Quantum Dynamics

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Goals:

- 1. Partial differential equation
- 2. Spectral method (Fourier transform)



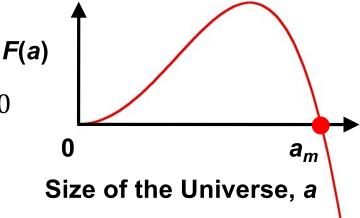


Quantum Universe

Wheeler-deWitt equation

$$\left[-\hbar^2 \frac{d^2}{da^2} + \left(\frac{3\pi c^3}{2G} \right)^2 \left(a^2 - \frac{a^4}{a_m^2} \right) \right] \psi(a) = 0$$

$$F(a)$$



CREATION OF UNIVERSES FROM NOTHING

CREATION OF UNIVERSES FROM NOTHING

Alexander VILENKIN

Physics Department, Tufts University, Medford, MA 02155, USA

Received 11 June 1982

A cosmological model is proposed in which the universe is created by quantum tunneling from literally nothing into a de Sitter space. After the tunneling, the model evolves along the lines of the inflationary scenario. This model does not have a big-bang singularity and does not require any initial or boundary conditions.

IS IT POSSIBLE TO CREATE A UNIVERSE IN THE LABORATORY BY QUANTUM TUNNELING?

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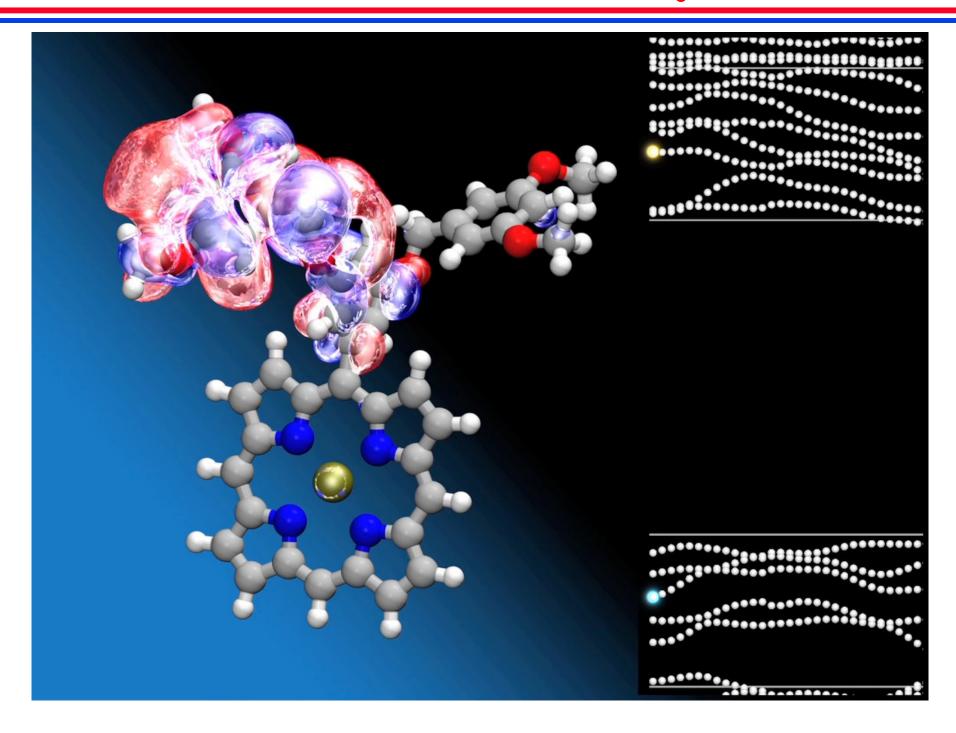
Jemal GUVEN

Instituto de Ciencias Nucleares, Universidad Nacional Autonoma de Mexico, Circuito Exterior C.U. A. Postal 70-543, 04510 Mexico D.F. Mexico

Nucl. Phys. B339, 417 ('90)

Phys. Lett. **117B**, 25 ('82)

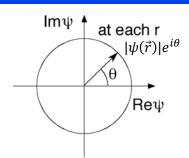
Photoexcited Electron Dynamics



Wave Equation

Complex wave function

$$\psi(\vec{r},t) = \text{Re}\psi(\vec{r},t) + i\text{Im}\psi(\vec{r},t) \in \mathbb{C}\left(i = \sqrt{-1}\right)$$



Probability

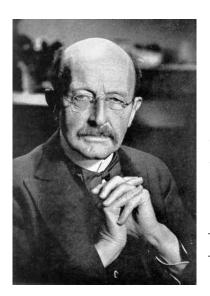
$$P(\vec{r},t) = \psi^*(\vec{r},t)\psi(\vec{r},t) = |\psi(\vec{r},t)|^2 = |\text{Re}\psi(\vec{r},t)|^2 + |\text{Im}\psi(\vec{r},t)|^2$$

Normalization

$$\int dx \int dy \int dz \, |\psi(\vec{r},t)|^2 = 1$$

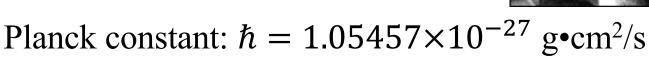
 $\psi^*\psi = (\psi_0 - i\psi_1)(\psi_0 + i\psi_1) = \psi_0^2 + \psi_1^2 \ge 0$

• Schrödinger (partial differential) equation



$$i\hbar \frac{\partial}{\partial t} \psi(\vec{r}, t) = \left[-\frac{\hbar^2}{2m} \nabla^2 + V(\vec{r}) \right] \psi(\vec{r}, t)$$

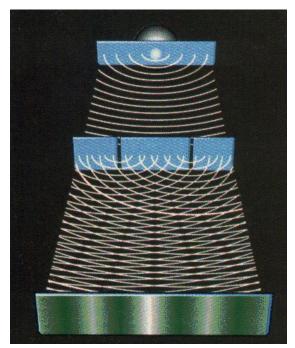
Laplacian:
$$\nabla^2 = \frac{\partial^2}{\partial x^2} + \frac{\partial^2}{\partial y^2} + \frac{\partial^2}{\partial z^2}$$

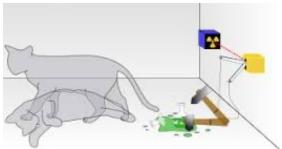


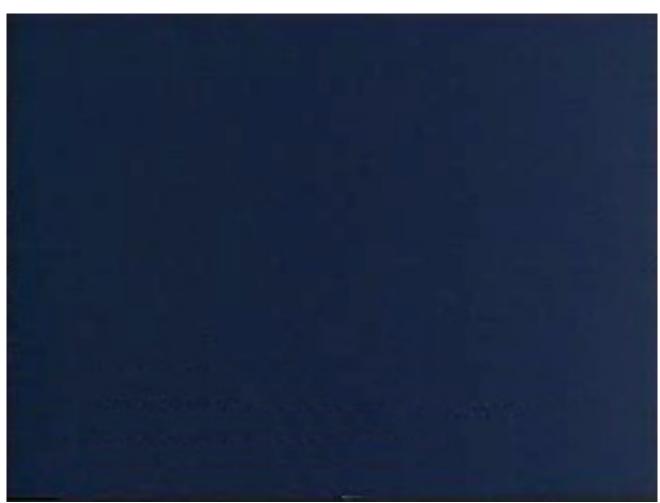


Single-Electron Double-Slit Experiment

What wave?







http://rdg.ext.hitachi.co.jp/rd/moviee/doubleslite-n.mpeg
Akira Tonomura (Hitachi, Ltd.)

