

Power Grid Reduction by Sparse Convex Optimization

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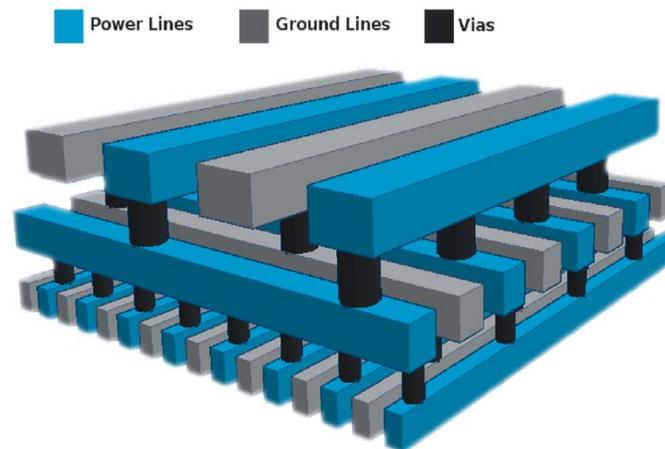
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On-chip Power Delivery Network

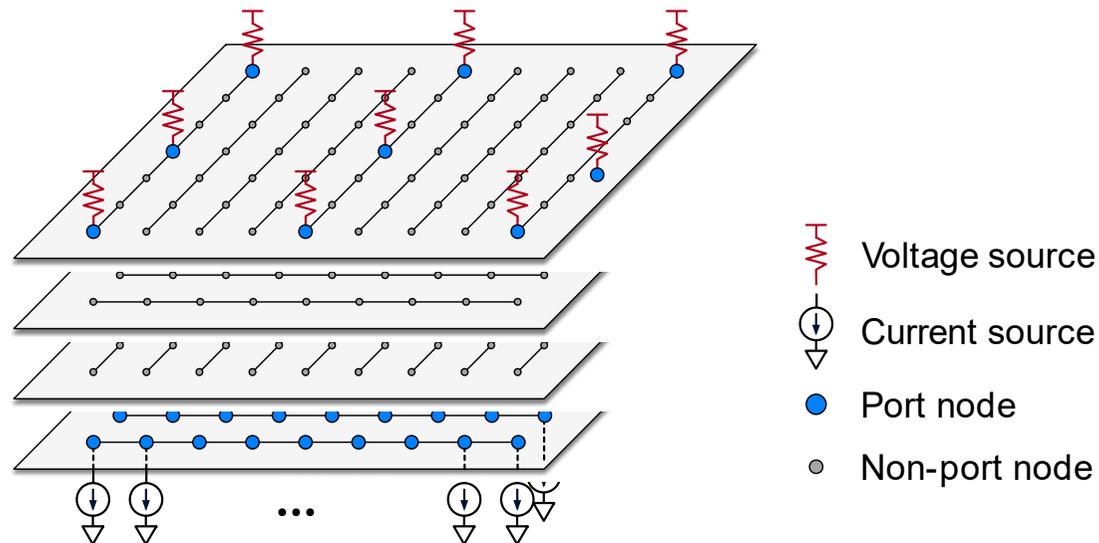
- ◆ Power grid
 - › Multi-layer mesh structure
 - › Supply power for on-chip devices
- ◆ Power grid verification
 - › Verify current density in metal wires (EM)
 - › Verify **voltage drop** on the grids
 - › More **expensive** due to increasing sizes of grids
 - › e.g., 10M nodes, >3 days



[Yassine+, ICCAD'16]

Modeling Power Grid

- ◆ Circuit modeling
 - › Resistors to represent metal wires/vias
 - › Current sources to represent current drawn by underlying devices
 - › Voltage sources to represent external power supply
 - › Transient: capacitors are attached from each node to ground
- ◆ Port node: node attached current/voltage sources
- ◆ Non-port node: only has internal connection



Linear System of Power Grid

- ◆ Resistive grid model:

$$Lv = i$$

- › L is $n \times n$ **Laplacian** matrix (symmetric and diagonally-dominant):

$$L_{i,j} = \begin{cases} \sum_{k,k \neq i} g(i,k), & \text{if } i = j \\ -g(i,j), & \text{if } i \neq j \end{cases}$$

