



FastRW: An Efficient Random Walk Method for Steady-State Thermal Analysis

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Key Results

FastRW accelerates Feynman–Kac-based random walk thermal simulation through exact truncation error analysis and prior/posterior estimation, achieving **over 6.5× speedup** over prior state-of-the-art methods without sacrificing accuracy.

Motivation

Accurate thermal simulation is critical for modern IC design with increasing power density. Random walk methods based on the Feynman–Kac formula [4] enable efficient **localized** temperature estimation without solving the global field [5, 1, 3]. However:

- In practical scenarios **without Dirichlet boundaries**, long random walk paths are needed for convergence.
- Existing methods [2, 3] select truncation thresholds **empirically** with no principled error guarantee.
- **FastRW (ours)**: provides quantitative truncation error analysis and uses prior/posterior estimation to dramatically shorten paths.

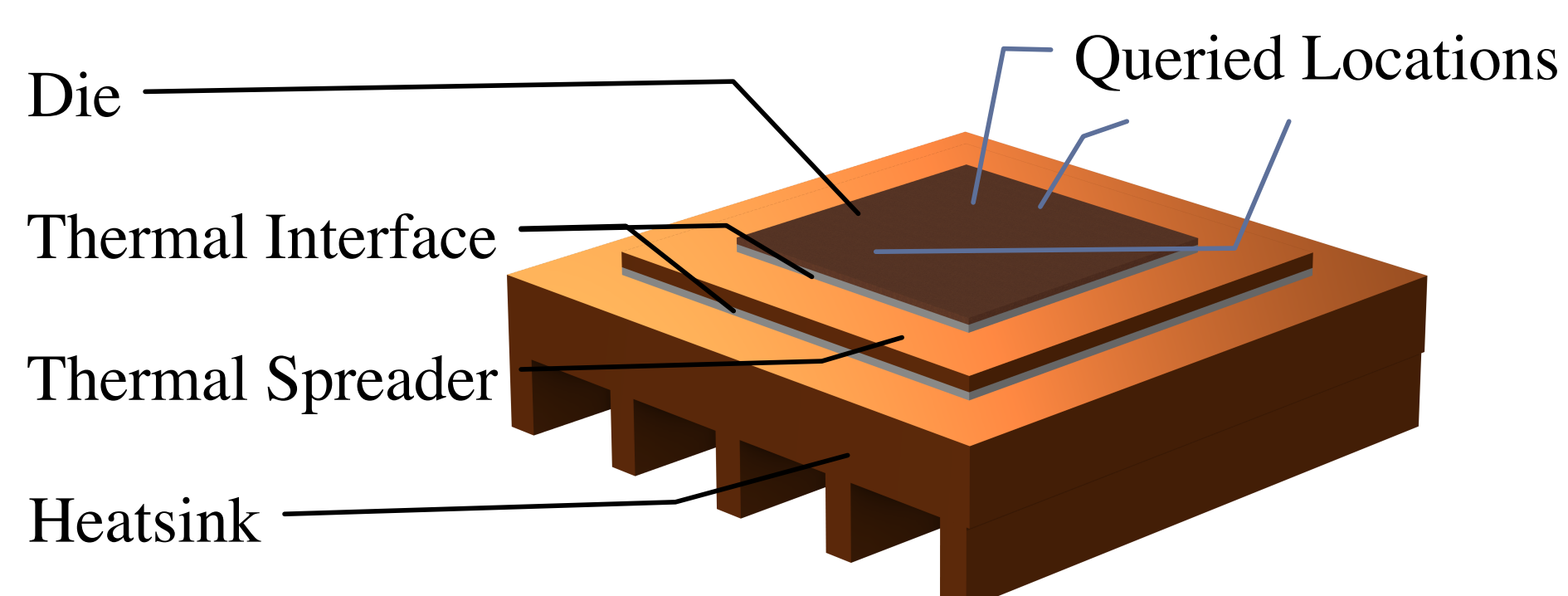


Figure 1. In many thermal design scenarios, only temperatures at a few queried hotspot locations are needed – ideal for random walk methods.

