# **Column Generation**

Because we cannot spend all day pricing

### **Column Generation**

### We now have a mechanism to price all variables not in the pool

...But we still need to handle an exponential number of them

■ I.e. enumerating paths would still be prohibitively expensive

What can we do about it?

### **Column Generation**

### We now have a mechanism to price all variables not in the pool

...But we still need to handle an exponential number of them

■ I.e. enumerating paths would still be prohibitively expensive

#### What can we do about it?

- There is no need to find all paths with negative  $\frac{\partial}{\partial x_j} f(x) < 0$
- We just need to determine whether one such path exists

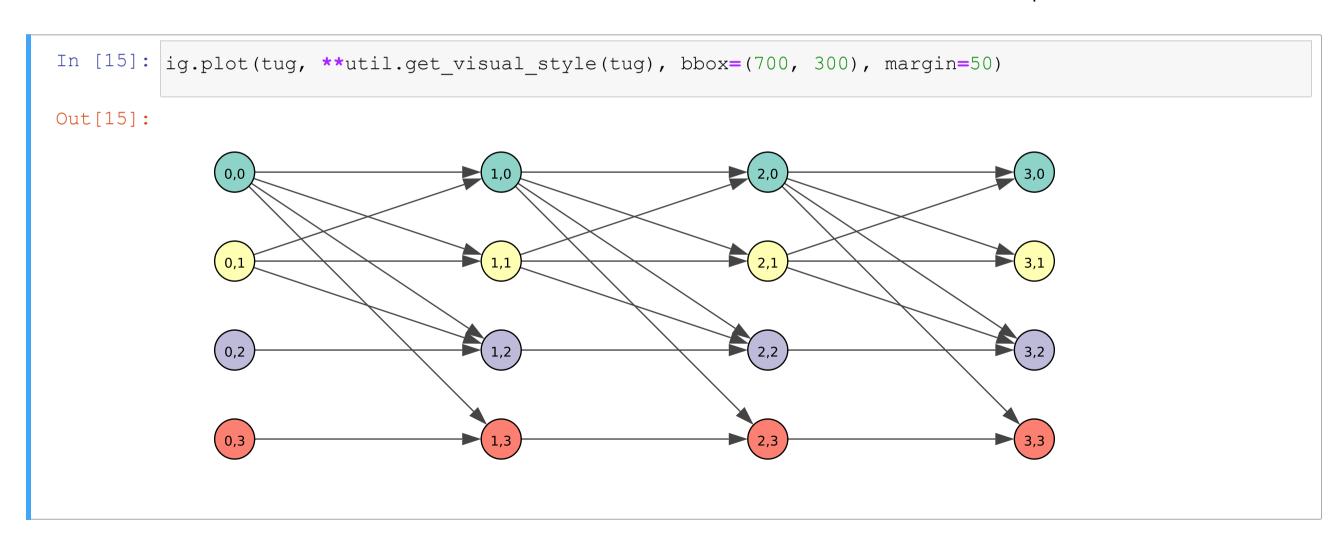
# Hence, we can build a variable with the most negative $\frac{\partial}{\partial x_j} f(x) < 0$

- Since variables correspond to columns in LP
- ...This approach is called Column Generation

# **Pricing ad Optimization**

### In practice, we view pricing as an optimization problem

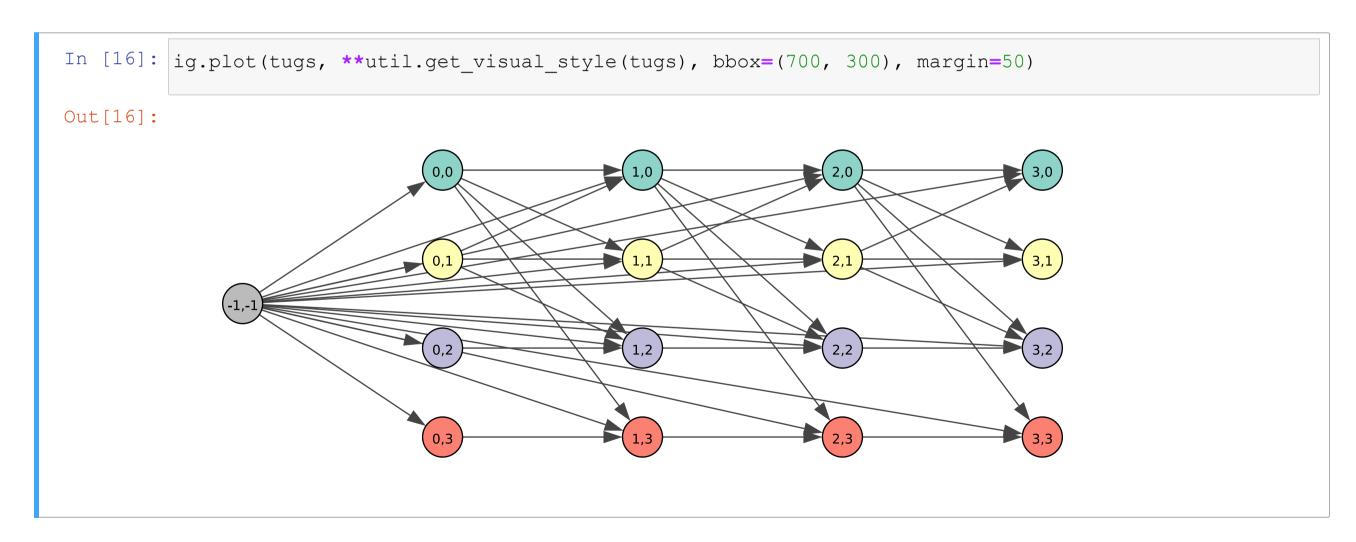
- In our case we are looking at paths
- ...So it make sense to visualize them on the Time Unfolded Graph



