

An NISQ Scheme for Generating LABS

The Born Rule Breakers

February 1, 2026

Background: The Low Autocorrelation Binary Sequence (LABS) Problem

Objective: Minimize energy $E(x)$ for $x \in \{-1, +1\}^N$

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Current Algorithmic Landscape:

Method	Complexity	Key Constraint
Memetic Tabu Search (MTS)	$O(1.37^N)$	<i>Current SOTA</i>
QAOA (Standard) [1]	$O(1.46^N)$	$O(1.21^N)$ requires non-NISQ
Adiabatic and QE-MTS [2]	$O(1.24^N)$	Trotter depth / High precision

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The Research Challenge

To demonstrate **heuristic advantage**, we must develop a scalable, shallow-depth algorithm that outperforms the classical benchmark.

Preview of our Results

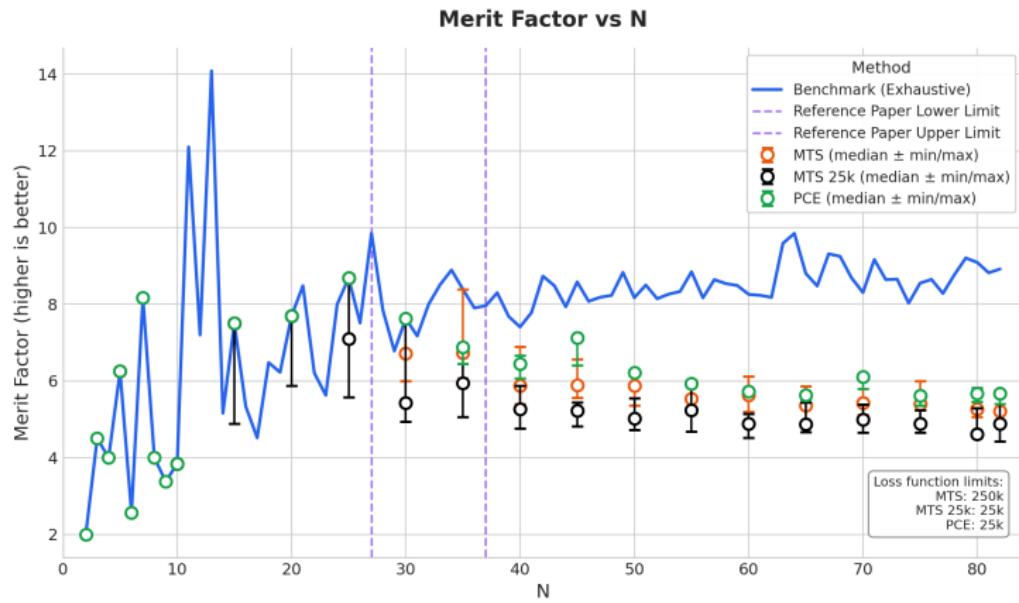


Figure: Merit Factor (MF) scaling across sequence sizes N . High MF indicates superior ground-state approximation for the LABS problem.

Our Approach

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- 1se **Optimization:** Update θ by minimizing the relaxed LABS loss:

$$L(\theta) = \sum_{\ell=1}^{N-1} \left(\sum_{i=1}^{N-\ell} \tilde{x}_i \tilde{x}_{i+\ell} \right)^2 - \beta \sum_{i=1}^N \tilde{x}_i^2$$

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- 1se **Decoding:** Upon convergence, use final expectations to assign the binary sequence: $x_i = \text{sign}(\langle \Pi_i \rangle)$.

Why PCE? Advantages for NISQ and Beyond

Variational approaches encode problems directly, while PCE maps variables to expectations of observables.

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It is a NISQ-friendly and scalable scheme, and PCE for the LABS problem demonstrates a runtime scaling of $\approx O(1.31^n)$, making it competitive with the best classical heuristics [3].