Feynman–Kac Representation

The temperature at \mathbf{x} is expressed via the Feynman–Kac formula as an expectation over reflecting Brownian motion (SRBM) paths:

$$T(\mathbf{x}) = \mathbb{E} \left\{ \int_0^\infty \hat{e}_c(t) f(X_t) dt \right\} + \mathbb{E} \left\{ \hat{e}_c(t_D) \phi_D(X_{t_D}) \right\} + \mathbb{E} \left\{ \int_0^\infty \hat{e}_c(t) \phi_N(X_t) dL(t) \right\} + \mathbb{E} \left\{ \int_0^\infty \hat{e}_c(t) \phi_R(X_t) dL(t) \right\}, \quad (1)$$

where $\hat{e}_c(t) = \exp\left(-\int_0^t c(X_s) dL(s)\right)$ and $L(t)$ is the local time at the boundary.

Truncation Error Analysis

For a threshold $\Lambda \in (0, 1)$, we truncate the path at t_Λ when $\hat{e}_c(t) \leq \Lambda$. The key result:

$$\mathbb{E} \left[\mathbb{P}_{t_\Lambda}^\infty \right] = \hat{e}_c(t_\Lambda) T(X_{t_\Lambda}) = \Lambda T(X_{t_\Lambda}). \quad (2)$$

Corollary: The expected truncation error equals the **weighted temperature** at the truncation location X_{t_Λ} . This provides a principled guideline for choosing Λ .

Single-Point Acceleration via Prior Estimation

We use a coarse FEM prior $\tilde{T}(\mathbf{x}) = T(\mathbf{x}) + \epsilon$ to compensate the truncation error:

$$T_i(\mathbf{x}) = \mathbb{P}_0^t + \hat{e}_c(t) \tilde{T}(\mathbf{x}) = \mathbb{P}_0^t + \mathbb{E} \left[\mathbb{P}_t^\infty \right] + \hat{e}_c(t) \epsilon. \quad (3)$$

Since $\epsilon \ll T(\mathbf{x})$, the threshold Λ can be **350× larger**, cutting path length by more than half.

