

SUPER-RESOLUTION

FOR IMAGERY ENHANCEMENT USING VARIATIONAL QUANTUM EIGENSOLVER

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Super-Resolution for Imagery Enhancement Using Variational Quantum Eigensolver

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Agenda

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- 2. Related Method
- 3. Proposed Method
- 4. Experimental Results
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Introduction

Introduction

Super-Resolution



INPUT (1080p)

OUTPUT (4K)

Figure 1: Estimating a High-Resolution (HR) version of an image based on a Low-Resolution (LR) input is a complex task described as Single Image Super Resolution (SISR) [1].

- Face recognition;
- Surveillance systems;
- Medical imaging;
- Remote sensing;

- Astronomical images;
- Forensics;
- Multimedia industry and video enhancement.

- RAISR: Rapid and Accurate Image Super Resolution (ROMANO; ISIDORO; MILANFAR, 2017).
- VQE: Variational Quantum Eigensolver (PERUZZO et al., 2014).

Related Method

RAISR: Rapid and Accurate Image Super Resolution (ROMANO; ISIDORO; MILANFAR, 2017).

Global Filter Learning

$$\min_{\mathbf{h}} \sum_{i=1}^{L} \|\mathbf{A}_{i}\mathbf{h} - \mathbf{b}_{i}\|_{2}^{2}$$
(1)
$$\min_{\mathbf{h}} \|\mathbf{Q}\mathbf{h} - \mathbf{V}\|_{2}^{2}$$
(2)



Figure 2: Learning stage.

Related Method // Global Filter Learning



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Related Method // Global Filter Learning



Figure 4: Speed performance comparison for traditional interpolation algorithms at several scale factors when resizing an input image from Set5 [51]. Nearest-neighbor consistently presents the best time performance for all scaling factors.

Related Method // Hashing-based Learning



Figure 5: Visual representation for a single pixel type of a learned filter for a $2 \times$ upscaling factor.

Related Method // Local Gradient Statistics

- Local Gradient Statistics
 - Angle θ_k , Strength λ_1^k , and Coherence μ_k .

$$\mathbf{G}_{k} = \begin{bmatrix} g_{\mathbf{X}_{k_{1}}} & g_{\mathbf{y}_{k_{1}}} \\ \vdots & \vdots \\ g_{\mathbf{X}_{k_{n}}} & g_{\mathbf{y}_{k_{n}}} \end{bmatrix}$$
(3)

$$\mathbf{G}_{k}^{T}\mathbf{W}_{k}\mathbf{G}_{k} \tag{4}$$

$$\boldsymbol{\theta}_{k} = \arctan(\phi_{1,y}^{k}, \phi_{1,x}^{k}), \quad \sqrt{\boldsymbol{\lambda}_{1}^{k}}, \quad \boldsymbol{\mu}_{k} = \frac{\sqrt{\lambda_{1}^{k}} - \sqrt{\lambda_{2}^{k}}}{\sqrt{\lambda_{1}^{k}} + \sqrt{\lambda_{2}^{k}}}.$$
 (5)

Classical Filter

 Berkeley Segmentation Data Set and Benchmarks 500 (BSDS500) [50].

 \mathbf{h}_{lpha}

Proposed Method

VQE: Variational Quantum Eigensolver (PERUZZO et al., 2014).

Ground State Energy of the Hamiltonian H

$$\min_{\vec{\theta}} = \left\langle \psi(\vec{\theta}) \middle| \mathbf{H} \middle| \psi(\vec{\theta}) \right\rangle$$
(6)

Proposed Method // Variational Quantum Eigensolver



Figure 6: VQE Overview [19].

Numerical Reference

$$\textbf{H} = \begin{cases} (0,0) & 1.345\,149\,691\,971\,852\,6\times10^{-4} \\ (0,1) & 2.421\,287\,807\,653\,396\,7\times10^{-5} \\ (1,0) & 2.421\,287\,807\,653\,396\,7\times10^{-5} \\ (1,1) & 2.436\,686\,682\,060\,184\,6\times10^{-4} \end{cases}$$