Atomic Unit

Length, energy & time in atomic unit

$$\frac{e^{2}}{r} = \frac{\hbar^{2}}{mr^{2}} \begin{cases} \vec{r} = \frac{\hbar^{2}}{me^{2}} \vec{r}' & \frac{\hbar^{2}}{me^{2}} = 0.529177 \text{ Å} & \text{Bohr} \\ E = \frac{e^{2}}{r} & V = \frac{me^{4}}{\hbar^{2}} V' & \frac{me^{4}}{\hbar^{2}} = 27.2116 \text{ eV} & \text{Hartree} \\ E = \frac{\hbar}{t} & t = \frac{\hbar^{3}}{me^{4}} t' & \frac{\hbar^{3}}{me^{4}} = 0.0241889 \text{ fs} \end{cases}$$

Gaussian unit

Time-dependent Schrödinger equation in atomic unit

$$i\hbar \frac{\partial}{\partial t} \psi(\vec{r}, t) = \left[-\frac{\hbar^2}{2m} \nabla^2 + V(\vec{r}) \right] \psi(\vec{r}, t)$$



$$i\frac{\partial}{\partial t'}\psi(\vec{r}',t') = \left[-\frac{\nabla'^2}{2} + V(\vec{r}')\right]\psi(\vec{r}',t')$$





Two-Dimensional Electron

Schrödinger equation (in atomic unit)

$$i\frac{\partial}{\partial t}\psi(x,y,t) = H\psi(x,y,t)$$

Hamiltonian operator

$$H = -\frac{1}{2} \frac{\partial^2}{\partial x^2} - \frac{1}{2} \frac{\partial^2}{\partial y^2} + V(x, y)$$
$$= T_x + T_y + V$$

Periodic boundary condition

 $\begin{cases} \psi(x + L_x, y) = \psi(x, y) \\ \psi(x, y + L_y) = \psi(x, y) \end{cases}$



The Nobel Prize in Physics 1985

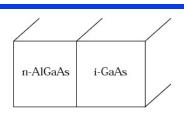
"for the discovery of the quantized Hall effect"

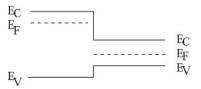


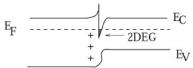
Klaus von Klitzing

Federal Republic of Germany

Max-Planck-Institut für Festkörperforschung Stuttgart, Federal Republic of Germany b. 1943









The Nobel Prize in Physics 1998

"for their discovery of a new form of quantum fluid with fractionally charged excitations"



Robert B. Laughlin

1/3 of the prize

Stanford University Stanford, CA, USA b. 1950



Horst L. Störmer

1/3 of the prize Federal Republic of Germany

New York, NY, USA b. 1949



Daniel C. Tsui

1/3 of the prize

Columbia University Princeton University Princeton, NJ, USA b. 1939 (in Henan, China)

Layered Materials Genome

• Atomically-thin layered materials will dominate materials science in this

century

Geim & Grigorieva, *Nature* **499**, 419 ('13)

The Nobel Prize in Physics 2010

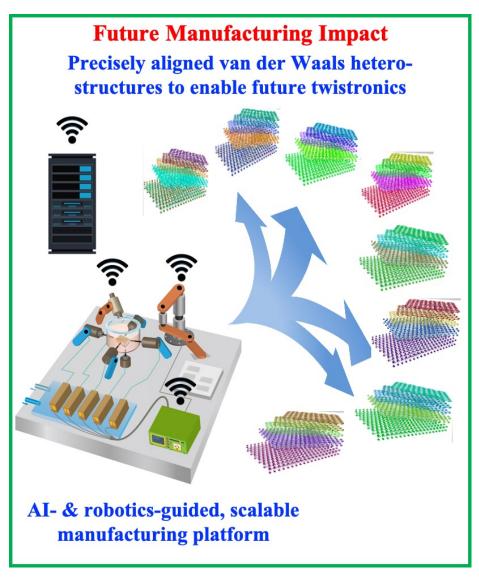


© The Nobel Foundat Photo: U. Montan Andre Geim Prize share: 1/2



© The Nobel Foundation. Photo: U. Montan Konstantin Novoselov Prize share: 1/2

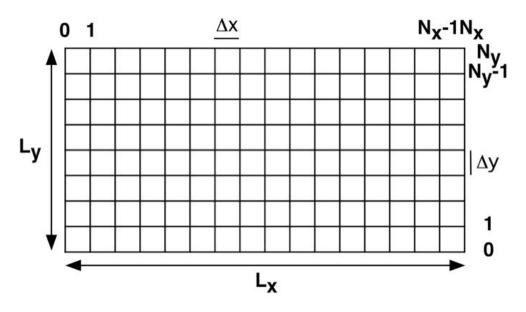
 Tuning material properties in desired ways by building heterostructures composed of unlimited combinations of atomically thin layers in a way similar to genomics



https://aiqma.netlify.app

Spatial Discretization

• Regular 2D mesh: $\psi_{jk} = \psi(j\Delta x, k\Delta y) \ (\Delta x = L_x/N_x \& \Delta y = L_y/N_y)$



Finite differencing

$$\begin{cases} (T_{x}\psi)_{j,k} = -\frac{1}{2} \frac{\psi_{j-1,k} - 2\psi_{j,k} + \psi_{j+1,k}}{(\Delta x)^{2}} \\ (T_{y}\psi)_{j,k} = -\frac{1}{2} \frac{\psi_{j,k-1} - 2\psi_{j,k} + \psi_{j,k+1}}{(\Delta y)^{2}} \\ (V\psi)_{j,k} = V_{j,k}\psi_{j,k} \end{cases}$$

$$\frac{\partial^{2}}{\partial x^{2}} \psi(x, y, t)$$

$$= \frac{\partial \psi(x + \Delta/2, y, t) / \partial x - \partial \psi(x - \Delta/2, y, t) / \partial x}{\Delta}$$

$$= \frac{\psi(x + \Delta, y, t) - \psi(x, y, t)}{\Delta} - \frac{\psi(x, y, t) - \psi(x - \Delta, y, t)}{\Delta}$$

$$= \frac{\psi(x + \Delta, y, t) - 2\psi(x, y, t) + \psi(x - \Delta, y, t)}{\Delta^{2}}$$

Temporal Propagation

- Formal solution to the Schrödinger equation: $\frac{\partial}{\partial t}\psi(t) = -iH\psi(t)$ $\psi(t + \Delta t) = \exp(-iH\Delta t)\psi(t)$ $\exp(-i\widehat{H}t) = \sum_{n=0}^{\infty} \frac{(-it)^n}{n!} \widehat{H}^n$
- Split-operator method (Trotter-expansion): unitary!

$$\psi(t + \Delta t) = \exp(-i(T_x + T_y + V)\Delta t)\psi(t)^{\exp(-i(T_x + T_y)\Delta t)} \stackrel{\exp(-iT_x\Delta t)\exp(-iT_y\Delta t)}{\Longrightarrow} \exp(-iT_x\Delta t)\exp(-iT_y\Delta t)$$

$$= \exp(-iV\Delta t/2)\exp(-iT_x\Delta t)\exp(-iT_y\Delta t)\exp(-iV\Delta t/2)\psi(t) + O([\Delta t]^3)$$
Split in a way each operator is easily exponentiated

• Potential propagator (mesh point-by-point complex-number multiplications) $(u_0 + iu_1)(\psi_0 + i\psi_1) = (u_0\psi_0 - u_1\psi_1) + i(u_1\psi_0 + u_0\psi_1)$

$$(\exp(-iV\Delta t/2)\psi)_{jk} = \exp(-iV_{jk}\Delta t/2)\psi_{jk}$$

$$= [\cos(V_{jk}\Delta t/2) - i\sin(V_{jk}\Delta t/2)][\operatorname{Re}\psi_{jk} + i\operatorname{Im}\psi_{jk}]$$

$$= [\cos(V_{jk}\Delta t/2)\operatorname{Re}\psi_{jk} + \sin(V_{jk}\Delta t/2)\operatorname{Im}\psi_{jk}]$$

$$+i[-\sin(V_{jk}\Delta t/2)\operatorname{Re}\psi_{jk} + \cos(V_{jk}\Delta t/2)\operatorname{Im}\psi_{jk}]$$

$$\exp(ia) = \sum_{n=0}^{\infty} \frac{(ia)^n}{n!} = \left(1 - \frac{a^2}{2!} + \frac{a^4}{4!} - \cdots\right) + i\left(a - \frac{a^3}{3!} + \frac{a^5}{5!} - \cdots\right) \qquad \begin{bmatrix} \cos(V_{jk}\Delta t/2) & \sin(V_{jk}\Delta t/2) \\ -\sin(V_{jk}\Delta t/2) & \cos(V_{jk}\Delta t/2) \end{bmatrix} \begin{bmatrix} \operatorname{Re}\psi_{jk} \\ \operatorname{Im}\psi_{jk} \end{bmatrix}$$