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Methodology and Benchmarking

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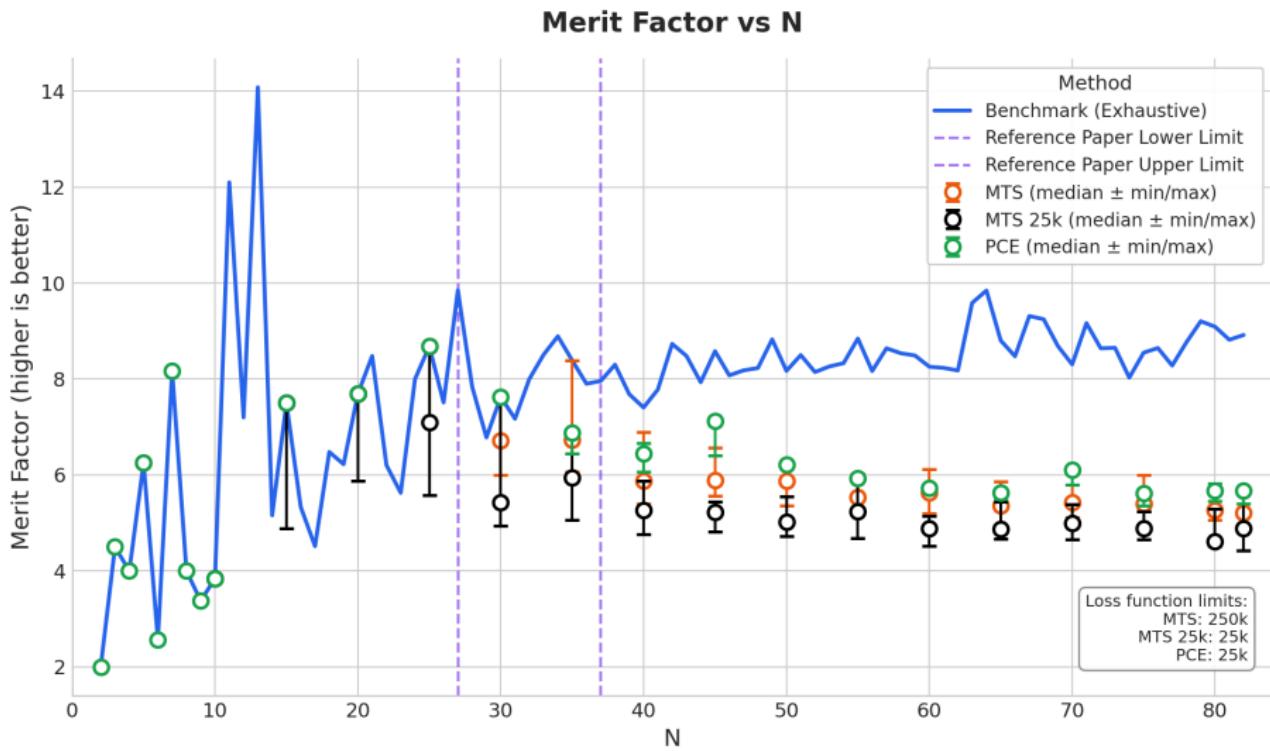
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Verification and Scaling

