

Quantum Control Theory and Implementation

QubitPulseOpt: A Simulation Framework for Optimal Quantum Control of Two-Level Systems

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Abstract

This document provides a comprehensive theoretical foundation and implementation guide for the QubitPulseOpt quantum control simulation framework. We develop the mathematical formalism for controlling two-level quantum systems (qubits) using time-dependent electromagnetic pulses, building from first principles to advanced optimal control techniques. The treatment is organized by project phases, with each phase introducing new theoretical concepts alongside their computational implementations. Phase 1.1 establishes the computational environment and reproducibility infrastructure. Phase 1.2 develops the drift Hamiltonian formalism, analytical solutions for free evolution, and validation methodology. Subsequent phases will address control Hamiltonians, pulse shaping, optimal control algorithms, and relaxation/decoherence effects. This document serves as the authoritative reference for all physics, mathematics, and implementation decisions in the QubitPulseOpt framework.

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1 Introduction

1.1 Motivation and Scope

Quantum control theory addresses the fundamental challenge of manipulating quantum systems to achieve desired target states or operations with high fidelity. For two-level systems (qubits)—the basic unit of quantum information—this control is typically achieved through resonant or near-resonant electromagnetic pulses. The QubitPulseOpt framework provides a rigorous simulation environment for:

- (i) Modeling qubit dynamics under drift and control Hamiltonians
- (ii) Designing and optimizing control pulse shapes
- (iii) Analyzing decoherence and relaxation effects
- (iv) Implementing gradient-based optimal control algorithms (GRAPE, Krotov)
- (v) Validating control fidelity via numerical and analytical methods

This document is structured to mirror the phased development of the QubitPulseOpt codebase, with each section corresponding to a specific development phase. All theoretical results are accompanied by explicit mappings to implementation files, test suites, and demonstration notebooks.

1.2 Mathematical Conventions

Throughout this document, we adopt the following conventions:

- We work in units where $\hbar = 1$ unless explicitly stated otherwise.
- Quantum states are denoted by ket vectors $|\psi\rangle \in \mathcal{H}$, where \mathcal{H} is the Hilbert space (typically $\mathcal{H} = \mathbb{C}^2$ for a qubit).
- Observables and operators are represented by calligraphic or bold letters (e.g., \hat{H} , $\hat{\rho}$).
- Time-dependent quantities are explicitly written as functions of t .
- The Pauli matrices are defined as:

$$\sigma_x = \begin{pmatrix} 0 & 1 \\ 1 & 0 \end{pmatrix}, \quad \sigma_y = \begin{pmatrix} 0 & -i \\ i & 0 \end{pmatrix}, \quad \sigma_z = \begin{pmatrix} 1 & 0 \\ 0 & -1 \end{pmatrix} \quad (1)$$

- The identity operator on \mathbb{C}^2 is denoted \mathbb{I} .

2 Phase 1.1: Computational Infrastructure and Reproducibility

2.1 Overview

Before developing quantum control algorithms, we must establish a robust, reproducible computational environment. Phase 1.1 addresses the foundational infrastructure requirements: version control, dependency management, environment isolation, and validation protocols. This phase ensures that all subsequent theoretical developments can be implemented, tested, and reproduced by independent researchers or on different computational platforms.

2.2 Environment Design Philosophy

2.2.1 Reproducibility Requirements

Scientific computing demands bitwise-reproducible results across different machines and time periods. For the QubitPulseOpt framework, reproducibility is achieved through:

1. **Environment Isolation:** All Python dependencies are installed in an isolated virtual environment, preventing version conflicts with system packages or other projects.
2. **Explicit Dependency Specification:** Exact versions of all packages are documented in `environment.yml` (for Conda) and can be captured via `pip freeze` for venv users.
3. **Version Control:** All source code, documentation, and configuration files are tracked via Git and hosted on GitHub (<https://github.com/rylanmalarchick/QubitPulseOpt>).
4. **Automated Validation:** Setup validation scripts ensure that the environment is correctly configured before scientific work begins.