• Exact Eigensolver (EE):

$$\mathbf{EE}\boldsymbol{\lambda}_{1}^{k} = 0.0002487985684589, \ \mathbf{EE}\boldsymbol{\phi}_{1}^{k} = \begin{bmatrix} -0.2072658478952684\\ -0.9782846560670648 \end{bmatrix}.$$
(8)
$$\mathbf{EE}\boldsymbol{\lambda}_{2}^{k} = 0.0001293850689443, \ \mathbf{EE}\boldsymbol{\phi}_{2}^{k} = \begin{bmatrix} -0.9782846560670648\\ 0.2072658478952684 \end{bmatrix}.$$
(9)

$$\min_{\vec{\theta}} = |\mathbf{H}\rangle_{|\psi(\vec{\theta})\rangle} \ge \lambda_{\min} \tag{10}$$

$$-\min_{\mathbf{x}} -f(x) = \max_{\mathbf{x}} f(x) \tag{11}$$

Proposed Method // Classical Optimization

- Classical Optimization:
 - ADAM;
 - Analytic Quantum Gradient Descent (AQGD);
 - Conjugate Gradient (CG) Method;
 - Constrained Optimization BY Linear Approximation (COBYLA);
 - Limited-memory Broyden-Fletcher-Goldfarb-Shanno Bound (L-BFGS-B);
 - Nelder-Mead (NM);
 - Sequential Least SQuares Programming (SLSQP);
 - Simultaneous Perturbation Stochastic Approximation (SPSA);
 - Truncated Newton (TNC).

Proposed Method // Classical Optimization



Figure 7: Energy convergence for local optimizers towards $\mathbf{EE}\lambda_k^2$.

Proposed Method // Classical Optimization



Figure 8: Divergence from reference value $\mathbf{EE}\lambda_k^2$.

COBYLA:

$$COBYLA\lambda_{max} = -0.00024879856832430445,$$
(12)
$$COBYLA\phi_{max} = \begin{bmatrix} -0.2072330078108220\\ -0.9782916132082906 \end{bmatrix}.$$
(13)

• Exact Eigensolver:

$$EE\lambda_{1}^{k} = 0.0002487985684589,$$
(14)
$$EE\phi_{1}^{k} = \begin{bmatrix} -0.2072658478952684\\ -0.9782846560670648 \end{bmatrix}.$$
(15)

COBYLA:

$$COBYLA\lambda_{min} = 0.00012938506911698078,$$
(16)
$$COBYLA\phi_{min} = \begin{bmatrix} 0.9782925363087142 \\ -0.2072286500526972 \end{bmatrix}.$$
(17)

• Exact Eigensolver:

$$EE\lambda_2^k = 0.0001293850689443,$$
(18)
$$EE\phi_2^k = \begin{bmatrix} -0.9782846560670648\\ 0.2072658478952684 \end{bmatrix}.$$
(19)

• Evolution Time (ET) in seconds for convergence to ground state considering 21 evaluations with COBYLA:

$$COBYLA\lambda_{max}ET = 4.131\,388\,664\,245\,605\,5\times10^{-2},\qquad(20)$$

$$COBYLA\lambda_{min}ET = 4.220\,700\,263\,977\,051\times10^{-2}.$$
 (21)

It took approximately 1 hour for all evaluations to be completed using a real device, with 3504 seconds for total Evaluation Time, considering 18 evaluations with COBYLA.

$$COBYLA\lambda_{min}^{\dagger} = 0.00013446128817024897.$$
 (22)

Numerical References:

$$COBYLA\lambda_{min} = 0.00012938506911698078.$$
 (23)
 $EE\lambda_2^k = 0.0001293850689443.$ (24)

Quantum-Classical Filter

Berkeley Segmentation Data Set and Benchmarks 500 (BSDS500) [50].

 $\mathbf{h}_{m{eta}}$

Experimental Results

Table 1: Quantitative comparison of the enhancement with several objective methods for the average performance result measured on the test images of Set5 [51].

Filter (2×)	MSE	PSNR	SSIM	VIF	RECO
Reference Values	0.00	100.00	1.00	1.00	1.00
\mathbf{h}_{α} (Classical)	102.85	33.55	0.904	0.91	1.07
$\mathbf{h}_{\boldsymbol{eta}}$ (Quantum-Classical)	102.85	33.55	0.904	0.91	1.07

Remarks and Conclusion

As with any nascent technology, there are limitations to the quantum computers and the simulators [24]. Time requirement to solve the complete operation using VQE could be decreased employing a generalization that reduces the number of samples [25] and the inherent parallelizability of the hybrid approach [26].

The classical optimization and the quantum computation are both parallelizable, with potential to overcome the more protracted runtime verified when comparing to the state-of-the-art, linearly improving per additional qubit. The concept of the VQE algorithm was first introduced in [2] with an implementation using integrated quantum photonics technology [27]. The VQE has the notable property that it can run on any quantum device [28], and it has already been experimentally demonstrated in other quantum platforms, including trapped ions [29] and superconducting circuits [30, 31].

This work benefits from the resources in the IBM Q Experience and the Quantum Information Science Kit – Qiskit [24], which conveniently includes the VQE algorithm.

Implementations of the VQE are also currently available in additional quantum computing and software platforms, such as in the ones offered by Rigetti Computing and Xanadu.

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