# **Pricing ad Optimization**

### When you want to consider all paths in a directed graph

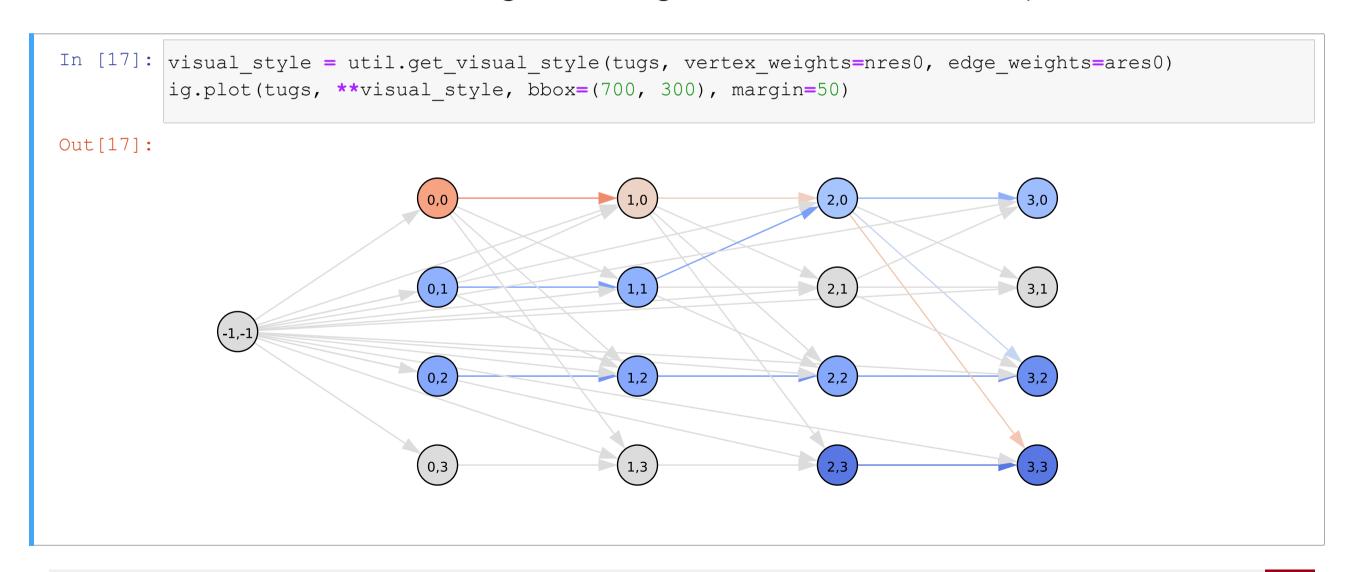
- ...It is convenient to add a fake source node
- Then you can assume that all paths start from that node



# **Pricing as Optimization**

### We can treat our residual as node and arc weights

- In the plot, a grey shade corresponds to near-zero weight
- A blue shade is used for negative weights, and a red share for positive ones



# **Shortest Path Algorithms**

### **Optimization problems over graphs**

...Are often amenable to dedicated, very efficient, algorithms

What about our graph?