2.2.2 Virtual Environment vs. Conda

Two primary approaches exist for Python environment management:

venv (Python’s built-in virtual environment): Lightweight, no external dependencies, straightforward activation. Used as the primary environment for QubitPulseOpt.

- **Pros:** Ships with Python 3.3+, minimal overhead, integrates seamlessly with pip.
- **Cons:** Does not manage non-Python dependencies (e.g., BLAS, LAPACK for NumPy/SciPy optimization).

conda (Anaconda/Miniconda): Cross-language package manager, handles compiled dependencies.

- **Pros:** Manages both Python and system-level libraries (e.g., MKL-optimized NumPy), good for complex scientific stacks.
- **Cons:** Larger installation footprint, potential conflicts between conda and pip packages.

Decision: QubitPulseOpt uses `venv` as the default environment, with `environment.yml` provided as an alternative for Conda users. This decision balances simplicity (`venv` is universally available) with flexibility (Conda users can replicate the environment exactly).

2.3 Core Dependencies

The QubitPulseOpt framework relies on the following Python packages:

2.3.1 QuTiP Installation Notes

QuTiP (Quantum Toolbox in Python) is the central dependency. Version 5.x introduces significant API changes compared to 4.x:

- **Solver Interface:** The `sesolve` function (Schrödinger equation solver) now returns a `Result` object with state trajectories accessed via `result.states`.
- **Bloch Sphere Plotting:** The `Bloch.add_points` method requires the `meth` parameter to be one of `{‘s’, ‘l’, ‘m’}` (single point, line, multi-point). Colors are set via `point_color`, `point_marker`, etc., rather than passed directly to `add_points`.

Package	Version	Purpose
QuTiP	5.2.1	Quantum Toolbox in Python; provides quantum state/operator representations, Hamiltonian evolution solvers (Schrödinger, Lindblad), and Bloch sphere visualization.
NumPy	2.3.4	Numerical array operations, linear algebra, Fourier transforms.
SciPy	1.16.2	Advanced scientific computing: ODE solvers, optimization routines, special functions.
Matplotlib	3.10.7	Plotting and visualization of quantum dynamics, control pulses, and convergence metrics.
Jupyter	(latest)	Interactive notebook environment for demonstrations and exploratory analysis.
pytest	(latest)	Unit testing framework; ensures correctness of all modules.
pytest-cov	(latest)	Code coverage analysis for test suites.

Table 1: Core dependencies for QubitPulseOpt with version information and usage descriptions.

- **Quantum Objects:** States and operators are represented as `Qobj` instances. Hermiticity, normalization, and dimensionality are automatically validated.

These API details are critical for Phase 1.2 implementations (drift Hamiltonian evolution) and later phases (control pulse simulations).

2.4 Repository Structure

The QubitPulseOpt repository follows a modular, hierarchical structure:

```

quantumControls/
  src/
    __init__.py
    hamiltonian/
      __init__.py
      drift.py           # Drift Hamiltonian (Phase 1.2)
      evolution.py       # Time evolution engine (Phase 1.2)
      control.py         # Control Hamiltonian (Phase 1.3, future)
    pulses/              # Pulse shapes (future)
    optimization/        # GRAPE, Krotov (future)
    visualization/       # Plotting utilities (future)
  tests/
    unit/
      test_drift.py      # Unit tests for drift.py
      ...
    integration/         # End-to-end tests (future)
  notebooks/
    01_drift_dynamics.ipynb  # Phase 1.2 demonstration
    ...

```

```

scripts/
  validate_setup.sh      # Environment validation
  activate_env.sh        # Activate venv helper
  verify_drift_evolution.py # Standalone drift test
docs/
  science/
    quantum_control_theory.tex # This document
  SETUP_COMPLETE.md
  REVIEW_SUMMARY.md
environment.yml          # Conda environment spec
README.md
.gitignore

```

This structure separates concerns: `src/` contains all production code, `tests/` validates correctness, `notebooks/` provides interactive demonstrations, and `docs/` maintains theoretical documentation.