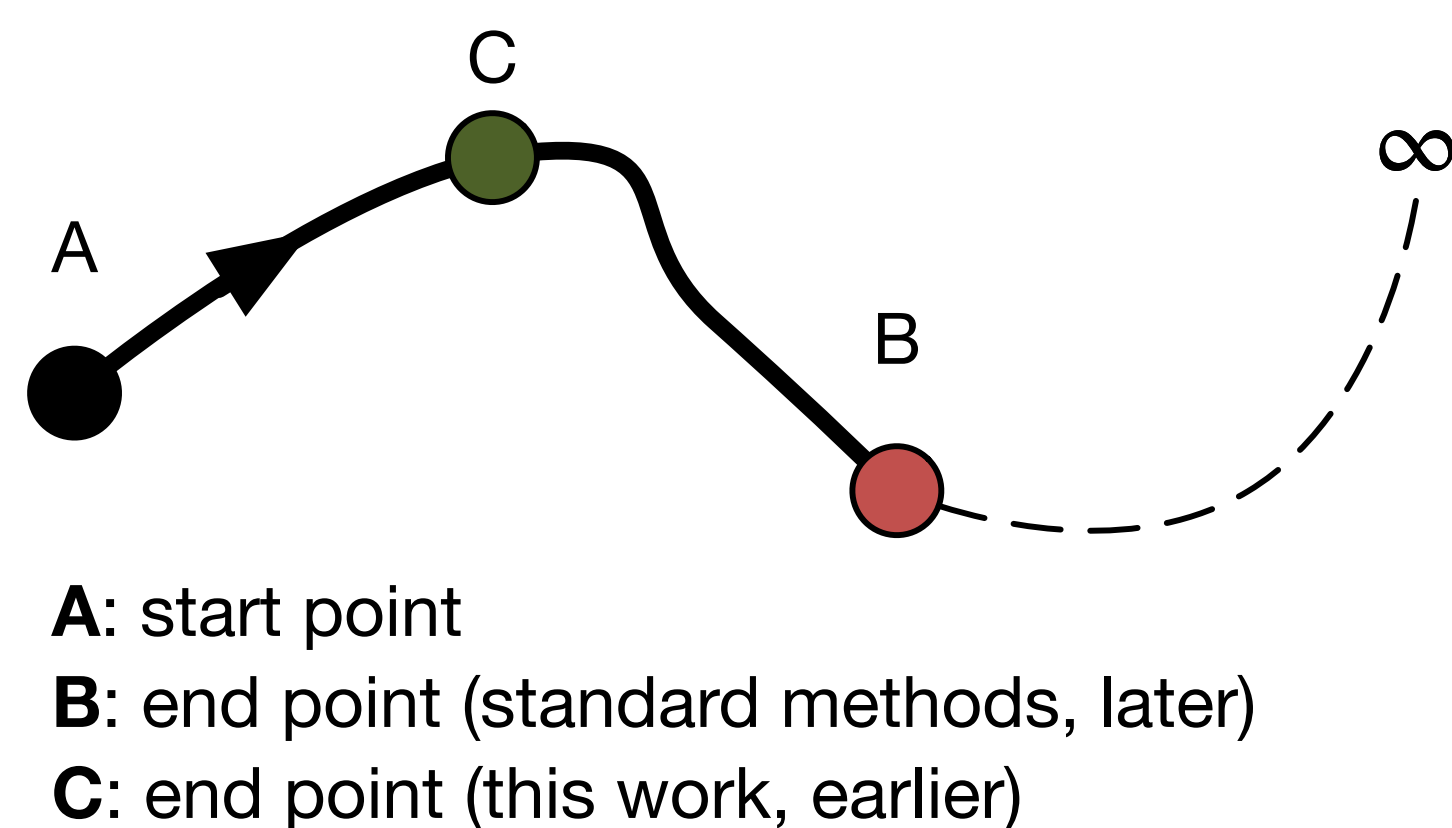


Figure 2. Standard methods compute the full path A→B. FastRW truncates at C and estimates the remaining contribution using the prior.

Multi-Point Acceleration via Posterior Observation

When estimating temperatures at M points $\{\mathbf{x}_q\}_{q=1}^M$, paths starting at \mathbf{x} that visit \mathbf{y} yield a noisy linear observation:

$$b_k = \mathbf{T}_{i_k} - \alpha_k \mathbf{T}_{j_k} + \eta_k, \quad \text{stacked as } \mathbf{B} = \mathbf{A} \mathbf{T} + \boldsymbol{\eta}. \quad (4)$$

With Bayesian inference (Gaussian prior from single-point estimates):

$$\hat{\mathbf{T}} = \boldsymbol{\Sigma}_{\text{post}} \left(\mathbf{A}^T \mathbf{R}^{-1} \mathbf{B} + \boldsymbol{\Sigma}_0^{-1} \boldsymbol{\mu}_0 \right), \quad (5)$$

cross-point information further reduces the number of paths needed.