# **Shortest Path Algorithms**

### Optimization problems over graphs

...Are often amenable to dedicated, very efficient, algorithms

### What about our graph?

### Weights can be negative

■ ...So one may think of using the Bellman-Ford algorithm, which runs in  $O(n_v n_e)$ 

### ...But this is a Direct, Acyclic Graph (DAG)

- Meaning that we can process the nodes in topological order
- lacktriangleright ...And apply Dijkstra algorithm, which runs in  $O(n_e)$

# Dijkstra's Algorithm for DAGs

### Intuitively, we proceed as follows

- $\mathbf{Q} = [0]$  # We enqueue the fake source node
- $lacksquare sp_i = [], \forall i = 0..n_v \# All shortest paths are empy$
- while |Q| > 0:
  - $\blacksquare$  pop a node *i* from Q
  - **append** i to  $sp_i$ # Extend the shortest path
  - $\blacksquare$  for *j* successor of *i*:
    - $\blacksquare$  mark the arc (i, j) as visited
    - $\blacksquare$  if the shortest path passing for *i* is shorter than  $sp_j$ 
      - lacksquare update  $sp_i$ # Keep only the shortest path to j
    - if all ingoing arcs for j have been visited
      - $\blacksquare$  append j in Q

# **Pricing via Shortest Paths**

### The approach is implemented in the solve\_pricing\_problem function

...Which returns shortest paths to all TUG nodes

```
In [19]: ncosts a, npaths a = util.solve pricing problem(tug, rflows0, rpaths0,
                                                      node counts, arc counts, filter paths=False)
         print('COST: PATH')
         util.print solution(tug, ncosts a, npaths a, sort=None)
         COST: PATH
         -39.66: 0,2 > 1,2 > 2,2 > 3,2
         -31.37: 0,1 > 1,1 > 2,0 > 3,0
         -31.37: 0.1 > 1.1 > 2.0 > 3.3
         -27.36: 0.2 > 1.2 > 2.2
         -23.18: 0,1 > 1,1 > 2,0
         -23.18: 0,1 > 1,1 > 2,0 > 3,1
         -16.42: 0.2 > 1.2
         -14.68: 0.1 > 1.1
         -14.68: 0,1 > 1,1 > 2,1
         -11.77: 0,1 > 1,0 > 2,3
         -5.47:0.2
         -4.89:0.1
         -3.60: 0,1 > 1,0
         0.00: 0,3
         0.00: 1,3
         4.61: 0,0
```

# **Pricing via Shortest Paths**

### We can ask for paths with a negative weighs/gradient term

```
In [20]: ncosts, npaths = util.solve pricing problem(tug, rflows0, rpaths0,
                                                      node counts, arc counts, filter paths=True)
         print('COST: PATH')
         util.print solution(tug, ncosts, npaths, sort=None)
         COST: PATH
         -39.66: 0,2 > 1,2 > 2,2 > 3,2
         -31.37: 0,1 > 1,1 > 2,0 > 3,0
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         -14.68: 0,1 > 1,1 > 2,1
         -11.77: 0,1 > 1,0 > 2,3
         -5.47: 0,2
         -4.89: 0,1
         -3.60: 0,1 > 1,0
```

- Returning multiple paths is usually a good idea
- ...Since it typically speeds up the convergence of our dynamic method

# Let's Loop!

Time to start iterating

### Every complex endeavor is worth a double (or triple) check

Let's check again our baseline result:

```
In [21]: rflows0, rpaths0 = util.solve_path_selection_full(tug, node_counts, arc_counts, initial_paths=path_pool, verbose=0)
print('FLOW: PATH')
util.print_solution(tug, rflows0, rpaths0, sort='descending')
sse = util.get_reconstruction_error(tug, rflows0, rpaths0, node_counts, arc_counts)
print(f'RSSE: {np.sqrt(sse):.2f}')

FLOW: PATH
1.96: 0,0 > 1,0 > 2,0 > 3,2
1.86: 0,0 > 1,0 > 2,0 > 3,3
0.79: 0,0 > 1,0 > 2,0 > 3,0
RSSE: 25.58
```

### Every complex endeavor is worth a double (or triple) check

Let's try adding paths with non-negative gradient terms

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Let's try adding paths with non-negative gradient terms

- This is perfectly useless
- ...Just as expected!

### Every complex endeavor is worth a double (or triple) check

Now, let's try again with paths having negative gradient terms

```
In [23]: n paths = [p for p, c in zip(npaths a, ncosts a) if c < 0]
         path pool1 = path pool + n paths
         rflows1, rpaths1 = util.solve path selection full(tug, node counts, arc counts,
                                                            initial paths=path pool1, verbose=0)
         print('FLOW: PATH')
         util.print solution(tug, rflows1, rpaths1, sort='descending')
         sse = util.get reconstruction error(tug, rflows1, rpaths1, node counts, arc counts)
         print(f'RSSE: {np.sqrt(sse):.2f}')
         FLOW: PATH
         5.79: 0.2 > 1.2 > 2.2 > 3.2
         3.05: 0.1 > 1.1 > 2.0 > 3.3
         2.52: 0.1 > 1.1 > 2.0 > 3.0
         1.70: 0.1 > 1.0 > 2.3
         1.13: 0,0 > 1,0 > 2,0 > 3,2
         0.87: 0.0 > 1.0 > 2.0 > 3.3
         0.34: 0,0 > 1,0 > 2,0 > 3,0
         RSSE: 15.18
```