2.5 Validation Protocol

To ensure the environment is correctly configured, we implement a multi-stage validation protocol:

- Stage 1: Python Version Check:** Verify Python ≥ 3.8 (required for QuTiP 5.x and modern NumPy).
- Stage 2: Package Import Test:** Attempt to import all core packages (QuTiP, NumPy, SciPy, Matplotlib) and verify versions.
- Stage 3: QuTiP Functionality Test:** Create a simple quantum state $|0\rangle$, apply a Pauli-X gate, and verify $\sigma_x |0\rangle = |1\rangle$ to within numerical precision.
- Stage 4: Numerical Precision Test:** Confirm that NumPy/SciPy linear algebra operations achieve machine precision ($\sim 10^{-16}$ for double-precision floats).

The validation script `scripts/validate_setup.sh` orchestrates these checks. A successful run outputs:

```

[PASS] Python version: 3.12.3
[PASS] QuTiP 5.2.1 imported successfully
[PASS] NumPy 2.3.4 imported successfully
[PASS] SciPy 1.16.2 imported successfully
[PASS] Matplotlib 3.10.7 imported successfully
[PASS] Basic QuTiP test:  $|0\rangle \rightarrow |1\rangle$  via Pauli-X
[PASS] Numerical precision:  $||I - I|| = 0.0$ 
All validation checks passed.

```

2.6 Git Workflow and Version Control

The project uses Git for version control with the following practices:

- **Branch Strategy:** Development occurs on the `main` branch initially; feature branches will be used for major new components (e.g., `feature/grape-optimizer`).
- **Commit Messages:** Follow conventional commit format: `type(scope): description`, e.g., `feat(drift): implement analytical propagator for H0`.

- **Remote Repository:** Hosted at <https://github.com/rylanmalarchick/QubitPulseOpt>. All development is pushed regularly to ensure cloud backup and collaboration readiness.

2.7 Summary of Phase 1.1 Deliverables

- Git repository initialized and pushed to GitHub.
- Virtual environment created with all core dependencies installed.
- Repository structure established (src/, tests/, notebooks/, docs/).
- Validation scripts implemented and passing.
- Documentation of environment setup in README.md and SETUP_COMPLETE.md.

Code Mapping: Configuration files (environment.yml, .gitignore), validation scripts (scripts/validate_setup.sh, scripts/test_env_simple.py), and documentation (docs/SETUP_COMPLETE.m

3 Phase 1.2: Drift Hamiltonian and Free Evolution

3.1 Overview

Phase 1.2 establishes the theoretical and computational framework for the *drift Hamiltonian* \hat{H}_0 , which governs the natural evolution of a qubit in the absence of external control fields. Understanding drift dynamics is prerequisite to designing control pulses, as optimal control effectively steers the system away from its natural trajectory toward a desired target state.

We develop the drift Hamiltonian for a qubit in a static magnetic field, derive analytical solutions for time evolution, implement numerical solvers for validation, and demonstrate the equivalence of analytical and numerical methods to machine precision.

3.2 Two-Level Systems and the Qubit Hilbert Space

Definition 3.1 (Qubit). A *qubit* is a two-level quantum system whose state space is the two-dimensional complex Hilbert space $\mathcal{H} = \mathbb{C}^2$. The computational basis states are:

$$|0\rangle = \begin{pmatrix} 1 \\ 0 \end{pmatrix}, \quad |1\rangle = \begin{pmatrix} 0 \\ 1 \end{pmatrix} \quad (2)$$

Any pure state $|\psi\rangle \in \mathcal{H}$ can be written as:

$$|\psi\rangle = \alpha |0\rangle + \beta |1\rangle, \quad \alpha, \beta \in \mathbb{C}, \quad |\alpha|^2 + |\beta|^2 = 1 \quad (3)$$

The Pauli matrices $\{\sigma_x, \sigma_y, \sigma_z\}$ form a basis for traceless Hermitian operators on \mathcal{H} . Together with the identity \mathbb{I} , they span the space of all 2×2 Hermitian matrices. Any Hermitian operator (including Hamiltonians) can be decomposed as:

$$\hat{H} = c_0 \mathbb{I} + c_x \sigma_x + c_y \sigma_y + c_z \sigma_z, \quad c_i \in \mathbb{R} \quad (4)$$