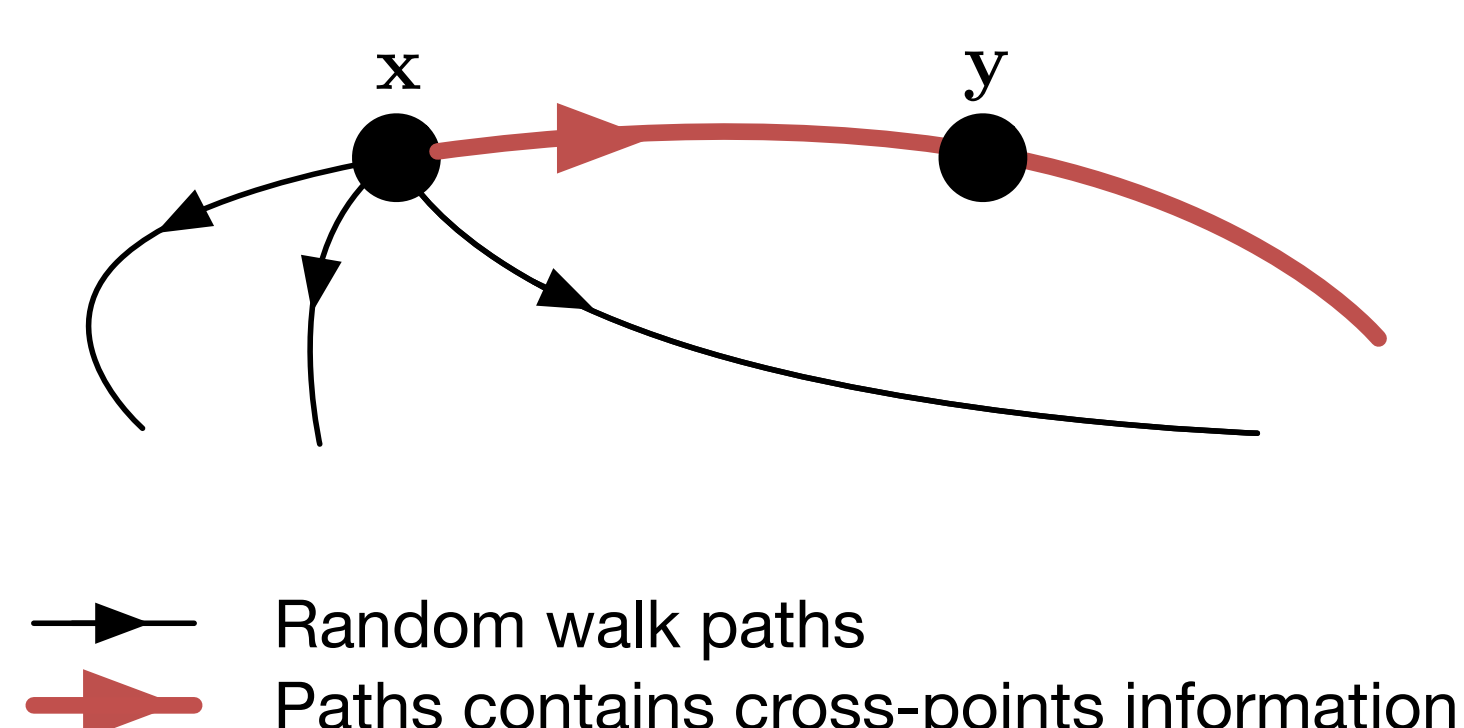


Figure 3. Paths may intersect other query points, providing cross-point linear observations for joint Bayesian refinement.

Experimental Setup

- **Test cases:** 3D chip model [5, 1] (3 layers: source-free / heat-source / source-free), 3 thickness configs, 3 power maps (**power6**, **4-core**, **16-core**).
- **Baseline:** PIRW [5], a Feynman–Kac method with standard truncation ($\Lambda=0.0001$, following [3]).
- **Hardware:** 2× Intel Xeon Silver 4214 (24 cores), 128 GB RAM.