### Every complex endeavor is worth a double (or triple) check

Now, let's try again with paths having negative gradient terms

```
In [23]: n paths = [p for p, c in zip(npaths_a, ncosts_a) if c < 0]</pre>
         path pool1 = path pool + n paths
         rflows1, rpaths1 = util.solve path selection full(tug, node counts, arc counts,
                                                            initial paths=path pool1, verbose=0)
         print('FLOW: PATH')
         util.print solution(tug, rflows1, rpaths1, sort='descending')
         sse = util.get reconstruction error(tug, rflows1, rpaths1, node counts, arc counts)
         print(f'RSSE: {np.sqrt(sse):.2f}')
         FLOW: PATH
         5.79: 0.2 > 1.2 > 2.2 > 3.2
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         0.34: 0.0 > 1.0 > 2.0 > 3.0
         RSSE: 15.18
```

This time we have a better solution!

### **The Column Generation Code**

### Our CG code can be found in the trajectory\_extraction\_cg function

First, we define our initial path pool

```
paths = [[v.index] for v in tug.vs]
```

■ We use one path per node, consisting of the node itself

Then we start looping:

```
for it in range(max_iter):
    # Solve the master problem
    # Solve the pricing problem
    ...
```

■ We control the total run-time via an iteration limit

### The Column Generation Code

### Our CG code can be found in the trajectory\_extraction\_cg function

When we add the new paths, we take care of discarding duplicates

```
old_as_set = set([tuple(p) for p in paths])
found_as_set = set([tuple(p) for p in np])
new_as_set = old_as_set.union(found_as_set)
```

- Duplicates should not theoretically arise
- ...But they may in practice due to numerical errors
- ...Or when we use approximate solvers

We trigger an eary stop if no new path can be added:

```
if nnew == 0: break
```

### **Column Generation in Action**

### Let's test CG on that graph that took ~10 sec with the baseline

```
In [24]: g8_5, t8_5, f8_5, p8_5, nc8_5, ac8_5 = util.get_default_benchmark_graph(nnodes=8, eoh=5, seed=42
%time f8_5, p8_5 = util.trajectory_extraction_cg(t8_5, nc8_5, ac8_5, max_iter=30, verbose=1)

It.0, sse: 310.26, #paths: 45, new: 5
It.1, sse: 91.33, #paths: 46, new: 1
It.2, sse: 0.00, #paths: 46, new: 0
CPU times: user 66.6 ms, sys: 17.1 ms, total: 83.7 ms
Wall time: 69.4 ms
```

### What if the graph is bigger?

```
In [25]: g20_7, t20_7, f20_7, p20_7, nc20_7, ac20_7 = util.get_default_benchmark_graph(nnodes=20, eoh=7,
%time f20_7, p20_7 = util.trajectory_extraction_cg(t20_7, nc20_7, ac20_7, max_iter=30, verbose=1)

It.0, sse: 160.78, #paths: 176, new: 36

It.1, sse: 11.04, #paths: 177, new: 1

It.2, sse: 7.59, #paths: 189, new: 12

It.3, sse: 0.00, #paths: 189, new: 0

CPU times: user 372 ms, sys: 3.68 ms, total: 375 ms
Wall time: 370 ms
```

### **Column Generation in Action**

### Let's scale up even more

With this graph, the total number of paths is  $O(40^{10})$ 

```
In [26]: g40_10, t40_10, f40_10, p40_10, nc40_10, ac40_10 = util.get_default_benchmark_graph(nnodes=40, e%time f40_10, p40_10 = util.trajectory_extraction_cg(t40_10, nc40_10, ac40_10, max_iter=30, verk)

It.0, sse: 947.53, #paths: 572, new: 172
It.1, sse: 450.66, #paths: 676, new: 104
It.2, sse: 134.47, #paths: 765, new: 89
It.3, sse: 9.58, #paths: 774, new: 9
It.4, sse: 8.88, #paths: 943, new: 169
It.5, sse: 8.00, #paths: 1088, new: 145
It.6, sse: 0.00, #paths: 1088, new: 0
CPU times: user 7.42 s, sys: 20.2 ms, total: 7.44 s
Wall time: 7.42 s
```

- The adds a small fraction of the total paths
- Convergence is fast
- ...And we manage to prove optimality!

### **Some Considerations**

### Column Generation is not easy approach to setup

...But when it works, it can provide many advantages

- The master can stay remarkably clean
- Complicated constraints can be moved in the variable definitions
- ...And tackled in the pricing problem
- Scalability is pretty good, given the humongous search space

CG makes you want to write models with massive number of variables

#### Some caveats

- A heuristic may still be faster (no sound mathematical theory, though...)
- It works well when you can put all its advantages to use
- ...In particular, the master problem structure should be very clean