- › $g_{i,j}$ denotes a physical conductance between two nodes i and j

- ◆ A power grid is **safe**, if $\forall i$:

$$v_i \leq V_{th}$$

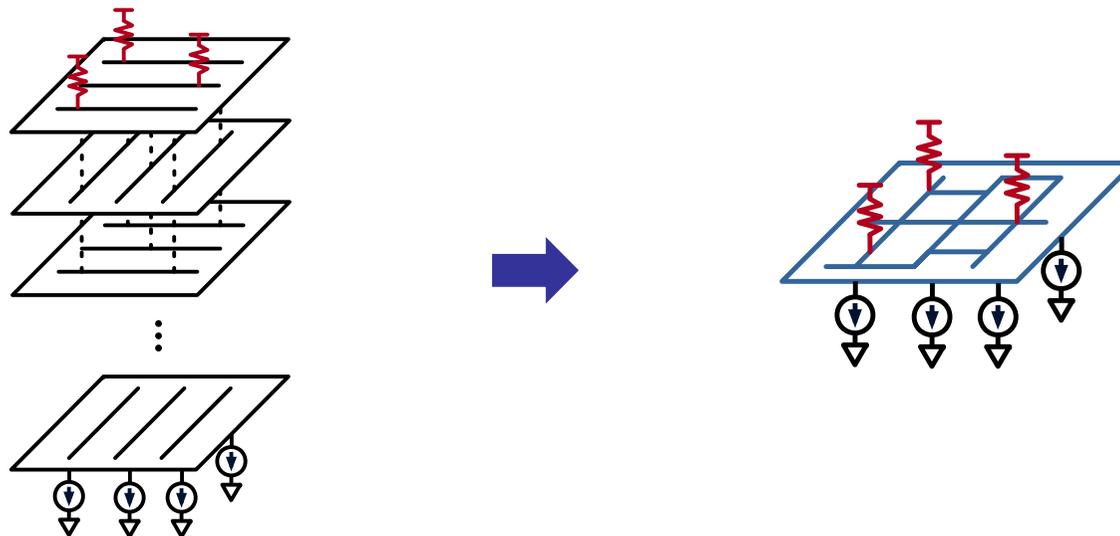
- ◆ Long runtime to solve $Lv = i$ for large linear systems

Previous Work

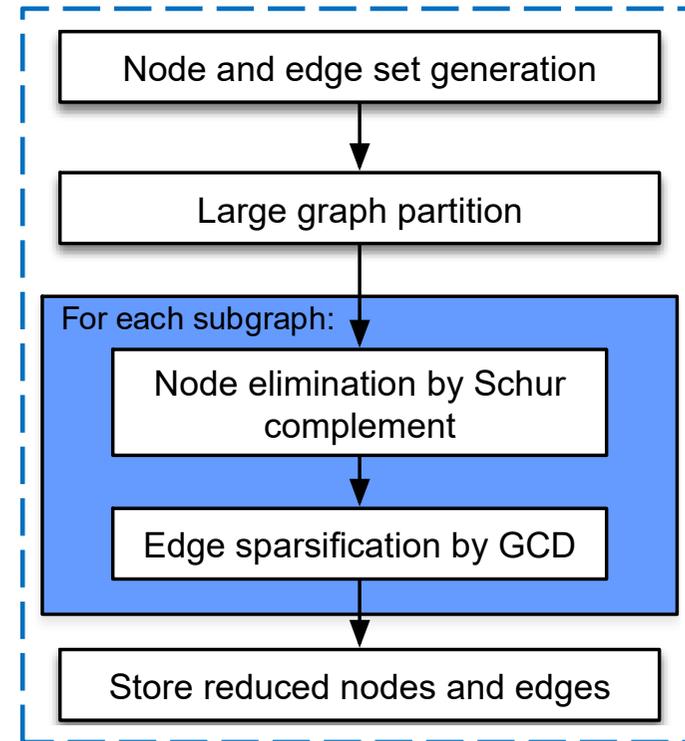
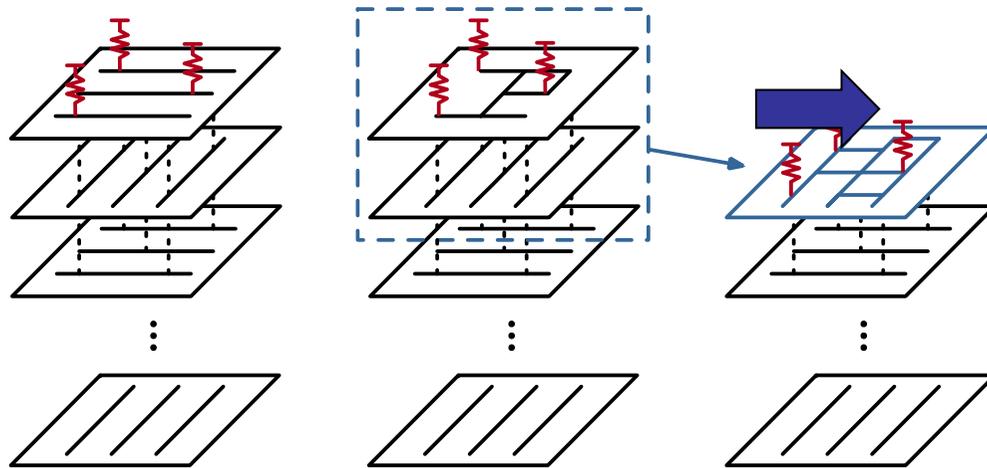
- ◆ Power grid reduction
 - › Reduce the size of power grid while preserving input-output behavior
 - › Trade-off between **accuracy** and reduction **size**
- ◆ Topological methods
 - › TICER [Sheehan+, ICCAD'99]
 - › Multigrid [Su+, DAC'03]
 - › Effective resistance [Yassine+, ICCAD'16]
- ◆ Numerical methods
 - › PRIMA [Odabasioglu+, ICCAD'97]
 - › Random sampling [Zhao+, ICCAD'14]
 - › Convex optimization [Wang+, DAC'15]