Kinetic Propagator

Mesh-point coupling

$$T_x \psi_{j,k} = b \psi_{j-1,k} + 2a \psi_{j,k} + b \psi_{j+1,k}$$

• Tridiagonal matrix representation

$$T_{x} = \begin{bmatrix} 2a & b & & & & & \\ b & 2a & b & & & \\ & b & 2a & b & & \\ & & \ddots & \ddots & \ddots & \\ & & b & 2a & b & \\ & & & b & 2a & b \\ & & & b & 2a \end{bmatrix}$$



Note the periodic boundary condition

$$\begin{cases} a = 1/2(\Delta x)^2 \\ b = -1/2(\Delta x)^2 \end{cases}$$

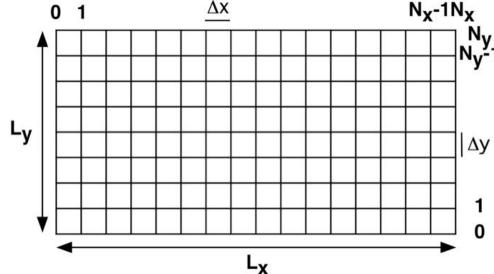
Space Splitting Method (SSM)

• 2×2 block-diagonal decomposition & split-operator exponentiation

$$T_{x} = \begin{bmatrix} 2a & b & & & & & b \\ b & 2a & b & & & & \\ b & 2a & b & & & & \\ b & 2a & b & & & & \\ b & 2a & b & & & \\ b & 2a & b & & & \\ b & 2a & b & & & \\ b & 2a & b & & \\ c_{n} & = \frac{1}{2} \left[\exp\left(-\frac{i\Delta t}{n}(a+b)\right) + \exp\left(-\frac{i\Delta t}{n}(a-b)\right) \right] \\ = \frac{1}{2} \begin{bmatrix} a & b & & & \\ b & a & & & \\ c_{n} & = \frac{1}{2} \left[\exp\left(-\frac{i\Delta t}{n}(a+b)\right) + \exp\left(-\frac{i\Delta t}{n}(a-b)\right) \right] \\ = \frac{1}{2} \begin{bmatrix} a & b & & & \\ b & a & & & \\ b & a & & & \\ c_{n} & = \frac{1}{2} \left[\exp\left(-\frac{i\Delta t}{n}(a+b)\right) + \exp\left(-\frac{i\Delta t}{n}(a-b)\right) \right] \\ = \frac{1}{2} \begin{bmatrix} a & b & & & \\ b & a & & & \\ b & a & & & \\ c_{n} & = \frac{1}{2} \left[\exp\left(-\frac{i\Delta t}{n}(a+b)\right) + \exp\left(-\frac{i\Delta t}{n}(a-b)\right) \right] \\ = \frac{1}{2} \begin{bmatrix} a & b & & & \\ b & a & & & \\ b & a & & & \\ c_{n} & = \frac{1}{2} \left[\exp\left(-\frac{i\Delta t}{n}(a+b)\right) + \exp\left(-\frac{i\Delta t}{n}(a-b)\right) \right] \\ = \frac{1}{2} \begin{bmatrix} a & b & & & \\ b & a & & & \\ b & a & & & \\ b & a & & & \\ c_{n} & = \frac{1}{2} \left[\exp\left(-\frac{i\Delta t}{n}(a+b)\right) + \exp\left(-\frac{i\Delta t}{n}(a-b)\right) \right] \\ = \frac{1}{2} \begin{bmatrix} a & b & & & & \\ b & a & & & \\ b & a & & & \\ b & a & & & \\ c_{n} & = \frac{1}{2} \left[\exp\left(-\frac{i\Delta t}{n}(a+b)\right) + \exp\left(-\frac{i\Delta t}{n}(a-b)\right) \right] \\ = \frac{1}{2} \begin{bmatrix} a & b & & & & \\ b & a & & & \\ b & a & & & \\ b & a & & & \\ c_{n} & = \frac{1}{2} \left[\exp\left(-\frac{i\Delta t}{n}(a+b)\right) + \exp\left(-\frac{i\Delta t}{n}(a-b)\right) \right] \\ = \frac{1}{2} \begin{bmatrix} a & b & & & & \\ b & a & & & \\ b & a & & & \\ b & a & & & \\ c_{n} & = \frac{1}{2} \left[\exp\left(-\frac{i\Delta t}{n}(a+b)\right) + \exp\left(-\frac{i\Delta t}{n}(a-b)\right) \right] \\ = \frac{1}{2} \begin{bmatrix} a & b & & & & \\ b & a & & & \\ b & a & & & \\ b & a & & & \\ c_{n} & = \frac{1}{2} \left[\exp\left(-\frac{i\Delta t}{n}(a+b)\right) + \exp\left(-\frac{i\Delta t}{n}(a-b)\right) \right] \\ = \frac{1}{2} \begin{bmatrix} a & b & & & & \\ b & a & & & \\ b & a & & & \\ b & a & & & \\ c_{n} & = \frac{1}{2} \left[\exp\left(-\frac{i\Delta t}{n}(a+b)\right) + \exp\left(-\frac{i\Delta t}{n}(a-b)\right) \right] \\ = \frac{1}{2} \begin{bmatrix} a & b & & & & \\ c_{n} & a & b & & \\ c_{n} & a & b & & \\ c_{n} & a & b & & \\ c_{n} & c_{n} & c_{n} & c_{n} \\ c_{n} & c_{$$

Data Structures in Program qd.c

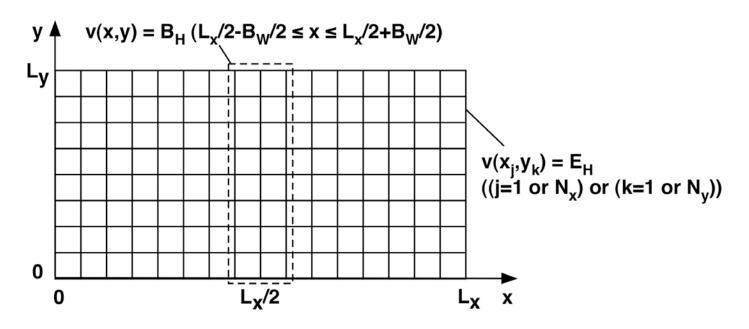
- Wave function: psi[NX+2][NY+2][2]
- Periodic boundary condition by auxiliary elements



https://aiichironakano.github.io/phys516/src/QD/

Potential Propagator in qd.c

• Potential barrier: v[NX+2][NY+2]



- Potential propagator: $\exp(-iV\Delta t/2)$, u[NX+2][NY+2][2]
- Potential propagation: $\psi \leftarrow \exp(-iV\Delta t/2) \psi$

```
for (sx=1; sx<=NX; sx++)
  for (sy=1; sy<=NY; sy++) {
    wr=u[sx][sy][0]*psi[sx][sy][0]-u[sx][sy][1]*psi[sx][sy][1];
    wi=u[sx][sy][0]*psi[sx][sy][1]+u[sx][sy][1]*psi[sx][sy][0];
    psi[sx][sy][0]=wr;
    psi[sx][sy][1]=wi;}</pre>
```