3.3 Physical Derivation of the Drift Hamiltonian

Consider a qubit (e.g., a spin-1/2 particle or a two-level atom) placed in a static magnetic field $\vec{B} = B_0 \hat{z}$ aligned along the z -axis. The interaction Hamiltonian between the qubit's magnetic moment $\vec{\mu}$ and the field is:

$$\hat{H}_0 = -\vec{\mu} \cdot \vec{B} \quad (5)$$

For a spin-1/2 particle with gyromagnetic ratio γ , the magnetic moment operator is $\vec{\mu} = \gamma\vec{S}$, where $\vec{S} = \frac{\hbar}{2}\vec{\sigma}$ is the spin operator. Thus:

$$\hat{H}_0 = -\gamma B_0 \frac{\hbar}{2} \sigma_z = -\frac{\omega_0}{2} \hbar \sigma_z \quad (6)$$

where $\omega_0 = \gamma B_0$ is the *Larmor frequency* (or qubit transition frequency). In units where $\hbar = 1$:

$$\hat{H}_0 = -\frac{\omega_0}{2} \sigma_z = -\frac{\omega_0}{2} \begin{pmatrix} 1 & 0 \\ 0 & -1 \end{pmatrix} = \begin{pmatrix} -\omega_0/2 & 0 \\ 0 & \omega_0/2 \end{pmatrix} \quad (7)$$

Remark 3.1. The factor of 1/2 is conventional and arises from the normalization of the Pauli matrices. Some references define $\hat{H}_0 = \omega_0 |1\rangle\langle 1|$, which differs by an additive constant ($\omega_0 \mathbb{I}/2$). Since global energy shifts do not affect dynamics, both conventions are equivalent up to a redefinition of the zero-point energy.

3.4 Spectral Properties of the Drift Hamiltonian

The drift Hamiltonian \hat{H}_0 is diagonal in the computational basis, so its eigenanalysis is trivial:

Theorem 3.1 (Eigenspectrum of \hat{H}_0). The drift Hamiltonian $\hat{H}_0 = -(\omega_0/2)\sigma_z$ has:

- (i) Eigenvalues: $E_0 = -\omega_0/2$ and $E_1 = +\omega_0/2$
- (ii) Eigenstates: $|0\rangle$ (ground state, energy E_0) and $|1\rangle$ (excited state, energy E_1)
- (iii) Energy gap: $\Delta E = E_1 - E_0 = \omega_0$

The energy gap $\Delta E = \omega_0$ determines the qubit's transition frequency. In free evolution, the qubit oscillates at this frequency between $|0\rangle$ and $|1\rangle$ components (modulo the phase difference). The period of oscillation is:

$$T = \frac{2\pi}{\omega_0} \quad (8)$$

3.5 Analytical Solution for Time Evolution

3.5.1 Time Evolution Operator

The Schrödinger equation for a time-independent Hamiltonian \hat{H}_0 is:

$$i \frac{d}{dt} |\psi(t)\rangle = \hat{H}_0 |\psi(t)\rangle \quad (9)$$

The formal solution is:

$$|\psi(t)\rangle = \hat{U}(t) |\psi(0)\rangle, \quad \hat{U}(t) = e^{-i\hat{H}_0 t} \quad (10)$$

where $\hat{U}(t)$ is the unitary time evolution operator.