Problem Definition

- ◆ Input:
 - › Large power grid
 - › Current source values
- ◆ Output: reduced power grid
 - › Small
 - › Sparse (as input grid)
 - › Keep all the port nodes
 - › Preserve the accuracy in terms of voltage drop error



Overall Flow



Node Elimination

- ◆ Linear system: $Lv = i$
- ◆ L can be represented as a 2×2 block-matrix:

$$L = \begin{bmatrix} L_{11} & L_{12} \\ L_{12}^T & L_{22} \end{bmatrix}$$

- ◆ v and i can be represented as follows:

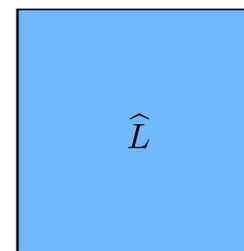
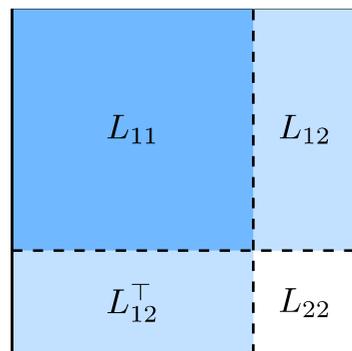
$$v = \begin{bmatrix} v_1 \\ v_2 \end{bmatrix} \text{ and } i = \begin{bmatrix} i_1 \\ 0 \end{bmatrix}$$

- ◆ Applying **Schur complement** on the DC system:

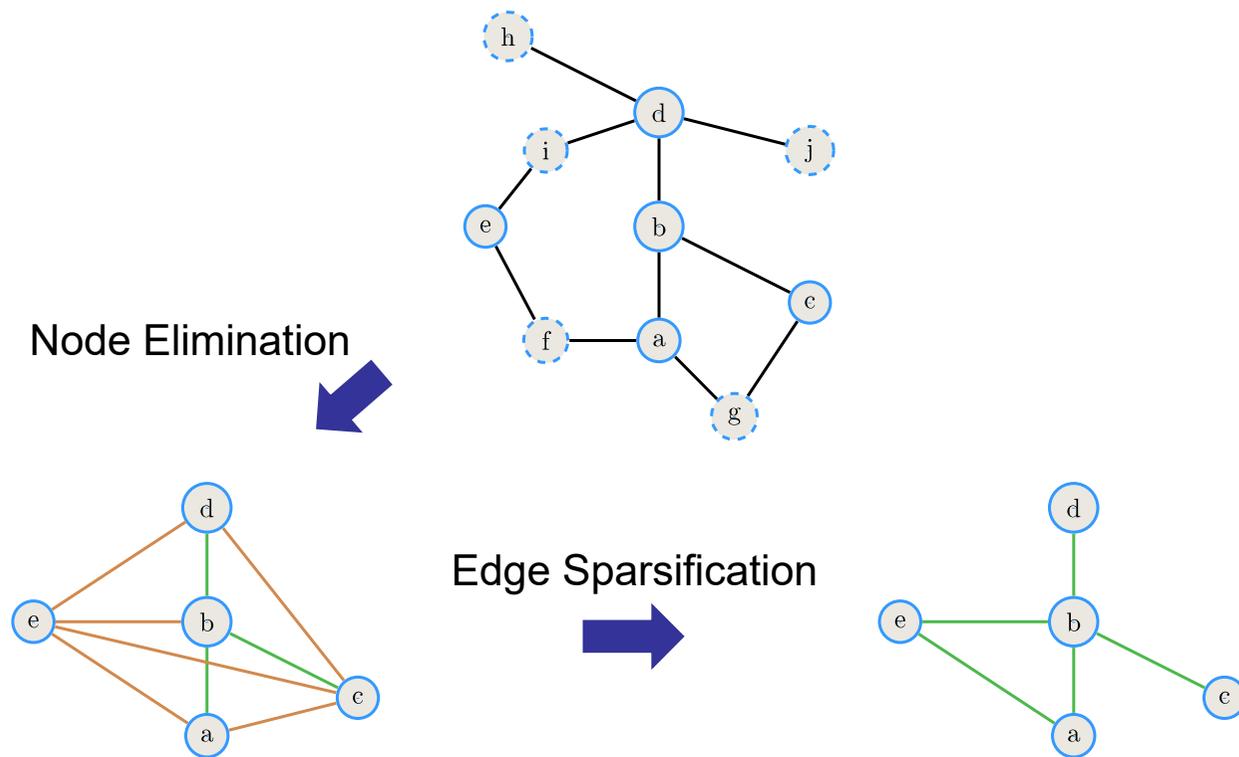
$$\hat{L} = L_{11} - L_{12}L_{22}^{-1}L_{12}^T$$