Kinetic Propagator in qd.c

```
\left(U_{\chi}^{(\text{half})}\psi\right)_{i,j} = \varepsilon_2^- \delta_{\text{mod}(i,2),0} \psi_{i-1,j} + \varepsilon_2^+ \psi_{i,j} + \varepsilon_2^- \delta_{\text{mod}(i,2),1} \psi_{i+1,j}
   \left(U_{\chi}^{(\text{full})}\psi\right)_{i,i} = \varepsilon_1^- \delta_{\text{mod}(i,2),1}\psi_{i-1,j} + \varepsilon_1^+ \psi_{i,j} + \varepsilon_1^- \delta_{\text{mod}(i,2),0}\psi_{i+1,j}
/* WRK | PSI holds the new | old wave function */
for (sx=1; sx\leq NX; sx++)
  for (sy=1; sy<=NY; sy++) {</pre>
     wr=al[d][t][0]*psi[sx][sy][0]-al[d][t][1]*psi[sx][sy][1];
     wi=al[d][t][0]*psi[sx][sy][1]+al[d][t][1]*psi[sx][sy][0];
     if (d==0) {
       wr+=(blx[t][sx][0]*psi[sx-1][sy][0]-blx[t][sx][1]*psi[sx-1][sy][1]);
       wi+=(blx[t][sx][0]*psi[sx-1][sy][1]+blx[t][sx][1]*psi[sx-1][sy][0]);
       wr+=(bux[t][sx][0]*psi[sx+1][sy][0]-bux[t][sx][1]*psi[sx+1][sy][1]);
       wi+=(bux[t][sx][0]*psi[sx+1][sy][1]+bux[t][sx][1]*psi[sx+1][sy][0]);
     else if (d==1) {
       wr+=(bly[t][sy][0]*psi[sx][sy-1][0]-bly[t][sy][1]*psi[sx][sy-1][1]);
       wi+=(bly[t][sy][0]*psi[sx][sy-1][1]+bly[t][sy][1]*psi[sx][sy-1][0]);
       wr+=(buy[t][sy][0]*psi[sx][sy+1][0]-buy[t][sy][1]*psi[sx][sy+1][1]);
       wi+=(buy[t][sy][0]*psi[sx][sy+1][1]+buy[t][sy][1]*psi[sx][sy+1][0]);
     wrk[sx][sy][0]=wr;
     wrk[sx][sy][1]=wi;}
/* Copy the new wave function back to PSI */
for (sx=1; sx\leq NX; sx++)
  for (sy=1; sy<=NY; sy++)</pre>
```

for (s=0; s<=1; s++) psi[sx][sy][s]=wrk[sx][sy][s];

Initial Wave Function

Gaussian wave packet

$$\psi(x, t = 0) = C \exp\left(-\frac{(x - x_0)^2}{4\sigma^2}\right) \exp(ik_0 x) \sin\left(\frac{\pi y}{L_y}\right)$$

Symbol	Variable in qd.c			
x_0 (packet center)	X0			
σ (packet spread)	S0			
$k_0^2/2$ (energy)	E0			

Quantum Dynamics—II One Dimensional System

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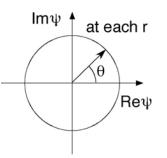
Goal: Understand qd1.c



Wave Equation

Complex wave function

$$\psi(x,t) = \text{Re}\psi(x,t) + i\text{Im}\psi(x,t) \in \mathbb{C} \left(i = \sqrt{-1}\right)$$



Normalization

$$\int dx |\psi(x,t)|^2 = 1$$

$$\psi^* \psi = (\psi_0 - i\psi_1)(\psi_0 + i\psi_1) = \psi_0^2 + \psi_1^2 \ge 0$$

Schrödinger equation (in atomic unit)

$$i\frac{\partial}{\partial t}\psi(x,t) = H\psi(x,t)$$

Hamiltonian operator

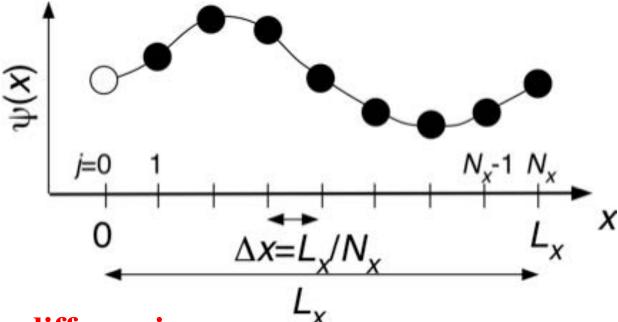
$$H = -\frac{1}{2} \frac{\partial^2}{\partial x^2} + V(x) = T_x + V$$

Periodic boundary condition

$$\psi(x + L_x) = \psi(x)$$

Spatial Discretization

• Regular 1D mesh: $\psi_i = \psi(j\Delta x) \ (\Delta x = L_x/N_x)$



• Finite differencing

$$\begin{cases} \left(T_{x}\psi\right)_{j} = -\frac{1}{2}\frac{\psi_{j-1} - 2\psi_{j} + \psi_{j+1}}{\left(\Delta x\right)^{2}} \\ \left(V\psi\right)_{j} = V_{j}\psi_{j} \end{cases}$$

$$= \frac{\frac{\partial^{2}}{\partial x^{2}}\psi(x,t)}{\frac{\Delta}{\partial x^{2}}}$$

$$= \frac{\frac{\partial\psi(x+\Delta/2,t)/\partial x - \partial\psi(x-\Delta/2,t)/\partial x}{\Delta}}{\frac{\Delta}{\Delta}}$$

$$= \frac{\frac{\psi(x+\Delta,t) - \psi(x,t)}{\Delta} - \frac{\psi(x,t) - \psi(x-\Delta,t)}{\Delta}}{\frac{\Delta}{\Delta}}$$

$$= \frac{\psi(x+\Delta,t) - 2\psi(x,t) + \psi(x-\Delta,t)}{\frac{\Delta}{\Delta}}$$

$$\frac{\partial^{2}}{\partial x^{2}} \psi(x,t)$$

$$= \frac{\partial \psi(x+\Delta/2,t)/\partial x - \partial \psi(x-\Delta/2,t)/\partial x}{\Delta}$$

$$= \frac{\frac{\psi(x+\Delta,t) - \psi(x,t)}{\Delta} - \frac{\psi(x,t) - \psi(x-\Delta,t)}{\Delta}}{\Delta}$$

$$= \frac{\psi(x+\Delta,t) - 2\psi(x,t) + \psi(x-\Delta,t)}{\Delta^{2}}$$

Temporal Propagation

• Formal solution to the Schrödinger equation: $\frac{\partial}{\partial t}\psi(t) = -iH\psi(t)$

$$\psi(t + \Delta t) = \exp(-iH\Delta t)\psi(t) \quad \exp(-i\widehat{H}t) = \sum_{n=0}^{\infty} \frac{(-it)^n}{n!} \widehat{H}^n$$

Split-operator method: unitary!

$$\psi(t + \Delta t) = \exp(-i(T_x + V)\Delta t)\psi(t)$$
$$= \exp(-iV\Delta t/2)\exp(-iT_x\Delta t)\exp(-iV\Delta t/2)\psi(t) + O([\Delta t]^3)$$

Split in a way each operator is easily exponentiated

 Potential propagator (mesh point-by-point complex-number multiplications)

$$(u_0 + iu_1)(\psi_0 + i\psi_1) = (u_0\psi_0 - u_1\psi_1) + i(u_1\psi_0 + u_0\psi_1)$$

$$(\exp(-iV\Delta t/2)\psi)_{j} = \exp(-iV_{j}\Delta t/2)\psi_{j}$$

$$= [\cos(V_{j}\Delta t/2) - i\sin(V_{j}\Delta t/2)][\operatorname{Re}\psi_{j} + i\operatorname{Im}\psi_{j}]$$

$$= [\cos(V_{j}\Delta t/2)\operatorname{Re}\psi_{j} + \sin(V_{j}\Delta t/2)\operatorname{Im}\psi_{j}]$$

$$+i[-\sin(V_{j}\Delta t/2)\operatorname{Re}\psi_{j} + \cos(V_{j}\Delta t/2)\operatorname{Im}\psi_{j}]$$

$$\exp(ia) = \sum_{n=0}^{\infty} \frac{(ia)^n}{n!} = \left(1 - \frac{a^2}{2!} + \frac{a^4}{4!} - \cdots\right) + i\left(a - \frac{a^3}{3!} + \frac{a^5}{5!} - \cdots\right) \qquad \begin{bmatrix} \cos(V_{jk}\Delta t/2) & \sin(V_{jk}\Delta t/2) \\ -\sin(V_{jk}\Delta t/2) & \cos(V_{jk}\Delta t/2) \end{bmatrix} \begin{bmatrix} \operatorname{Re}\psi_{jk} \\ \operatorname{Im}\psi_{jk} \end{bmatrix}$$

Kinetic Propagator: It's a Matrix!