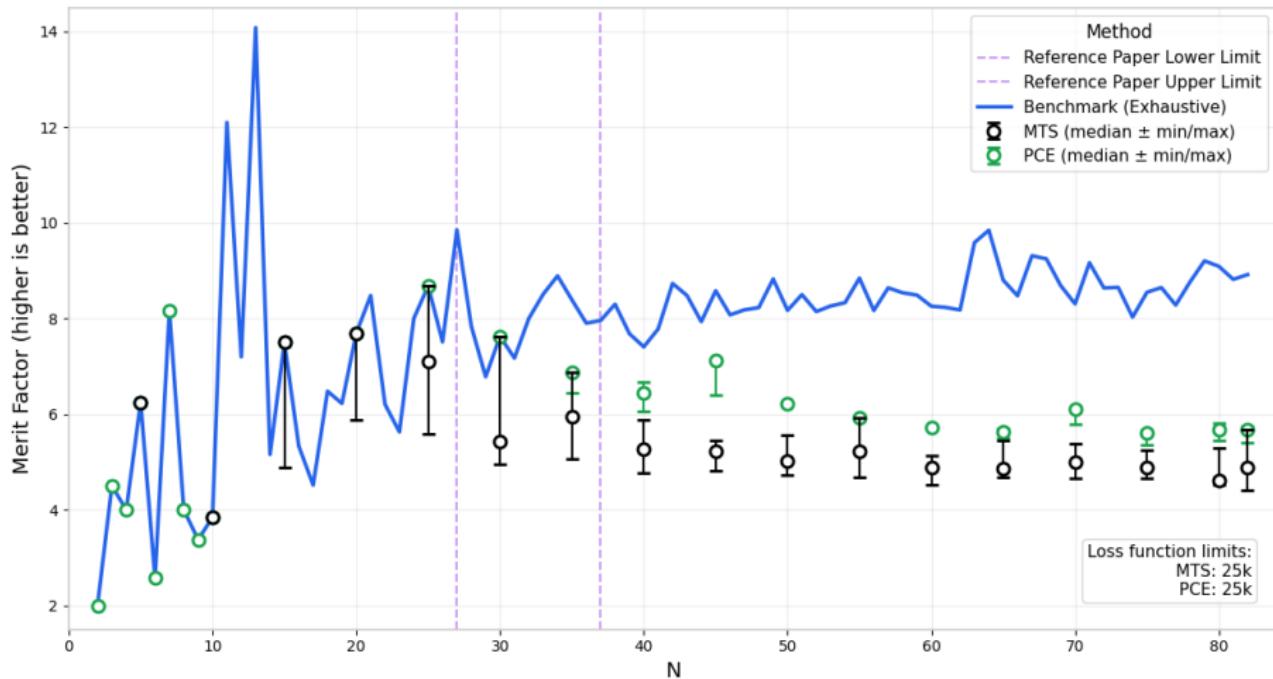
- **Validation:** MTS implementation verified against known optimal MF values up to $N = 82$. Each MTS was seeded and run 10 times.
- **Comparison:** Benchmarked PCE-MTS vs. MTS to identify where quantum seeding accelerates convergence toward global optima.