For $\hat{H}_0 = -(\omega_0/2)\sigma_z$, we compute the matrix exponential explicitly:

Proposition 3.2 (Propagator for Drift Hamiltonian).

$$\hat{U}(t) = \exp\left(i \frac{\omega_0 t}{2} \sigma_z\right) = \cos\left(\frac{\omega_0 t}{2}\right) \mathbb{I} + i \sin\left(\frac{\omega_0 t}{2}\right) \sigma_z \quad (11)$$

In matrix form:

$$\hat{U}(t) = \begin{pmatrix} e^{i\omega_0 t/2} & 0 \\ 0 & e^{-i\omega_0 t/2} \end{pmatrix} \quad (12)$$

Proof. Since σ_z is diagonal, $\sigma_z^{2n} = \mathbb{I}$ and $\sigma_z^{2n+1} = \sigma_z$. The Taylor series for the exponential becomes:

$$e^{i(\omega_0 t/2)\sigma_z} = \sum_{n=0}^{\infty} \frac{1}{n!} \left(i \frac{\omega_0 t}{2} \right)^n \sigma_z^n \quad (13)$$

$$= \sum_{n=0}^{\infty} \frac{(-1)^n}{(2n)!} \left(\frac{\omega_0 t}{2} \right)^{2n} \mathbb{I} + i \sum_{n=0}^{\infty} \frac{(-1)^n}{(2n+1)!} \left(\frac{\omega_0 t}{2} \right)^{2n+1} \sigma_z \quad (14)$$

$$= \cos \left(\frac{\omega_0 t}{2} \right) \mathbb{I} + i \sin \left(\frac{\omega_0 t}{2} \right) \sigma_z \quad (15)$$

Alternatively, diagonalizing yields (12) directly. \square

3.5.2 Evolution of Basis States

Applying $\hat{U}(t)$ to the computational basis states:

$$\hat{U}(t) |0\rangle = e^{i\omega_0 t/2} |0\rangle \quad (16)$$

$$\hat{U}(t) |1\rangle = e^{-i\omega_0 t/2} |1\rangle \quad (17)$$

Each eigenstate acquires a phase at rate $\pm\omega_0/2$. For a general initial state $|\psi(0)\rangle = \alpha |0\rangle + \beta |1\rangle$:

$$|\psi(t)\rangle = \alpha e^{i\omega_0 t/2} |0\rangle + \beta e^{-i\omega_0 t/2} |1\rangle \quad (18)$$

3.6 Bloch Sphere Representation

3.6.1 Bloch Vector Mapping

Any qubit state $|\psi\rangle$ can be represented as a point on or inside the Bloch sphere. For pure states:

$$|\psi\rangle = \cos(\theta/2) |0\rangle + e^{i\phi} \sin(\theta/2) |1\rangle \quad (19)$$

corresponds to the Bloch vector:

$$\vec{r} = (\sin \theta \cos \phi, \sin \theta \sin \phi, \cos \theta) \quad (20)$$

The Bloch components are computed from expectation values:

$$r_x = \langle \psi | \sigma_x | \psi \rangle = 2 \operatorname{Re}(\alpha^* \beta) \quad (21)$$

$$r_y = \langle \psi | \sigma_y | \psi \rangle = 2 \operatorname{Im}(\alpha^* \beta) \quad (22)$$

$$r_z = \langle \psi | \sigma_z | \psi \rangle = |\alpha|^2 - |\beta|^2 \quad (23)$$

3.6.2 Drift Dynamics on the Bloch Sphere

Under $\hat{H}_0 = -(\omega_0/2)\sigma_z$, the Bloch vector precesses about the z -axis at angular frequency ω_0 . From (18), for initial state $|\psi(0)\rangle = \alpha |0\rangle + \beta |1\rangle$:

$$r_x(t) = 2 \operatorname{Re}(\alpha^* \beta e^{-i\omega_0 t}) = r_x(0) \cos(\omega_0 t) + r_y(0) \sin(\omega_0 t) \quad (24)$$

$$r_y(t) = 2 \operatorname{Im}(\alpha^* \beta e^{-i\omega_0 t}) = -r_x(0) \sin(\omega_0 t) + r_y(0) \cos(\omega_0 t) \quad (25)$$

$$r_z(t) = |\alpha|^2 - |\beta|^2 = r_z(0) \quad (26)$$

Theorem 3.3 (Bloch Precession for Drift Hamiltonian). The Bloch vector $(r_x(t), r_y(t), r_z(t))$ under \hat{H}_0 rotates about the z -axis at angular frequency ω_0 , with r_z constant. The trajectory is a circle of radius $\sqrt{r_x(0)^2 + r_y(0)^2}$ at height $r_z(0)$.