• Mesh-point coupling

$$T_x \psi_j = b \psi_{j-1} + 2a \psi_j + b \psi_{j+1}$$

• Tridiagonal matrix representation

$$T_{x} = \begin{bmatrix} 2a & b & & & & & & \\ b & 2a & b & & & & \\ & b & 2a & b & & & \\ & & \ddots & \ddots & \ddots & & \\ & & b & 2a & b & \\ & & & b & 2a & b \\ & & & b & 2a \end{bmatrix}$$

Note the periodic boundary condition

$$\begin{cases} a = 1/2(\Delta x)^2 \\ b = -1/2(\Delta x)^2 \end{cases}$$





Space Splitting Method (SSM)

2×2 block-diagonal decomposition & split-operator exponentiation

$$T_{x} = \begin{bmatrix} 2a & b & & & & b \\ b & 2a & b & & & b \\ b & 2a & b & & & & \\ & & & \ddots & \ddots & \ddots & \\ & & & & b & 2a & b \\ b & & & & b & 2a \end{bmatrix}$$

$$= \sum_{n=0}^{\infty} \frac{1}{n!} \begin{bmatrix} \bullet & \bullet & \bullet \\ \bullet & \bullet & \bullet \\ \bullet & \bullet & \bullet \end{bmatrix} = \begin{bmatrix} e^{\bullet} & \bullet & \bullet \\ \bullet & \bullet & \bullet \\ \bullet & \bullet & \bullet \end{bmatrix}$$

$$= \sum_{n=0}^{\infty} \frac{1}{n!} \begin{bmatrix} \bullet & \bullet & \bullet \\ \bullet & \bullet & \bullet \\ \bullet & \bullet & \bullet \end{bmatrix}$$

Split-operator (Trotter expansion) again

How? Block diagonal → block-by-block exponentiation

Space Splitting Method (SSM)

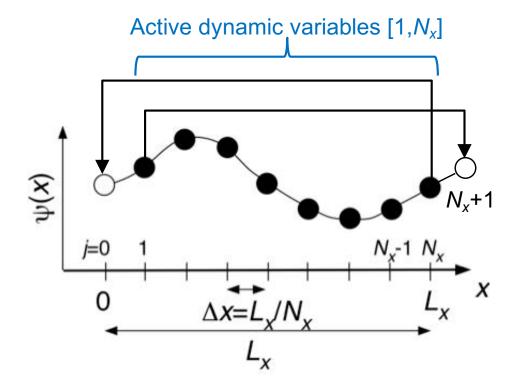
$$\begin{cases} \varepsilon_n^+ = \frac{1}{2} \left[\exp\left(-\frac{i\Delta t}{n}(a+b)\right) + \exp\left(-\frac{i\Delta t}{n}(a-b)\right) \right] & \textbf{Just need 2} \times \textbf{2 exponentiation} \\ \varepsilon_n^- = \frac{1}{2} \left[\exp\left(-\frac{i\Delta t}{n}(a+b)\right) - \exp\left(-\frac{i\Delta t}{n}(a-b)\right) \right] & \exp\left(-\frac{i\Delta}{2} \begin{bmatrix} a & b \\ b & a \end{bmatrix} \right) \end{cases}$$

Use eigen-decomposition & telescoping

Data Structures in Program qd1.c

- Wave function: psi[NX+2][2] $psi[j][0|1] = (Re|Im)\psi_{j\Delta x}$
- Periodic boundary condition by auxiliary elements

```
for (s=0; s<=1; s++) {
  psi[0][s] = psi[NX][s];
  psi[NX+1][s] = psi[1][s];
}</pre>
```



https://aiichironakano.github.io/phys516/src/QD/

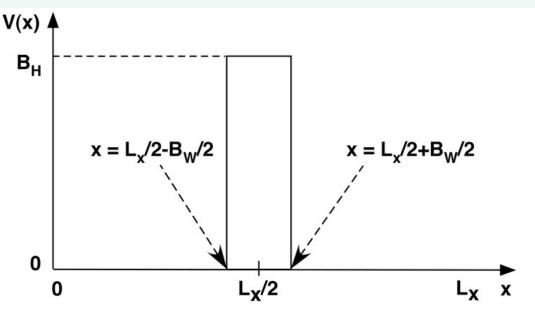
Potential Propagator in qd1.c

- Potential barrier: v[NX+2]
- Potential propagator: $\exp(-iV\Delta t/2)$, u[NX+2][2]
- Potential propagation: $\psi \leftarrow \exp(-iV\Delta t/2) \psi$

```
for (sx=1; sx<=NX; sx++)
  wr=u[sx][0]*psi[sx][0]-u[sx][1]*psi[sx][1];
  wi=u[sx][0]*psi[sx][1]+u[sx][1]*psi[sx][0];
  psi[sx][0]=wr;
  psi[sx][1]=wi;</pre>
```

$$\begin{cases} u[j][0] = \cos\left(-\frac{\Delta}{2}V_j\right) \\ u[j][1] = \sin\left(-\frac{\Delta}{2}V_j\right) \end{cases}$$

$$\exp\left(-\frac{iV_{j}\Delta}{2}\right)\psi_{j} \equiv u\psi = (u_{0} + iu_{1})(\psi_{0} + i\psi_{1}) = \underbrace{(u_{0}\psi_{0} - u_{1}\psi_{1})}_{\text{new }\psi_{1}} + i\underbrace{(u_{0}\psi_{1} + u_{1}\psi_{0})}_{\text{new }\psi_{1}}$$



Kinetic Propagator in qd1.c

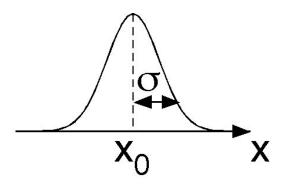
```
 \begin{pmatrix} U_{x}^{(\text{half})} \psi \end{pmatrix}_{i} = \varepsilon_{2}^{-} \delta_{\text{mod}(i,2),0} \psi_{i-1} + \varepsilon_{2}^{+} \psi_{i} + \varepsilon_{2}^{-} \delta_{\text{mod}(i,2),1} \psi_{i+1} 
 \begin{pmatrix} U_{x}^{(\text{full})} \psi \end{pmatrix}_{i} = \varepsilon_{1}^{-} \delta_{\text{mod}(i,2),1} \psi_{i-1} + \varepsilon_{1}^{+} \psi_{i} + \varepsilon_{1}^{-} \delta_{\text{mod}(i,2),0} \psi_{i+1} 
for (sx=1; sx<=NX; sx++) { // wrk[][]/psi[][] holds new old wave function
         wr=al[t][0]*psi[sx][0]-al[t][1]*psi[sx][1]; // al[0|1][]: \alpha_{half|full}
         wi=al[t][0]*psi[sx][1]+al[t][1]*psi[sx][0];
         wr += (bl[t][sx][0]*psi[sx-1][0]-bl[t][sx][1]*psi[sx-1][1]); // bl[0|1][][]: \beta_{table : table : tab
                                                                                                                                                                                                                                                                                                                                            half|full
         wi+=(bl[t][sx][0]*psi[sx-1][1]+bl[t][sx][1]*psi[sx-1][0]);
        wi+=(bu[t][sx][0]*psi[sx+1][1]+bu[t][sx][1]*psi[sx+1][0]);
         wrk[sx][0]=wr;
                                                                                                          \psi_i \leftarrow \beta_l \psi_{j-1} + \alpha \psi_j + \beta_u \psi_{j+1}
         wrk[sx][1]=wi;}
for (sx=1; sx<=NX; sx++) // Copy new wave function back to psi
         for (s=0; s<=1; s++)
                                                                                                                                                                                                                                               \left[\varepsilon_n^+ = \frac{1}{2} \left| \exp\left(-\frac{i\Delta t}{n}(a+b)\right) + \exp\left(-\frac{i\Delta t}{n}(a-b)\right) \right|
                   psi[sx][s]=wrk[sx][s];
                                                                                \exp(-i\Delta t T_x) = U_x^{\text{(half)}} U_x^{\text{(full)}} U_x^{\text{(half)}} + O([\Delta t]^3) \qquad \left| \varepsilon_n^- = \frac{1}{2} \left[ \exp\left(-\frac{i\Delta t}{n}(a+b)\right) - \exp\left(-\frac{i\Delta t}{n}(a-b)\right) \right] \right|
```

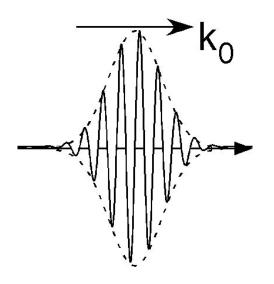
Initial Wave Function

• Gaussian wave packet

$$\psi(x, t = 0) = C \exp\left(-\frac{(x - x_0)^2}{4\sigma^2}\right) \exp(ik_0 x)$$