Benchmarking Results: Merit Factor vs. N



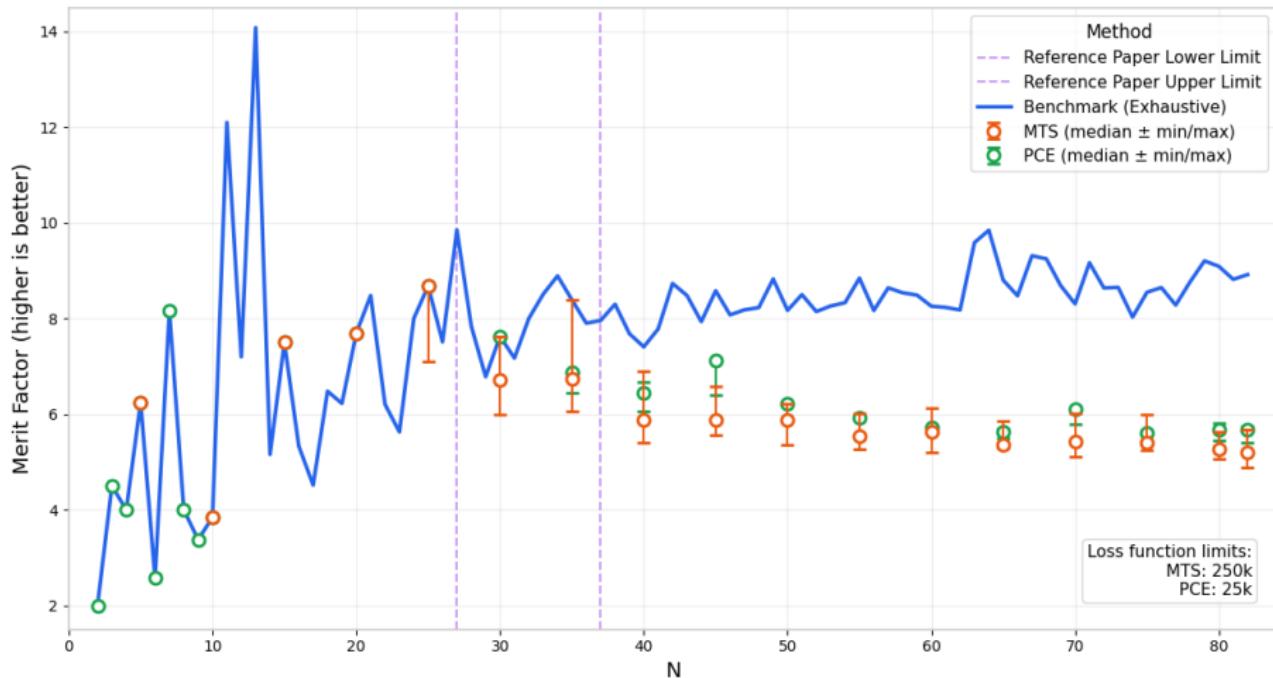
16.72% Improvement for 25k vs 25k LFCs

Merit Factor vs N



5.86% Improvement for 25k vs 250k LFCs

Merit Factor vs N



PCE Phase: Optimization of (Hyper)parameters

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The EGT-CG Approach

Optimizes on the **Fubini-Study metric/QGT** of the quantum parameter space [6]:

- **Scalability:** Reduces loss function calls for large sequences ($N \approx 50$).
- **Stability:** Suppresses **barren plateaus** via conjugate gradient (CG) method.
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Future improvement: could we tweak the encoding to better leverage quantum geometry?

