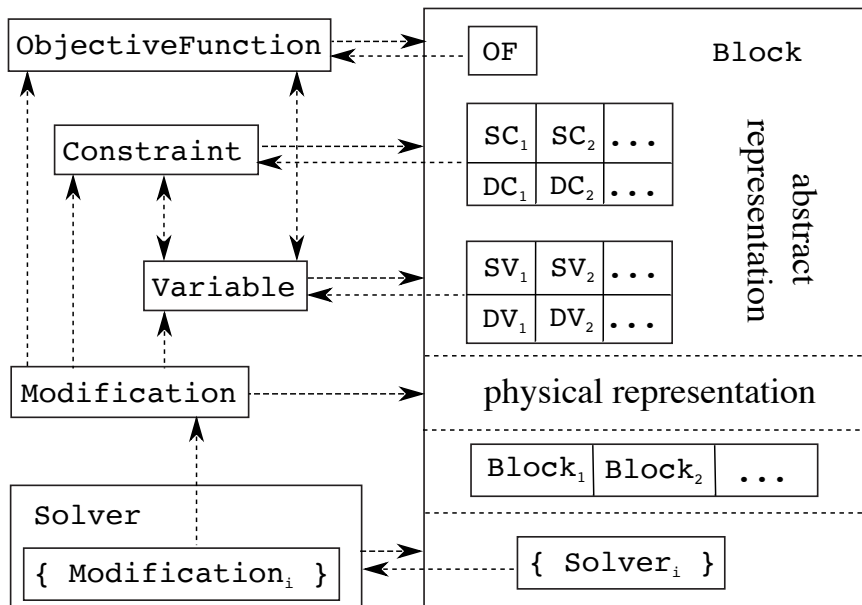
<sup>25</sup> Colombo et al. "A Structure-Conveying Modelling Language for Mathematical [...] Programming" *Mathe. Prog. Comp.*, 2009



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- A **modelling language/system** which:
  - explicitly supports the notion of **block  $\equiv$  nested structure**
  - separately provides “semantic” information from “syntactic” details (list of constraints/variables)
  - allows exploiting specialised solvers on blocks with specific structure
  - caters all needs of complex methods: dynamic generation of constraints/variables, modifications in the data, reoptimization
- C++ library: set of “core” classes, easily extendable
- Why C++? A number of reasons:
  - all serious solvers are written in C/C++
  - we all love it (especially C++11/14)
  - tried with Julia/JuMP, but could not handle well C++ interface



- **Block** = abstract class representing the general concept of “a part of a mathematical model with a well-understood identity”
- Each **Block::** a model with specific **structure** (e.g., **Block::BinKnapsackBlock** = a 0/1 knapsack problem)
- **Physical representation** of a Block: whatever data structure is required to describe the instance (e.g.,  $a$ ,  $b$ ,  $c$ )
- **Abstract representation** of a Block:
  - one (for now) **ObjectiveFunction**
  - any # of **groups** of (pointers) to **(static) Variable**
  - any # of **groups** of **std::list** of (pointers) to **(dynamic) Variable**
  - any # of **groups** of (pointers) to **(static) Constraint**
  - any # of **groups** of **std::list** of (pointers) to **(dynamic) Constraint**groups of Variable/Constraint can be single (**std::list**) or **std::vector (...)** or **boost::multi\_array** thanks to **boost::any**
- Any # of sub-Blocks (recursively), possibly of **specific type** (e.g., **Block::MMCFBlock** can have  $k$  **Block::MCFBlocks** inside)

- Abstract concept, thought to be extended (a matrix, a function, ...)
- Does **not even have a value**
- Knows which Block it belongs to
- Can be **fixed** and **unfixed** to/from its current value (whatever that is)
- Keeps the set of Constraint/ObjectiveFunction it **influences**
- **Fundamental design decision: “name” of a Variable = its memory address  $\implies$  copying a Variable makes a different Variable  $\implies$  dynamic Variables always live in `std::lists`**
- `Modification::VariableModification (fix/unfix)`