Symbol	Variable in qd1.c			
x_0 (packet center)	X0			
σ (packet spread)	S0			
$k_0^2/2$ (energy)	E0			





Quantum Dynamics—III Spectral Method

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Goal: Understand Fourier transform in the context of the orthonormal plane-wave basis set in a vector space

Resolution of identity:

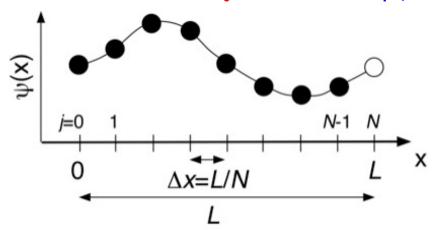


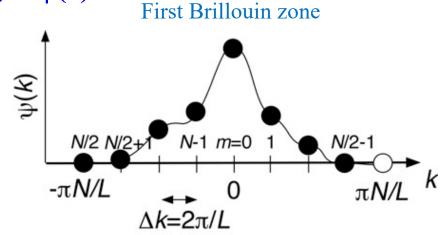
$$1 = \sum_{n} |n\rangle\langle n|$$



Discrete Fourier Transform

- Discretize $\psi(x) \in \mathbb{C}$ $(x \in [0, L])$ on N mesh points, $x_j = j\Delta x$ (j = 0, ..., N-1), with equal mesh spacing, $\Delta x = L/N$
- Periodic boundary condition: $\psi(x + L) = \psi(x)$



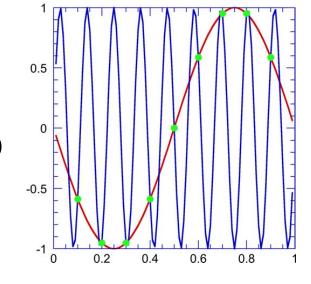


• Discrete Fourier transform: Represents $\psi(x)$ as a linear combination of $\exp(ikx) = \cos(kx) + i\sin(kx)$, with different wave numbers, k

$$\psi_{j} = \sum_{m=0}^{\infty} \tilde{\psi}_{m} \exp(ik_{m}x_{j})$$

$$k_{m} = \begin{cases} 2\pi m/L & (m = 0,1,...,N/2 - 1) \\ 2\pi (m - N)/L & (m = N/2,N/2 + 1,...,N - 1) \end{cases}$$

$$\tilde{\psi}_{m} = \frac{1}{N} \sum_{j=0}^{N-1} \psi_{j} \exp(-ik_{m}x_{j})$$



Wave Numbers

• Periodic boundary condition, $\psi(x+L) = \psi(x)$, is guaranteed by choosing $k_m = 2\pi m/L$

$$e^{i\left(\frac{2\pi m}{L}(\mathbf{x}+\mathbf{L})\right)} = e^{i\left(\frac{2\pi m}{L}\mathbf{x}+2\pi m\right)} = e^{i\frac{2\pi m}{L}\mathbf{x}} \underbrace{e^{i2\pi m}}_{\mathbf{k}} = e^{i\frac{2\pi m}{L}\mathbf{x}} = e^{i\frac{2\pi m}{L}\mathbf{x}}$$

• Folding back the latter half of wave numbers by $\frac{2\pi N}{L} = 2\pi/\Delta x$ $\Delta x = \frac{L}{N}$ (cf. first Brillouin zone):

From
$$\left[0, \frac{2\pi N}{L}\right] = \left[0, \frac{2\pi}{\Delta x}\right]$$
 to $\left[-\frac{2\pi}{L} \cdot \frac{N}{2}, \frac{2\pi}{L} \cdot \frac{N}{2}\right] = \left[-\frac{\pi}{\Delta x}, \frac{\pi}{\Delta x}\right]$

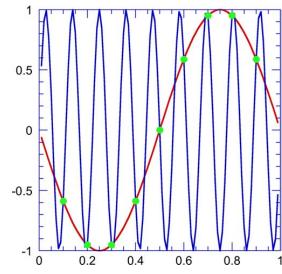
The shift won't change wave-function value on any grid point

$$e^{i\left(k_{m}-\frac{2\pi N}{L}\right)x_{j}}$$

$$=e^{ik_{m}x_{j}}e^{-i\frac{2\pi N}{L}\cdot\frac{L}{N}j}$$

$$=e^{ik_{m}x_{j}}e^{-i2\pi j}$$

$$=e^{ik_{m}x_{j}}$$



Orthonormal Basis Set

- *N*-dimensional vector space: $|\psi\rangle = (\psi_0, \psi_1, ..., \psi_{N-1})$
- Plane-wave basis set: $\{ |m\rangle = b_m(x_j) = \frac{1}{\sqrt{N}} \exp(ik_m x_j) | m = 0,1,...,N-1 \}$
- Orthonormality: $\langle m|n\rangle = \sum_{j=0}^{N-1} b_m^*(x_j)b_n(x_j) = \delta_{m,n} = \begin{cases} 1 & (m=n) \\ 0 & (m \neq n) \end{cases}$

$$\therefore \langle m | n \rangle = \frac{1}{N} \sum_{j=0}^{N-1} \exp(i(k_n - k_m)x_j) = \frac{1}{N} \sum_{j=0}^{N-1} \exp\left(i\frac{2\pi}{N}(n - m)j\right)$$

$$(k_n - k_m)x_j = \frac{2\pi(n - m)}{L} \cdot \frac{L}{N}j$$

$$= \begin{cases} \frac{1}{N} \frac{\exp(i2\pi(n-m))-1}{\exp(i\frac{2\pi}{N}(n-m))-1} = 0 & (m \neq n) \end{cases}$$

$$= \begin{cases} \frac{1}{N} \frac{\exp(i2\pi(n-m))-1}{\exp(i\frac{2\pi}{N}(n-m))-1} = 0 & (m \neq n) \end{cases}$$

$$S = \sum_{j=0}^{N-1} \left(e^{i\frac{2\pi}{N}(n-m)}\right)^{j} = 1 + \dots + r^{N-1}$$

$$\frac{1}{N} \cdot N = 1 \qquad (m = n) \end{cases}$$

$$\therefore (r-1)S = (r^{N}-1) \qquad |\psi\rangle = \sum_{m} c_{m}|m\rangle$$

$$\langle n|\times \psi \rangle$$

- Completeness: $|\psi\rangle = \sum_{m=0}^{N-1} |m\rangle\langle m|\psi\rangle$ or $1 = \sum_{m=0}^{N-1} |m\rangle\langle m|$ $\langle n|\psi\rangle = \sum_{m} c_m \stackrel{\delta_{nm}}{\langle n|m\rangle} = c_n$
- Fourier transform: $\psi_j = \sum_{m=0}^{N-1} \exp(ik_m x_j) \frac{1}{N} \sum_{l=0}^{N-1} \exp(-ik_m x_l) \psi_l$ $\tilde{\psi}_m$

Spectral Method

• Kinetic-energy operator is diagonal in the momentum space: $\tilde{\psi}_m \stackrel{T}{\to} \frac{k_m^2}{2} \tilde{\psi}_m$

$$-\frac{1}{2}\frac{\partial^2}{\partial x^2} \sum_{m=0}^{N-1} \tilde{\psi}_m \exp(ik_m x) = \sum_{m=0}^{N-1} \frac{k_m^2}{2} \tilde{\psi}_m \exp(ik_m x)$$

- Potential-energy operator is diagonal in the real space: $\psi_j \stackrel{v}{\to} V_j \; \psi_j$
- Split-operator technique & spectral method

$$\psi(t + \Delta t) = \exp\left(-\frac{iV\Delta t}{2}\right)^{-F} \exp(-iT\Delta t)^{-F-1} \exp\left(-\frac{iV\Delta t}{2}\right) \psi(t) + O\left((\Delta t)^{3}\right)$$

1.
$$\psi_j \xrightarrow{\exp(-iV\Delta t/2)} \exp(-iV_j\Delta t/2)\psi_j$$

2.
$$\psi_j \xrightarrow{F^{-1}} F^{-1} \psi_j = \tilde{\psi}_m = \frac{1}{N} \sum_{j=1}^N \psi_j \exp(-ik_m x_j)$$

3.
$$\tilde{\psi}_m \xrightarrow{\exp(-iT\Delta t)} \exp(-ik_m^2 \Delta t/2)\tilde{\psi}_m = \exp(\frac{i\Delta t}{2}\frac{\partial^2}{\partial x^2})\sum_m \tilde{\psi}_m e^{-ik_m x} = \sum_m \exp(-\frac{i\Delta t k_m^2}{2})\tilde{\psi}_m e^{-ik_m x}$$

4.
$$\tilde{\psi}_m \xrightarrow{F} F\tilde{\psi}_m = \psi_j = \sum_{m=1}^N \tilde{\psi}_m \exp(ik_m x_j)$$

5.
$$\psi_j \xrightarrow{\exp(-iV\Delta t/2)} \exp(-iV_j\Delta t/2)\psi_j$$

Exact exponentiation!