which satisfies:

$$\hat{L}v_1 = i_1$$



Node Elimination (cont'd)



- ◆ Output graph keeps all the nodes of interest
- ◆ Output graph is **dense**
- ◆ Edge sparsification: sparsify the reduced Laplacian without losing accuracy

Edge Sparsification

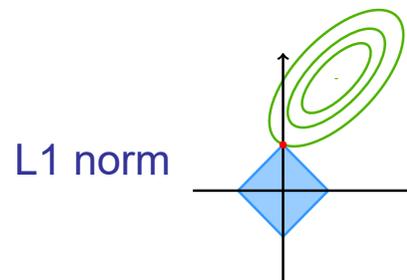
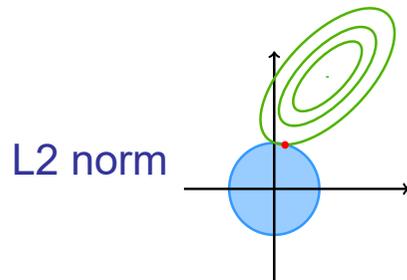
- ◆ Goal of edge sparsification
 - › **Accuracy**
 - › **Sparsity** reduce the nonzero elements off-the-diagonal in L
- ◆ Formulation (1):

$$\min_{X \in \mathbb{R}^{n \times n}} \frac{1}{2m} \sum_{k=1}^m \|(X - L)v_k\|_2^2 + \lambda \|X\|_0, \quad \text{s.t. } X \text{ is a Laplacian matrix}$$

- ◆ Formulation (2): [Wang+, DAC2014]

$$\min_{X \in \mathbb{R}^{n \times n}} \frac{1}{2m} \sum_{k=1}^m \|(X - L)v_k\|_2^2 + \lambda \|X\|_1, \quad \text{s.t. } X \text{ is a Laplacian matrix}$$

$$\min_{X \in \mathbb{R}^{n \times n}} \frac{1}{2m} \sum_{k=1}^m \|(X - L)v_k\|_2^2 + \lambda \sum_{i=1}^n X_{i,i}, \quad \text{s.t. } X \text{ is a Laplacian matrix}$$



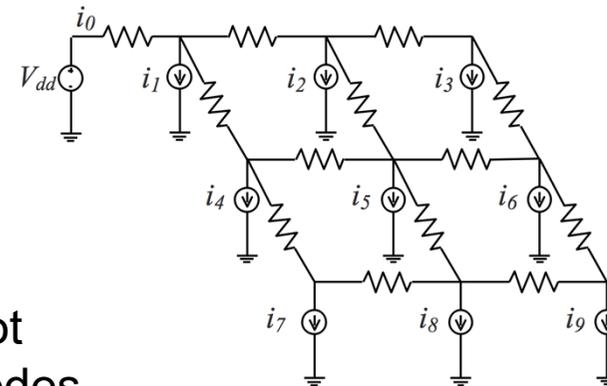
Edge Sparsification

◆ Formulation (2): [DAC2014 Wang+]

$$\min_{X \in \mathbb{R}^{n \times n}} \frac{1}{2m} \sum_{k=1}^m \|(X - L)v_k\|_2^2 + \lambda \sum_{i=1}^n X_{i,i}, \quad \text{s.t. } X \text{ is a Laplacian matrix}$$

$$|\Delta i_0|^2 + |\Delta i_1|^2 + \dots + |\Delta i_9|^2$$

$$i_0 \gg i_k, \forall i = 1, \dots, 9$$



Problem: accuracy on the V_{dd} node does not guarantee accuracy on the current source nodes