- Abstract concept, thought to be extended (any algebraic constraint, a matrix constraint, a PDE constraint, bilevel program, ...)
- Keeps the set of Variables it is **influenced from**
- Either **satisfied** or not by the current value of the Variables
- Knows which Block it belongs to
- Can be **relaxed** and **enforced**
- **Fundamental design decision: “name” of a Constraint = its memory address  $\Rightarrow$  copying a Constraint makes a different Constraint  $\Rightarrow$  dynamic Constraints always live in `std::lists`**
- `Modification::ConstraintModification (relax/enforce)`

- Abstract concept, perhaps to be extended (vector-valued ...)
- Either minimized or maximized
- Keeps the set of Variables it **depends from**
- Can be **evaluated** w.r.t. the current value of the Variables (but its value depends on the specific form)
- `ObjectiveFunction::RealObjectiveFunction` implements “value is an extended real”
- Knows which Block it belongs to
- Same fundamental design decision ... (but there is no such thing as a dynamic `ObjectiveFunction`)
- `Modification::OFModification` (change verse)

- Any # of **Solvers** attached to a Block to solve it
- **Solver::** for a **specific Block::** can use the physical representation  
⇒ no need for explicit Constraints  
⇒ abstract representation of Block only constructed on demand
- However, Variables are always present (interface with Solver)
- A **general-purpose Solver** uses the abstract representation
- **Dynamic Variable/Constraints** can be generated on demand (user cuts/lazy constraints/column generation)
- For a Solver attached to a Block:
  - Variables **not belonging to the Block** are **constants**
  - Constraints **not belonging to the Block** are **ignored**  
(belonging = declared there or in any sub-Block recursively)
- ObjectiveFunction of sub-Blocks **summed** to that of father Block if has same verse, but **min/max supported**



# Solver

- Solver = interface between a Block and algorithms solving it
- Each Solver attached to a single Block, from which it picks all the data, but any # of Solvers can be attached to the same Block
- Solutions are written directly into the Variables of the Block
- Individual Solvers can be attached to sub-Blocks of a Block
- Tries to cater for all the important needs:
  - optimal and sub-optimal solutions, provably unbounded/unfeasible
  - time/resource limits for solutions, but restarts (reoptimization)
  - any # of multiple solutions produced on demand
  - lazily reacts to changes in the data of the Block via Modifications
- Heavily slanted towards RealObjectiveFunction (optimality guarantees being upper and lower bounds)
- Derived CDASolver is “Convex Duality Aware”: bounds are associated to dual solutions (possibly, multiple)
- Something relevant may be missing, asynchronous calls not clear yet

# Block and Modification

- Most Block components can change, but not all:
  - set of sub-Blocks
  - number and shape of groups of Variables/Constraints
- Any change is communicated to each interested Solver (attached to the Block or any of its ancestor) via a Modification object
- `anyone_there()`  $\equiv \exists$  interested Solver (Modification needed)
- However, two different kinds of Modification (what changes):
  - physical Modification, only specialized Solvers concerned
  - abstract Modification, only Solvers using it concerned
- Abstract Modification on Variable/Constraint must always be issued, even if no Solver, to keep both representations in sync
- A single change may trigger more than one Modification
- A Solver will disregard a Modification it does not understand (there must always be another one it understands)
- A Block may refuse to support some changes (explicitly declaring it)

# Modification

- Almost empty base class, then everything has its own derived ones
- Each change to Block/Variable/Constraint ... produces a Modification, and a **smart pointer** is passed to the Block
- The Block funnels it to the **interested Solvers** (above, if any)
- **Heavy stuff** can be attached to a Modification (e.g., added/deleted dynamic Variable/Constraints)
- Each Solver has the **responsibility** of cleaning up its list of Modifications (smart pointers → memory will finally be released)
- Modifications **processed in the arrival order** to ensure consistency
- Solvers are supposed to **reoptimize** to improve efficiency, which is **easier if you can see all list of changes at once** (lazy update)
- A Solver may optimize the changes (Modifications may cancel each other out ...), but **its responsibility**

# Solution and Configuration

- Block produces one **Solution**, possibly using its sub-Blocks'
- A Solution can read() its own Block and write() itself back
- Solution is Block-specific rather than Solver-specific
- Solution may save dual information
- Solution may save only a specific subset of the primal/dual solution
- Block, Solution are **tree-structured complex objects**
- **Configuration** for them a (possibly) tree-structured complex object but also Configuration::SimpleConfiguration (an int)
- Configuration::BlockConfiguration sets (recursively):
  - which dynamic Variable/Constraints are generated, how (Solver, time limit ...)
  - which Solvers attached to each sub-Block
  - which Solution is produced ...