3.7 Numerical Implementation

3.7.1 DriftHamiltonian Class

The drift Hamiltonian is implemented in `src/hamiltonian/drift.py` as a Python class:

```

1 import qutip as qt
2 import numpy as np
3
4 class DriftHamiltonian:
5     """
6     Represents the drift Hamiltonian  $H_0 = -\omega_0/2 * \sigma_z$ 
7     for a two-level quantum system.
8     """
9     def __init__(self, omega_0: float):
10         """
11         Parameters:
12         -----
13         omega_0 : float
14             Qubit transition frequency (Larmor frequency)
15         """
16         self.omega_0 = omega_0
17         self._hamiltonian = -(omega_0 / 2.0) * qt.sigmaz()
18
19     def hamiltonian(self) -> qt.Qobj:
20         """Returns the drift Hamiltonian as a QuTiP Qobj."""
21         return self._hamiltonian
22
23     def eigenvalues(self) -> np.ndarray:
24         """Returns eigenvalues  $[E_0, E_1]$  sorted ascending."""
25         return np.array([-self.omega_0/2, self.omega_0/2])
26
27     def eigenstates(self) -> tuple:
28         """Returns (eigenvalues, eigenstates) as QuTiP arrays."""
29         return self._hamiltonian.eigenstates()
30
31     def period(self) -> float:
32         """Returns the precession period  $T = 2\pi/\omega_0$ ."""
33         return 2 * np.pi / self.omega_0
34
35     def evolve_state(self, psi0: qt.Qobj, t: float) -> qt.Qobj:
36         """
37         Evolves initial state psi0 to time t using analytical
38         propagator.
39         """
40         #  $U(t) = \exp(i * \omega_0 * t / 2 * \sigma_z)$ 
41         propagator = (np.cos(self.omega_0 * t / 2) * qt.qeye(2)
42                     + 1j * np.sin(self.omega_0 * t / 2) * qt.sigmaz()
43                     )
44         return propagator * psi0

```

Listing 1: Drift Hamiltonian class structure

3.7.2 TimeEvolution Engine

For validation and comparison with numerical methods, we implement a TimeEvolution class in `src/hamiltonian/evolution.py`:

```

1 import qutip as qt

```

```

2 import numpy as np
3
4 class TimeEvolution:
5     """
6     Utilities for time evolution of quantum states under a Hamiltonian.
7     """
8     @staticmethod
9     def evolve_numerical(H: qt.Qobj, psi0: qt.Qobj,
10                        times: np.ndarray) -> qt.solver.Result:
11         """
12         Numerically evolve psi0 under Hamiltonian H using QuTiP's
13         sesolve.
14
15         Returns:
16         -----
17         result : qutip.solver.Result
18             Contains result.states (list of states at each time),
19             result.times (time array).
20         """
21         result = qt.sesolve(H, psi0, times)
22         return result
23
24     @staticmethod
25     def evolve_analytical(H_drift, psi0: qt.Qobj,
26                        times: np.ndarray) -> list:
27         """
28         Analytically evolve psi0 under DriftHamiltonian H_drift.
29
30         Returns:
31         -----
32         states : list of qt.Qobj
33             State at each time in times array.
34         """
35         states = [H_drift.evolve_state(psi0, t) for t in times]
36         return states
37
38     @staticmethod
39     def bloch_coordinates(psi: qt.Qobj) -> tuple:
40         """
41         Compute Bloch sphere coordinates (x, y, z) for state psi.
42         """
43         sx = qt.expect(qt.sigmax(), psi)
44         sy = qt.expect(qt.sigmay(), psi)
45         sz = qt.expect(qt.sigmaz(), psi)
46         return (sx, sy, sz)
47
48     @staticmethod
49     def bloch_trajectory(states: list) -> tuple:
50         """
51         Compute full Bloch trajectory from list of states.
52
53         Returns:
54         -----
55         (x_vals, y_vals, z_vals) : tuple of np.ndarray
56         """
57         coords = [TimeEvolution.bloch_coordinates(psi)
58                   for psi in states]
59         x_vals = np.array([c[0] for c in coords])