◆ Formulation (3):

$$\min_{X \in \mathbb{R}^{n \times n}} \frac{1}{2m} \sum_{k=1}^m \|((X - L)v_k) \circ w\|_2^2 + \lambda \sum_{i=1}^n X_{i,i}, \quad \text{s.t. } X \text{ is a Laplacian matrix}$$

- › Weight vector: $w_0 = 1/n, w_i = 1, \forall i = 1, \dots, n$
- › Strongly convex and coordinate-wise Lipschitz smooth

Coordinate Descent (CD) Method

- ◆ Update **one coordinate** at each iteration

- ◆ Coordinate descent:

Set $t = 1$ and $X^1 = 0$

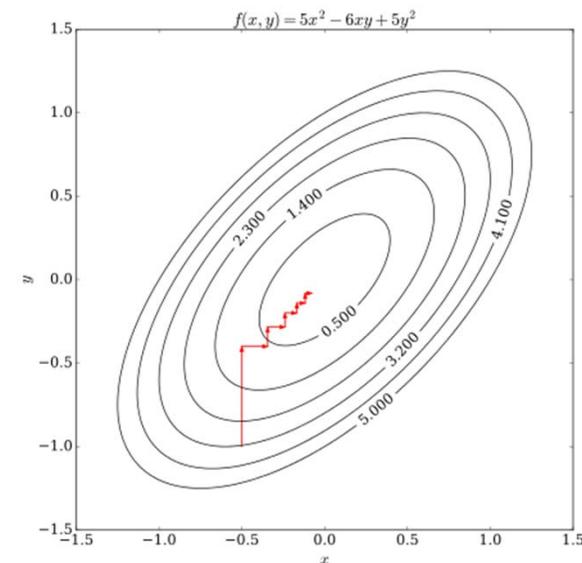
For a fixed number of iterations (or convergence is reached):

Choose a coordinate (i, j)

Compute the step size δ^* by minimizing
$$\operatorname{argmin}_{\delta} f(X + \delta e_{i,j})$$

Update $X_{i,j}^{t+1} \leftarrow X_{i,j}^t + \delta^*$

- ◆ How to decide the coordinate?
 - › Cyclic (CCD)
 - › Random sampling (RCD)
 - › **Greedy coordinate descent (GCD)**

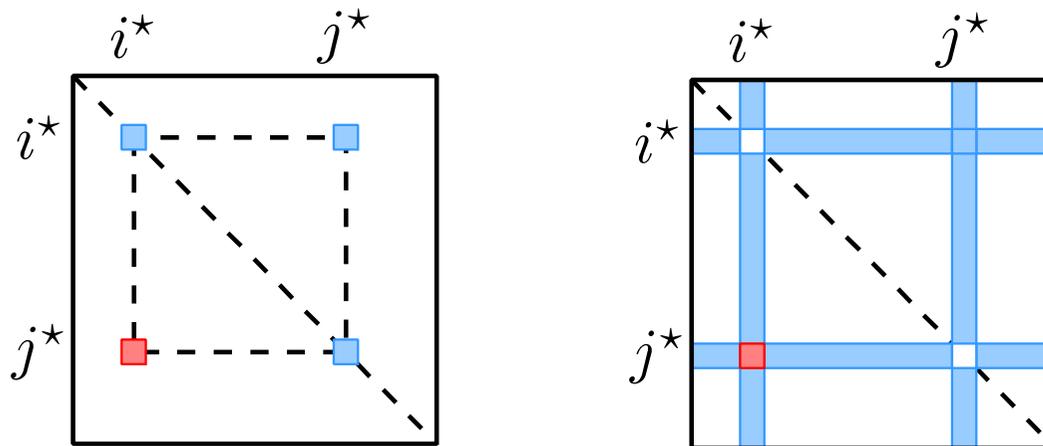


CD vs Gradient Descent

- ◆ Gradient descent (GD) algorithm:

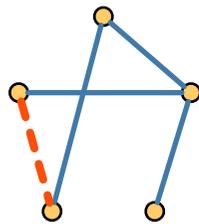
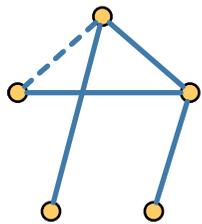
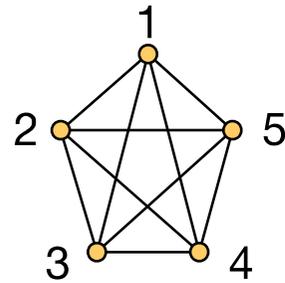
$$X^{t+1} \leftarrow X^t - \alpha \nabla f(X)$$

- ◆ GD/SGD update $O(n^2)$ elements in X and gradient matrix G at each iteration
- ◆ CD updates $O(1)$ elements in X (Laplacian property)
- ◆ CD proves to update $O(n)$ elements in G for Formulation (2) and (3).