- Often **reformulation** crucial, but also **relaxation** or **restriction**:  
`get_R3_Block()` produces one, possibly using sub-Blocks'
- Obvious special case: **copy** (clone), should always work
- Available R<sup>3</sup>Blocks `Block::`-specific, a Configuration needed
- R<sup>3</sup>Block **completely independent** (**new** Variable/Constraints),  
useful for algorithmic purposes (branch, fix, solve, ...)
- Solution of R<sup>3</sup>Block useful to Solvers for original Block:  
`map_back_solution()` (best effort in case of dynamic Variables)
- Sometimes **keeping R<sup>3</sup>Block in sync with original** necessary:  
`map_forward_modifications()`, **task of original Block**
- `map_forward_solution()` and `map_back_modifications()` useful,  
e.g., **dynamic generation of Variable/Constraints** in the R<sup>3</sup>Block
- **Block::** **is in charge** of all this, thus **decides what it supports**

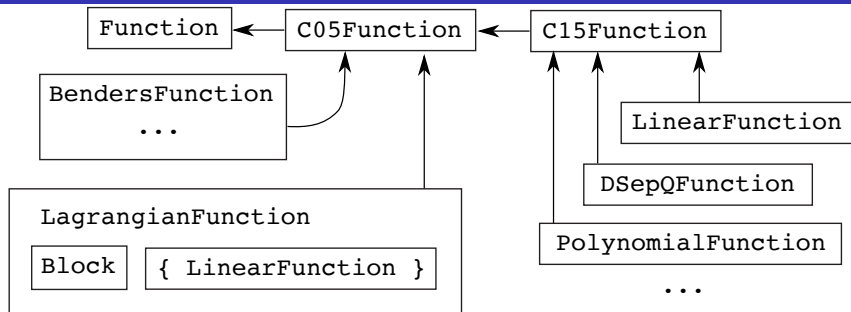
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# First Basic Implementations

- `Variable::ColVariable` implements “value = one single real”, possibly restricted to  $\mathbb{Z}$ , with (possibly infinite) bounds
- `Modification::ColVariableModification` (change bounds, type)
- `Constraint::RowConstraint` implements “ $l \leq \text{a real} \leq u$ ”
- Has dual variable attached to it (single real)
- `Modification::RowConstraintModification` (change  $l$ ,  $u$ )
- `RowConstraint::FRowConstraint`: “a real” given by a `Function`
- `RealObjectiveFunction::FRealObjectiveFunction`: “value” given by a `Function`

# Function



- Function only deals with (real) **values**
- **Approximate computation** supported in a quite general way<sup>26</sup>
- **Asynchronous evaluation** **still not defined**
- Handles set of Variables upon which it depends
- `FunctionModification[Variables]` for “easy” changes  $\implies$  **reoptimization** (shift, adding/removing “**quasi separable**” Variables)

<sup>26</sup> van Ackooij, F. “Incremental bundle methods using upper models” *SIOPT*, 2018



- C05Function/C15Function deal with 1<sup>st</sup>/2<sup>nd</sup> order information (not necessarily continuous)
- General concept of “linearization” (gradient, convex/concave subgradient, Clarke subgradient, ...)
- Multiple linearizations produced at each evaluation (local pool)
- Global pool of linearizations for reoptimization:
  - convex combination of linearizations
  - “important linearization” (at optimality)
- C05FunctionModification[Variables/LinearizationShift] for “easy” changes  $\implies$  reoptimization (linearizations shift, some linearizations entries changing in simple ways)
- C15Function supports Hessians, unclear how much reoptimization possible/useful