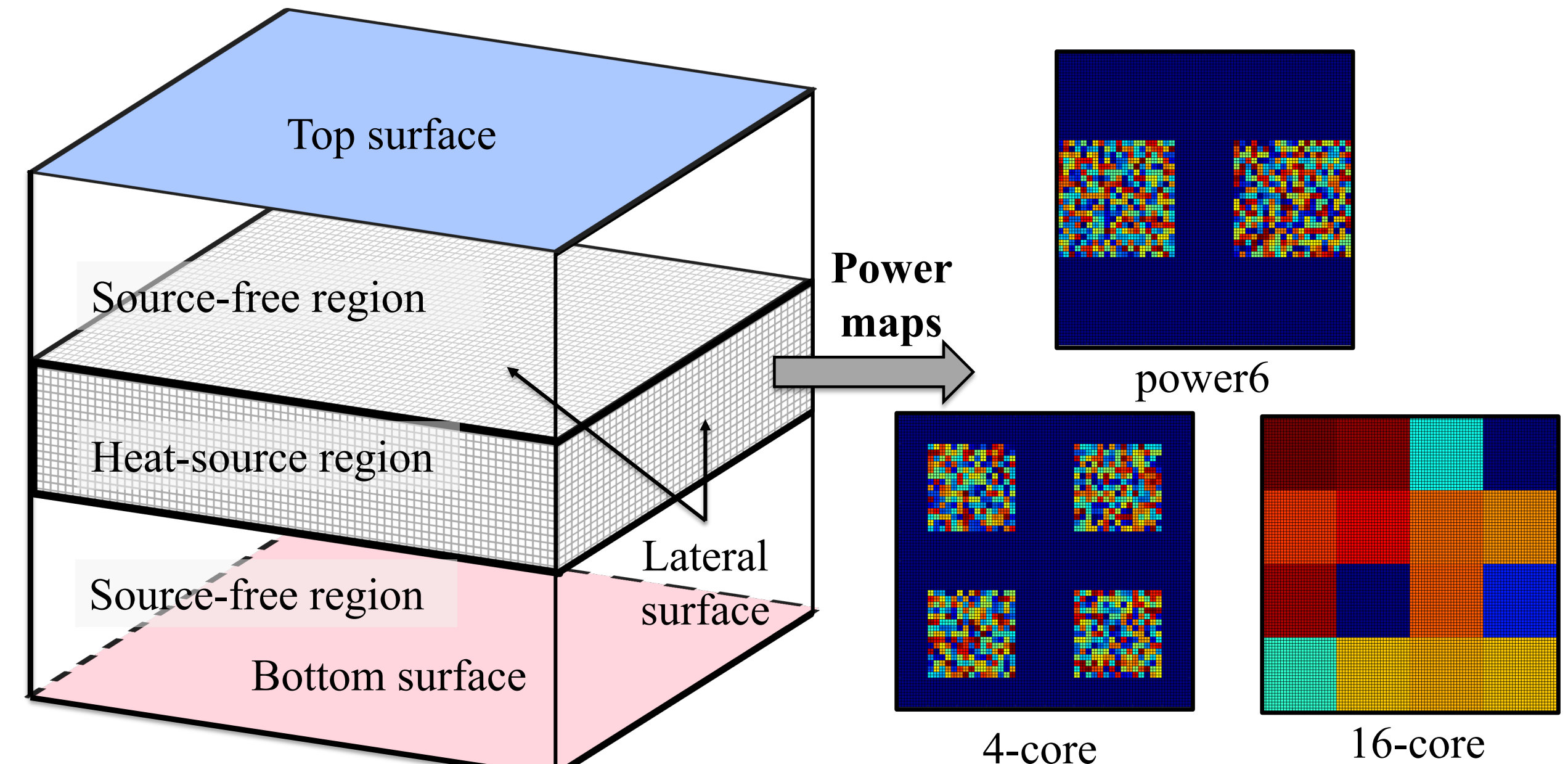


Figure 4. Test case geometry and power map configurations.

Single-Point Results

Table 1. Temperature estimation comparison. FastRW uses $\Lambda=0.03$ with coarse FEM prior.

Case	Power Map	PIRW [5]				FastRW				
		Error/K	StdDev/K	Steps	Paths	Error/K	StdDev/K	Steps	Paths	Speedup
1	power6	0.38	0.47	5.79e6	1000	0.36	0.45	2.20e6	400	6.58×
2	4-core	0.43	0.55	5.83e6	1000	0.42	0.54	2.22e6	400	6.57×
3	16-core	0.38	0.46	1.03e7	1000	0.32	0.41	3.93e6	400	6.55×

- FastRW achieves **6.5×+ speedup** with **improved accuracy** across all cases.
- Path length reduced by >50%; number of required paths reduced from 1000 to 400.

Multi-Point Acceleration

Table 2. Multi-point estimation (Case 1, power6).

Group Size	Error(K)	StdDev(K)	\hat{N}	Speedup
1	0.356	0.446	1000	1.00×
4	0.292	0.379	1385	1.39×
16	0.221	0.281	2519	2.52×

Table 3. Efficiency analysis: time per point (Case 1).

Group Size	Prior(s)	RW(s)	Posterior(s)	Extra Cost(%)
1	0.28	15.41	0.00	1.82
4	0.28	15.41	0.02	1.95
16	0.28	15.41	0.61	5.78

- Larger group sizes yield more cross-point observations \Rightarrow additional **2.52×** acceleration.
- Prior + posterior overhead is only **5.78%** even at group size 16.

Effect of Prior Accuracy

Table 4. Better FEM prior enables larger Λ and shorter paths.

DoF	Prior Max-Error (K)	Random Walk		
		Threshold Λ	Steps	Time (s)
1875	1.93	0.01	2.89e6	20.2
4800	0.88	0.02	2.45e6	17.1
12500	0.59	0.03	2.20e6	15.4

More accurate priors (finer FEM) reduce truncation error, allowing larger Λ and further runtime savings.

Conclusion

FastRW accelerates thermal analysis by:

1. **Exact truncation error analysis** – the expected error equals the weighted temperature at the truncation point.
2. **Prior estimation** – coarse FEM priors shorten paths by >50% and reduce variance, yielding **6.5×** overall single-point speedup.
3. **Posterior correction** – cross-point Bayesian inference provides up to **2.5×** additional acceleration with <6% overhead.

Future extensions: more accurate priors (learning-based), transient conduction, and GPU acceleration.

References

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