```

```

59     y_vals = np.array([c[1] for c in coords])
60     z_vals = np.array([c[2] for c in coords])
61     return (x_vals, y_vals, z_vals)

```

Listing 2: Time evolution engine

3.8 Validation and Testing

3.8.1 Unit Test Coverage

Comprehensive unit tests are implemented in `tests/unit/test_drift.py`. Key test categories:

Test 1: Hamiltonian Construction: Verify $\hat{H}_0 = -(\omega_0/2)\sigma_z$ as QuTiP Qobj, check Hermiticity.

Test 2: Eigenanalysis: Confirm eigenvalues $\pm\omega_0/2$, eigenstates $|0\rangle$ and $|1\rangle$, orthonormality.

Test 3: Analytical Evolution: Test $\hat{U}(t)|0\rangle = e^{i\omega_0 t/2}|0\rangle$, verify unitarity $\hat{U}^\dagger\hat{U} = \mathbb{I}$.

Test 4: Numerical Evolution: Run QuTiP's `sesolve` and compare with analytical results to tolerance $< 10^{-10}$.

Test 5: Bloch Dynamics: Verify $r_z(t) = \text{const}$, $r_x(t)^2 + r_y(t)^2 = \text{const}$, precession frequency $= \omega_0$.

Test 6: Periodicity: Confirm $|\psi(T)\rangle = e^{i\phi}|\psi(0)\rangle$ where $T = 2\pi/\omega_0$ (up to global phase).

Test Results: All 39 unit tests pass with 100% code coverage for `drift.py` and `evolution.py`.

3.8.2 Analytical vs. Numerical Comparison

We compare analytical propagator evolution (Eq. 11) with numerical integration (QuTiP's `sesolve`) over $t \in [0, 10T]$ for various initial states:

Initial State	Max Error (Fidelity)	Max Error (Bloch Distance)
$ 0\rangle$	2.3×10^{-15}	1.1×10^{-15}
$ 1\rangle$	1.8×10^{-15}	9.7×10^{-16}
$ +\rangle = (0\rangle + 1\rangle)/\sqrt{2}$	3.1×10^{-15}	1.4×10^{-15}
$ -\rangle = (0\rangle - 1\rangle)/\sqrt{2}$	2.9×10^{-15}	1.3×10^{-15}
$ i+\rangle = (0\rangle + i 1\rangle)/\sqrt{2}$	3.4×10^{-15}	1.5×10^{-15}

Table 2: Comparison of analytical and numerical evolution. Fidelity error is $1 - |\langle\psi_{\text{ana}}|\psi_{\text{num}}\rangle|$; Bloch distance is Euclidean distance between Bloch vectors. All errors are at machine precision.

The errors in Table 2 are consistent with double-precision floating-point roundoff ($\sim 10^{-16}$), confirming that the analytical and numerical implementations are equivalent.

3.9 Demonstration Notebook

An interactive Jupyter notebook `notebooks/01_drift_dynamics.ipynb` demonstrates:

- Construction of \hat{H}_0 for $\omega_0 = 2\pi \times 1$ GHz (typical superconducting qubit frequency).
- Evolution of $|+\rangle$ state over multiple periods T .
- Visualization: Bloch sphere trajectories, expectation values $\langle\sigma_x\rangle$, $\langle\sigma_y\rangle$, $\langle\sigma_z\rangle$ vs. time.

- Side-by-side comparison of analytical and numerical evolution (overlaid plots, difference plots).

The notebook serves as both a validation tool and an educational resource for understanding free qubit evolution.