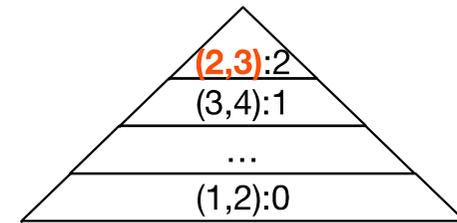
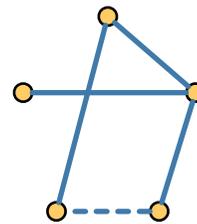


Greedy Coordinate Descent (GCD)

Input L

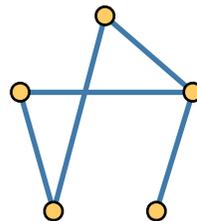


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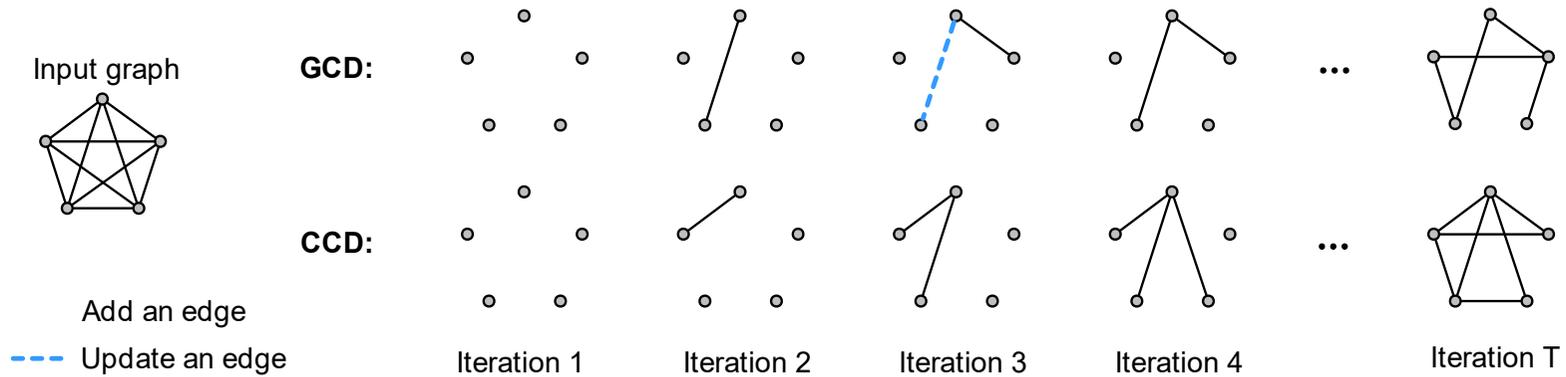


Max-heap

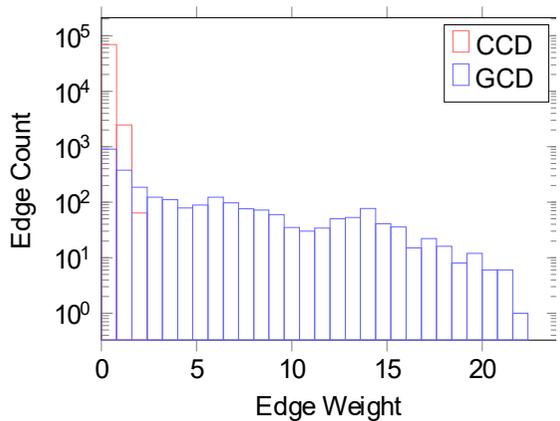
Output X



GCD vs CCD



- ◆ GCD produces sparser results
 - › CCD (RCD) goes through all coordinates repeatedly
 - › GCD selects the most significant coordinates to update

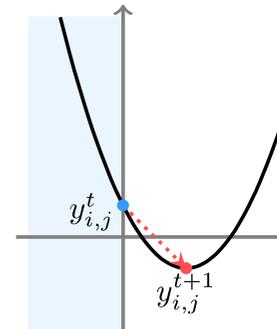
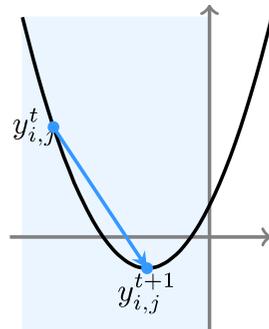


GCD Coordinate Selection

- ◆ General Gauss-Southwell Rule:

$$(i^*, j^*) = \arg \max_{(i,j) \in [n] \times [n]} |G_{i,j}|$$

- ◆ **Observation:** the objective function is quadratic w.r.t. the chosen coordinate
- ◆ GCD is stuck for some corner cases:

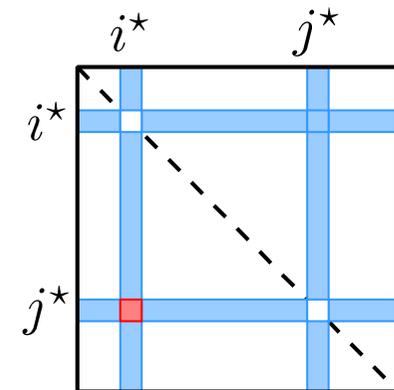


- ◆ A new coordinate selection rule:

$$(i^*, j^*) = \arg \max_{(i,j) \in [n] \times [n]} |G_{i,j}| \quad \text{s.t. } G_{i,j} > 0 \text{ or } y_{i,j} \neq 0$$

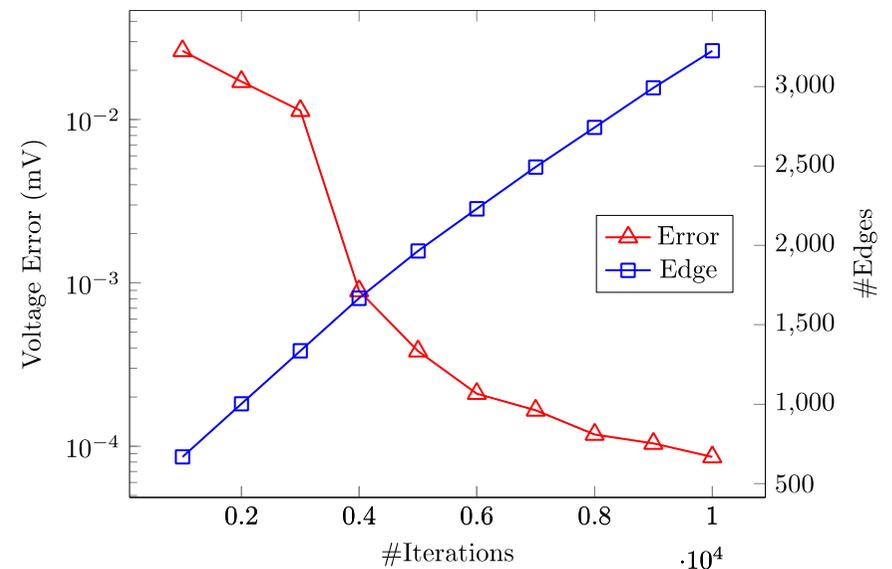
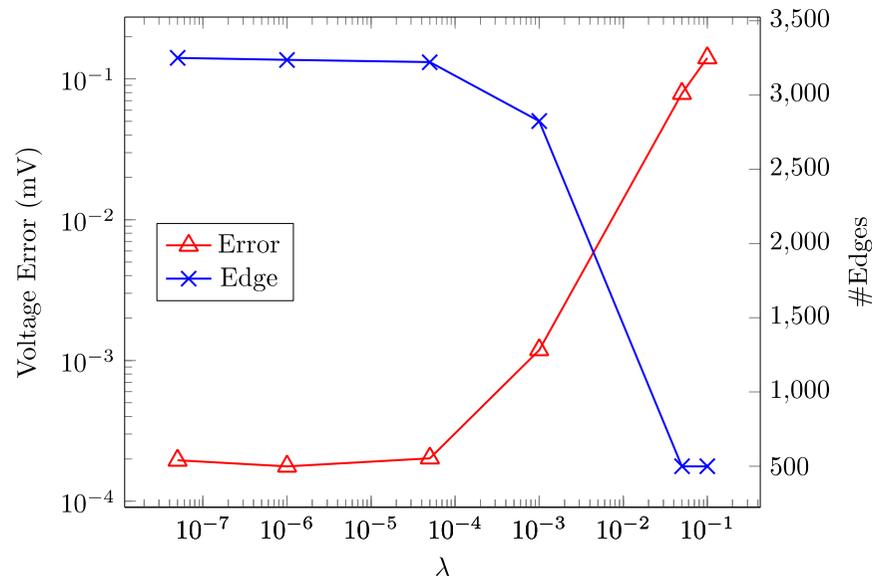
GCD Speedup

- ◆ Time complexity is $O(n^2)$ per iteration
 - › traverse $O(n^2)$ elements to get the best index
 - › As expensive as gradient descent
- ◆ **Observation:** each node has at most n neighbors \rightarrow heap
- ◆ Heap to store $O(n^2)$ elements in G :
 - › Pick the largest gradient, $O(1)$
 - › Update $O(n)$ elements, $O(n \log n)$
- ◆ Lookup table
 - › $O(n^2)$ space; $O(1)$ for each update
- ◆ **Improved** time complexity $O(n \log n)$