Numerical Recipes FFT: four1()

Spectral method requires

$$\psi_{j} \xrightarrow{F^{-1}} F^{-1}\psi_{j} = \tilde{\psi}_{m} = \frac{1}{N} \sum_{j=1}^{N} \psi_{j} \exp(-ik_{m}x_{j})$$

$$\tilde{\psi}_{m} \xrightarrow{F} F\tilde{\psi}_{m} = \psi_{j} = \sum_{m=1}^{N} \tilde{\psi}_{m} \exp(ik_{m}x_{j})$$

- four1(double data[], unsigned long nn, int isign)
 On input, the data[] array contains 2*nn elements that represent nn
 complex function values, such that data[2*j-1] & data[2*j] (j = 1, ..., nn) are
 the real & imaginary parts of the function value on the j-th grid point
- On output data[] is replaced by:

$$-isign = 1 data_{j} \leftarrow \sum_{m=1}^{N} data_{m} \exp(i2\pi mj/N) -isign = -1 data_{m} \leftarrow \sum_{i=1}^{N} data_{j} \exp(-i2\pi mj/N) k_{m}x_{j} = \frac{2\pi m}{L} \times \frac{L}{N}j = \frac{2\pi mj}{N}$$

• Note that four 1() does not perform the division by N in F^{-1}

See four1.c in the class home page

Using four1()

• Define double psi[2*N], where psi[2*j] & psi[2*j+1] (j = 0, ..., N-1) are the real & imaginary parts of ψ_i

```
/* Fourier transform */
four1(psi-1, (unsigned long) N, 1);
/* Inverse Fourier transform */
four1(psi-1, (unsigned long) N, -1);
for (int j=0; j<2*N; j++)
   psi[j] /= N;</pre>
```

• Note that four1() assumes 1 offset for the first argument but psi[] is 0 offset

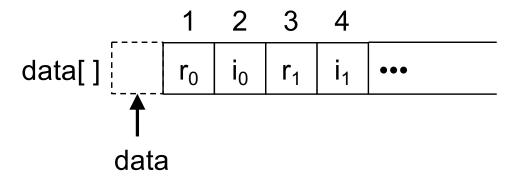
$$\psi_{j} \xrightarrow{F^{-1}} F^{-1}\psi_{j} = \tilde{\psi}_{m} = \frac{1}{N} \sum_{j=1}^{N} \psi_{j} \exp(-ik_{m}x_{j})$$

$$\tilde{\psi}_{m} \xrightarrow{F} F\tilde{\psi}_{m} = \psi_{j} = \sum_{m=1}^{N} \tilde{\psi}_{m} \exp(ik_{m}x_{j})$$

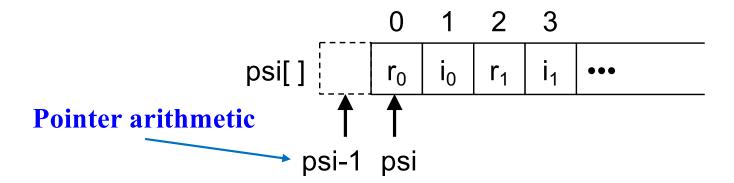
	0	1	2	3		2 <i>N</i> -2	2 <i>N</i> -1
psi[]	Re ψ_0	Im ψ_0	Re ψ_1	Im ψ_1	•••	${ m Re}\psi_{N-1}$	Im ψ_{N-1}

Array Offset

• four1() assumes 1-offset (because of its Fortran origin)



But psi[] uses 0-offset (C convention)



In C language, array name is a pointer to the zero-th element

Energy

$$\langle H \rangle = \langle T \rangle + \langle V \rangle$$

$$= \int dx \psi^*(x) \left(-\frac{1}{2} \frac{\partial^2}{\partial x^2} \right) \psi(x) + \int dx \psi^*(x) V(x) \psi(x)$$

$$\text{discretize} \quad \cong dx \sum_{j=0}^{N-1} \psi_j^* \left(-\frac{1}{2} \frac{\partial^2}{\partial x^2} \right) \psi_j + dx \sum_{j=0}^{N-1} \psi_j^* V_j \psi_j$$

$$= dx N \sum_{m=0}^{N-1} \frac{k_m^2}{2} \left| \tilde{\psi}_m \right|^2 + dx \sum_{j=0}^{N-1} V_j \left| \psi_j \right|^2 \text{ weighted sums}$$

In calc_energy():

- 1. $\tilde{\psi}_m \leftarrow F^{-1}[\psi_j]$
- 2. $E_{\text{kin}} \leftarrow dx N \sum_{m=0}^{N-1} \frac{k_m^2}{2} \left| \tilde{\psi}_m \right|^2$
- 3. $\psi_j \leftarrow F[\tilde{\psi}_m]$ // don't forget to come back
- 4. $E_{\text{pot}} \leftarrow dx \sum_{j=0}^{N-1} V_j |\psi_j|^2$

Kinetic Energy

$$\langle T \rangle = dx \sum_{j=0}^{N-1} \sum_{m=0}^{N-1} \tilde{\psi}_{m}^{*} \exp\left(-ik_{m}x_{j}\right) \left(-\frac{1}{2} \frac{\partial^{2}}{\partial x^{2}}\right) \sum_{n=0}^{N-1} \tilde{\psi}_{n} \exp\left(ik_{n}x_{j}\right)$$

$$= dx \sum_{j=0}^{N-1} \sum_{m=0}^{N-1} \tilde{\psi}_{m}^{*} \exp\left(-ik_{m}x_{j}\right) \sum_{n=0}^{N-1} \frac{k_{n}^{2}}{2} \tilde{\psi}_{n} \exp\left(ik_{n}x_{j}\right) \approx \frac{e^{-ik_{n}\Delta x} - 2 + e^{ik_{n}\Delta x}}{\Delta x^{2}} e^{ik_{n}x_{j}}$$

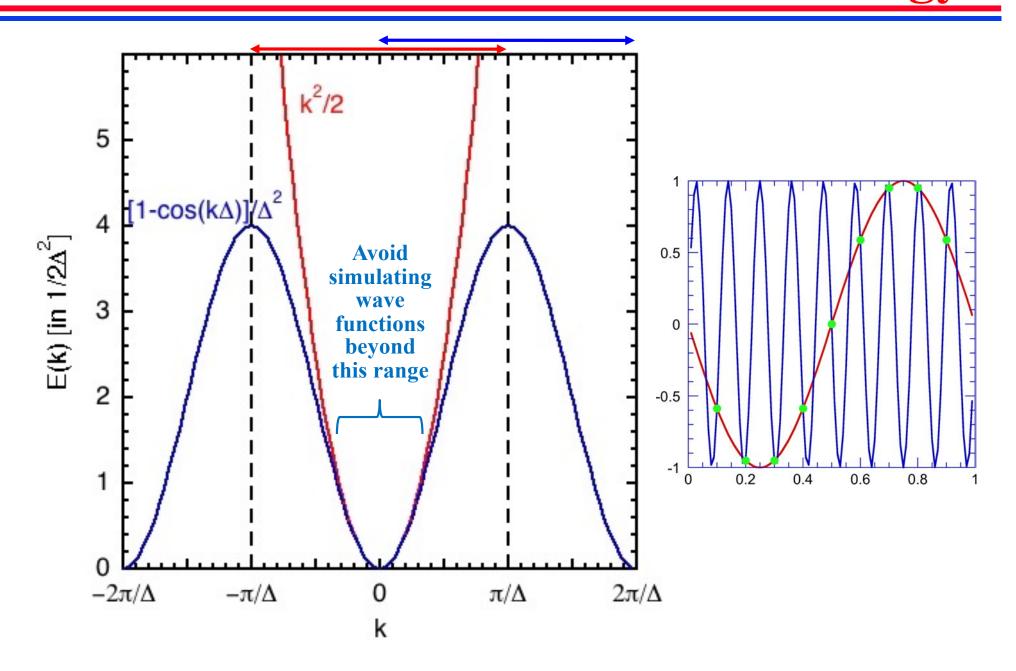
$$= dx \sum_{m=0}^{N-1} \tilde{\psi}_{m}^{*} \sum_{n=0}^{N-1} \frac{k_{n}^{2}}{2} \tilde{\psi}_{n} \sum_{j=0}^{N-1} \exp\left(i(k_{n} - k_{m})x_{j}\right) = \frac{2[\cos(k_{n}\Delta x) - 1]}{\Delta x^{2}} e^{ik_{n}x_{j}}$$