# LagrangianFunction

- `C05Function::LagrangianFunction` has one **isolated** Block + set of (so far) `LinearFunction` to define Lagrangian term
- `evaluate() = Block.get_registered_solvers()[ i ].solve(): asynchronous Solver  $\Rightarrow$  asynchronous Function`
- `Solutions` extracted from Block  $\equiv$  linearizations
- Solver provides local pool
- `LagrangianFunction` handles global pool
- All changes lead to reoptimization-friendly Modification
- `BendersFunction` should be quite similar

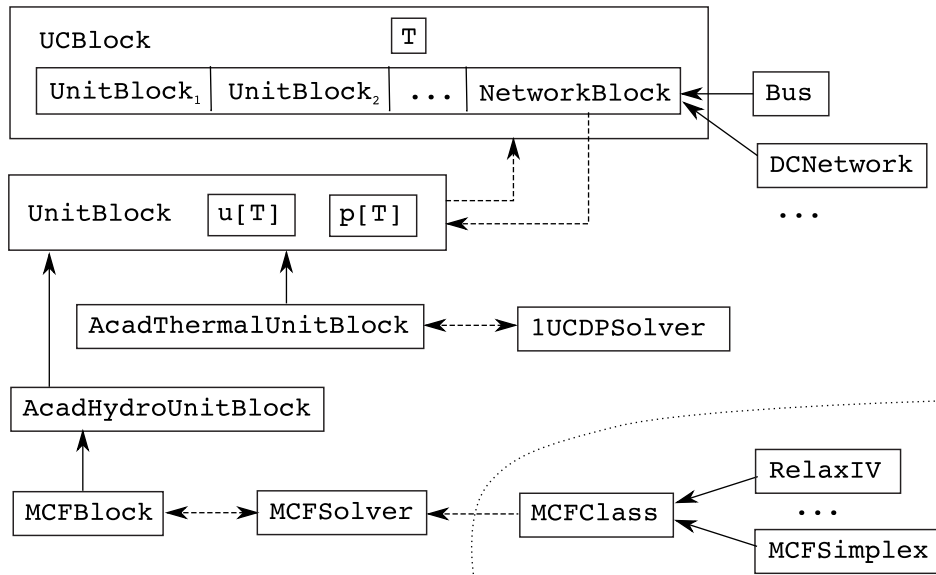
# Other useful stuff

- `un_any_thing()` template functions/macros to extract (`std::vector` or `boost::multi_array` of) (`std::list` of) `Variable/Constraints` out of a `boost_any` and work on that
- `Solution::ColVariableSolution` uses the abstract representation of any `Block` that only have (`std::vector` or `boost::multi_array` of) (`std::list` of) `ColVariables` to read/write the solution
- `Solution::RowConstraintSolution` uses the abstract representation of any `Block` that only have (...) `RowConstraints` to read/write the `dual` solution
- Of course, `Solution::CVFRRSolution` ...
- `Solver::MILPSolver` solves with `Cplex` any `Block` that only has (...) `ColVariables`, `FRowConstraints` and `FRealObjectiveFunction` with `LinearFunctions` (uses the abstract representation)

# Outline

- 1 Structure: top-down
- 2 Structure: bottom-up
- 3 The Core Elements of SMS++
- 4 The quasi-Core Elements of SMS++
- 5 Example: SMS++ for the Unit Commitment
- 6 Conclusions and (a Lot of) Future Work

# UCBlock and Companion Classes





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# A Lot of Work, Then Maybe Conclusions

- Alpha version, not all the features you have seen are complete
- Design principles have kept evolving, new ideas continue to crop up
- Core nicely general, but only success in applications validate it
- Heavily slanted towards optimization, useful for numerical analysis?
- Really  $\neq$  from all I've seen so far, had to invent almost everything
- Overhead still largely unknown (although C++ efficient)
- Asynchronous still to be figured out (but very relevant)
- Clearly not for the faint of heart ...



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but when it'll work it will be useful in many applications
- Implementing general, flexible methods for heterogeneous, multi-level structured problems is highly complex, have to make the tools first

We are trying. Someone cares to join?

# Acknowledgements

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