GPU Acceleration: CUDA-Q & CuPy Integration

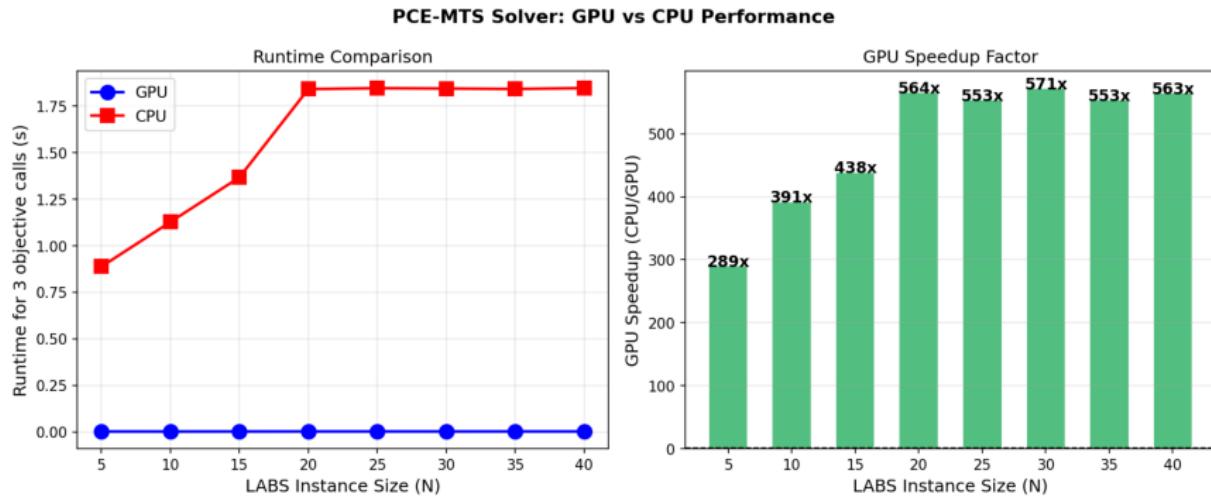


Figure: PCE-MTS Solver: GPU vs CPU Performance

GPU Acceleration: CUDA-Q & CuPy Integration

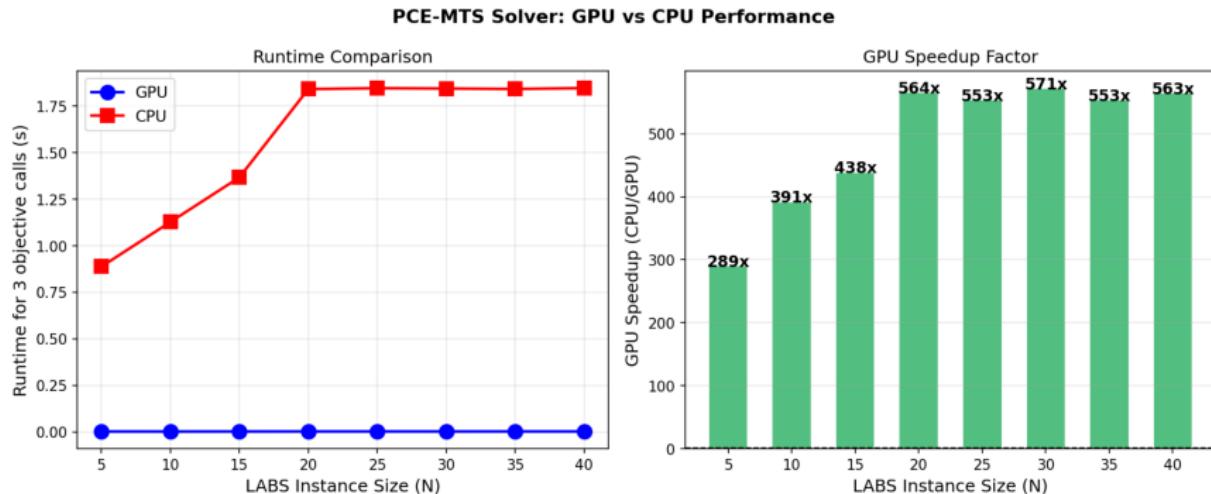


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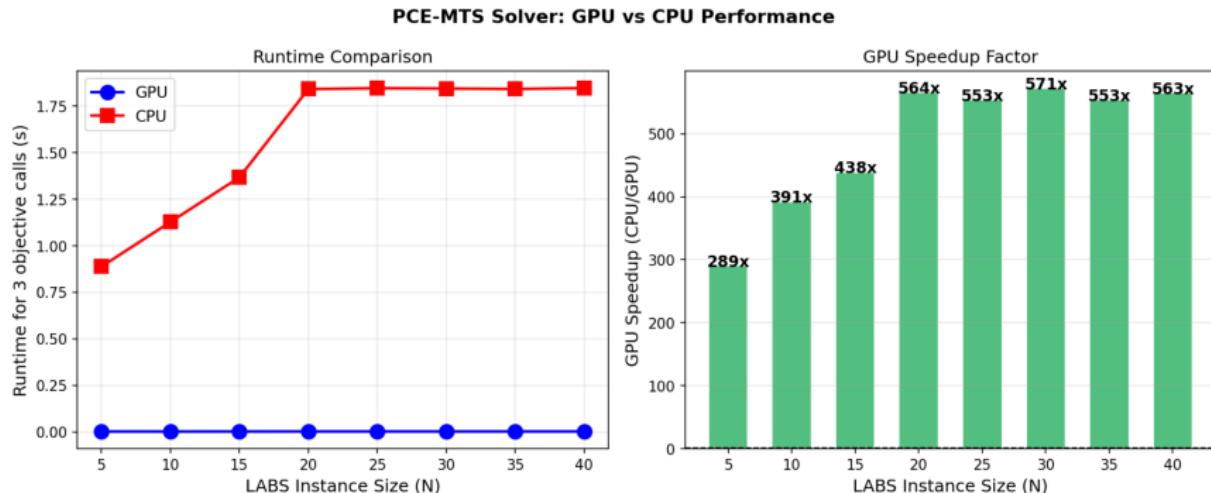


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- **End-to-End GPU Residency:** Statevectors (CUDA-Q) and MTS populations (CuPy) remain in GPU memory
- **Massive Parallelization:** Parallelizes parameter-shift gradients and energy evaluations via vectorized kernels and batched observables.

References I

- [1] M. Harrigan et al., *Evidence for scaling advantage of the quantum approximate optimization algorithm*, 2023. arXiv: 2308.02342 [quant-ph]. [Online]. Available: <https://arxiv.org/abs/2308.02342>.
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