### PREDICTIONS OF QUBITS-CANDIDATES USING MACHINE LEARNING

by

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#### Abstract

This is an abstract. First coffee.

### Acknowledgements

Acknowledgements. Coffe-time?

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### Introduction

This it The introduction. Another coffee.

Part I

Theory

# Part II Implementation

Part III

Results

## Part IV Conclusion

# Part V Appendices

#### COFFEEEEEEEEEE appendix

### Bibliography

1. Balachandran, P. V. *et al.* Predictions of new ABO3 perovskite compounds by combining machine learning and density functional theory. *Physical Review Materials* **2.** doi:10.1103/physrevmaterials.2.043802 (Apr. 2018).