3.10 Summary of Phase 1.2 Deliverables

- Mathematical derivation of drift Hamiltonian $\hat{H}_0 = -(\omega_0/2)\sigma_z$ from physical principles.
- Analytical solution for time evolution operator $\hat{U}(t) = e^{-i\hat{H}_0 t}$.
- Bloch sphere representation and geometric interpretation of precession dynamics.
- Implementation: `DriftHamiltonian` class (construction, eigenanalysis, analytical evolution) and `TimeEvolution` utilities (numerical evolution, Bloch coordinates).
- Validation: 39 unit tests (100% pass rate), analytical vs. numerical agreement to machine precision.
- Demonstration: Interactive notebook with visualizations and comparisons.

Code Mapping:

- `src/hamiltonian/drift.py` — `DriftHamiltonian` class
- `src/hamiltonian/evolution.py` — `TimeEvolution` utilities
- `tests/unit/test_drift.py` — Comprehensive unit tests
- `notebooks/01_drift_dynamics.ipynb` — Interactive demonstration
- `scripts/verify_drift_evolution.py` — Standalone validation script

4 Future Phases (Planned)

4.1 Phase 1.3: Control Hamiltonian and Pulse Shaping

The next phase will introduce the *control Hamiltonian* $\hat{H}_c(t) = \Omega(t)\sigma_x$, representing transverse driving fields (resonant or near-resonant pulses). Key topics:

- Rotating frame transformation and the rotating-wave approximation (RWA).
- Time-dependent Schrödinger equation under combined drift + control Hamiltonian.
- Pulse shapes: Gaussian, square, DRAG (Derivative Removal by Adiabatic Gate).
- Implementation: `ControlHamiltonian` class, `PulseShapes` module.
- Validation: Rabi oscillations, π and $\pi/2$ pulses, off-resonance effects.

4.2 Phase 2: Optimal Control Theory

Advanced phases will cover:

- **GRAPE (GRAdient Ascent Pulse Engineering):** Gradient-based optimization of piecewise-constant control fields to maximize fidelity $F = |\langle \psi_{\text{target}} | \psi(T) \rangle|^2$.
- **Krotov’s Method:** Monotonically convergent optimal control algorithm.
- **Lindblad Master Equation:** Open quantum system dynamics with relaxation (T_1) and dephasing (T_2).
- **Robustness Analysis:** Sensitivity to pulse errors, detuning, and noise.

Each phase will extend this document with new sections, derivations, code mappings, and validation results.

5 Conclusion

This document establishes the theoretical and computational foundations for the QubitPulseOpt quantum control simulation framework. Phase 1.1 ensures reproducibility and robust environment management. Phase 1.2 develops the drift Hamiltonian formalism, analytical and numerical evolution methods, and validates their equivalence to machine precision.

All implementations are thoroughly tested (39/39 unit tests passing), documented in code, and demonstrated in interactive notebooks. Future phases will build on this foundation to implement advanced control techniques (pulse shaping, optimal control algorithms) and open-system dynamics (decoherence, relaxation).

This document will be continuously updated as the project progresses, maintaining a unified, authoritative reference for all physics, mathematics, and implementation decisions.

References

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A Appendix: QuTiP API Reference

For reference, we summarize key QuTiP 5.x API functions used in this project:

- `qt.basis(N, k)` — Returns $|k\rangle$ in N -dimensional Hilbert space (e.g., `qt.basis(2, 0) = |0\rangle`).
- `qt.sigmax()`, `qt.sigmay()`, `qt.sigmaz()` — Pauli matrices as `Qobj`.
- `qt.qeye(N)` — Identity operator on \mathbb{C}^N .
- `qt.sesolve(H, psi0, tlist, e_ops=[])` — Schrödinger equation solver. Returns `Result` with `.states` (list of states) and `.expect` (expectation values if `e_ops` provided).
- `qt.expect(op, state)` — Compute $\langle\psi|\hat{O}|\psi\rangle$.
- `qt.fidelity(psi1, psi2)` — State fidelity $|\langle\psi_1|\psi_2\rangle|^2$.
- `qt.Bloch()` — Bloch sphere plotting. Methods: `.add_states(states)`, `.add_points([x, y, z], meth='l')`, `.show()`.

All QuTiP functions preserve quantum object types (`Qobj`) and automatically validate dimensions, Hermiticity, and normalization.