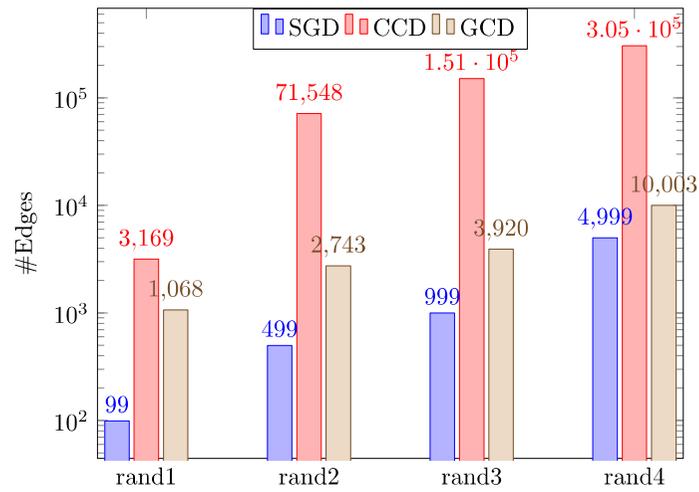


Experimental Results

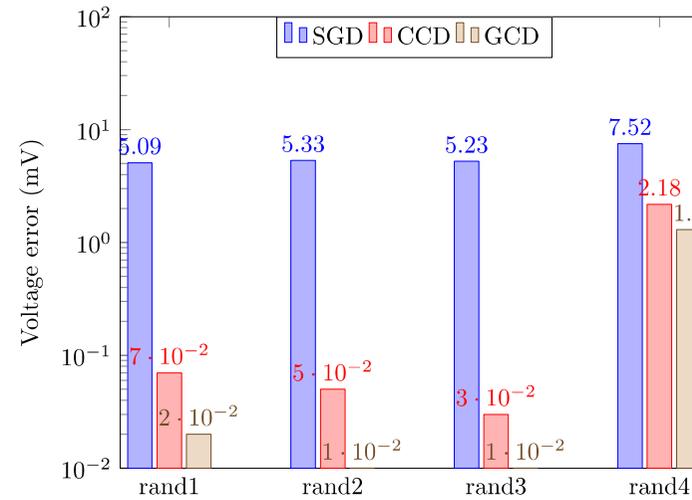
- ◆ Sparsity and accuracy trade-off
- ◆ Accuracy and runtime trade-off



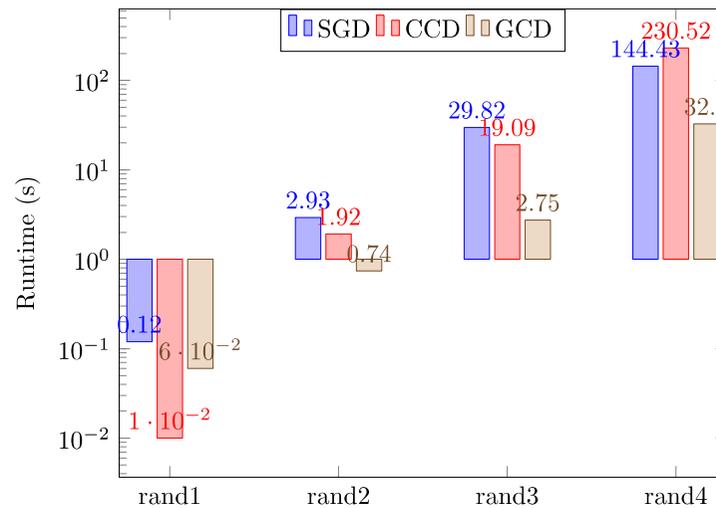
Gradient Descent Comparison



Sparsity



Accuracy



Runtime

Experimental Results

	CKT	ibmpg2	ibmpg3	ibmpg4	ibmpg5	ibmpg6
#Port Nodes	Before	19,173	100,988	133,622	270,577	380,991
	After	19,173	100,988	133,622	270,577	380,991
#Non-port Nodes	Before	46,265	340,088	345,122	311,072	481,675
	After	0	0	0	0	0
#Edges	Before	106,607	724,184	779,946	871,182	1283,371
	After	48,367	243,011	284,187	717,026	935,322
Error		1.2%	0.7%	4.8%	2.2%	2.0%
Runtime		38s	106s	132s	123s	281s

Conclusion

- ◆ Main Contributions:

- › An iterative power grid reduction framework
- › Weighted convex optimization-based formulation
- › A GCD algorithm with optimality guarantee and runtime efficiency for edge **sparsification**

- ◆ Future Work:

- › Extension to RC grid reduction

Thanks