$$= dx \sum_{m=0}^{N-1} \tilde{\psi}_{m}^{*} \sum_{n=0}^{N-1} \frac{k_{n}^{2}}{2} \tilde{\psi}_{n} N\delta_{m,n}$$

$$= dx \sum_{m=0}^{N-1} \tilde{\psi}_{m}^{*} \sum_{n=0}^{N-1} \frac{k_{n}^{2}}{2} |\tilde{\psi}_{m}|^{2}$$

$$= dx N \sum_{m=0}^{N-1} \frac{k_{m}^{2}}{2} |\tilde{\psi}_{m}|^{2}$$

Continuum vs. Discrete Kinetic Energy



Total Energy Conservation

- Energy eigenvalues & eigenvectors: $H|n\rangle = \varepsilon_n|n\rangle$ (n = 0, ..., N-1)
- Time evolution of a wave function

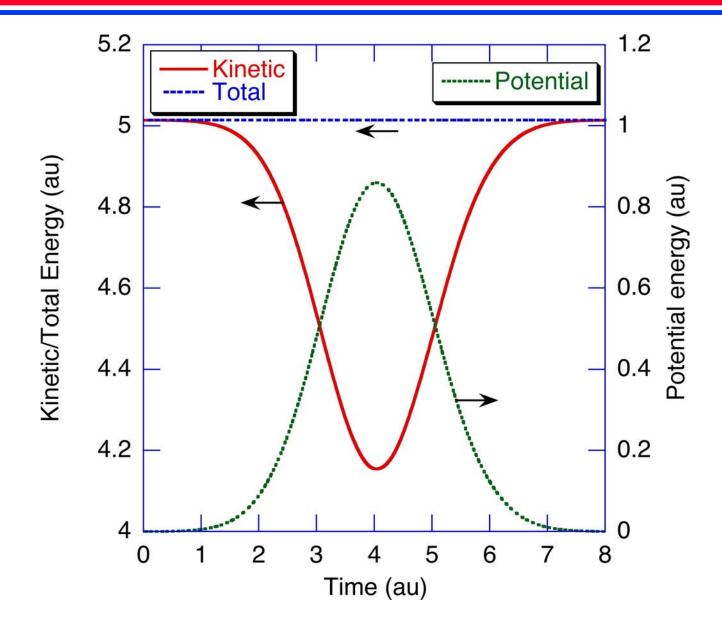
$$|\psi(t)\rangle = \exp(-iHt) \sum_{n=0}^{N-1} |n\rangle\langle n|\psi(0)\rangle = \sum_{n=0}^{N-1} \exp(-i\varepsilon_n t)|n\rangle\langle n|\psi(0)\rangle$$

= 1: completeness (resolution of identity)

Total energy

$$\begin{split} \langle \psi(t)|H|\psi(t)\rangle &= \left(\sum_{m=0}^{N-1} \langle \psi(0)|m\rangle \exp(i\varepsilon_m t)\langle m|\right) H\left(\sum_{n=0}^{N-1} \exp(-i\varepsilon_n t)|n\rangle\langle n|\psi(0)\rangle\right) \\ &= \left(\sum_{m=0}^{N-1} \langle \psi(0)|m\rangle \exp(i\varepsilon_m t)\langle m|\right) \left(\sum_{n=0}^{N-1} \exp(-i\varepsilon_n t)\varepsilon_n|n\rangle\langle n|\psi(0)\rangle\right) \\ &= \sum_{m=0}^{N-1} \sum_{n=0}^{N-1} \exp(i(\varepsilon_m - \varepsilon_n)t)\varepsilon_n\langle \psi(0)|m\rangle\langle n|\psi(0)\rangle\langle m|n\rangle \\ &= \sum_{n=0}^{N-1} \varepsilon_n|\langle n|\psi(0)\rangle|^2 = \text{constant} \end{split}$$

Energy Conservation for 1D Square Barrier



Energy conservation: good program verification

Initial Wave Function

Gaussian wave packet

$$\psi(x, t = 0) = C \exp\left(-\frac{(x - x_0)^2}{4\sigma^2}\right) \exp(ik_0 x)$$
SO E0 = $k_0^2/2$

$$x_0$$

$$i\frac{\partial}{\partial t}\exp\left(ik_0x - i\frac{k_0^2}{2}t\right) = \frac{k_0^2}{2}\exp\left(ik_0x - i\frac{k_0^2}{2}t\right)$$
Free-space solution

$$\left[-\frac{1}{2} \frac{d^2}{dx^2} + 0 \right] \exp\left(ik_0 x - i\frac{k_0^2}{2}t\right) = \frac{k_0^2}{2} \exp\left(ik_0 x - i\frac{k_0^2}{2}t\right)$$

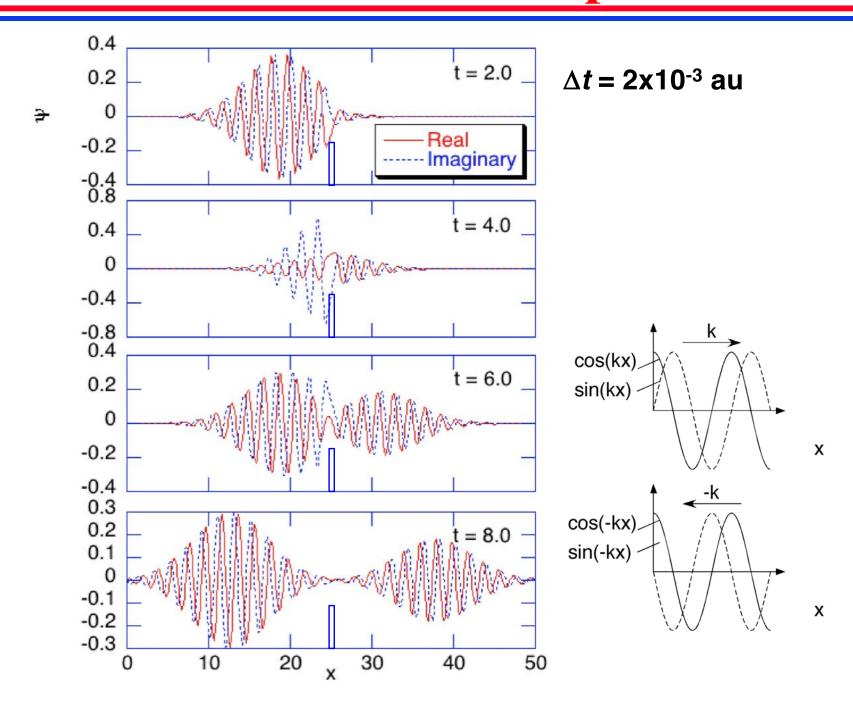
$$\psi(x,t) = \int dk \tilde{\psi}(k) \exp(ikx - i\omega(k)t) \quad \text{Here, } \omega(k) = k^2/2$$

$$= \int dk \tilde{\psi}(k) \exp\left(i \overset{k_0 + k - k_0}{\tilde{k}} x - i \overset{\cong \omega(k_0) + \frac{d\omega}{dk_0}(k - k_0)}{\tilde{\omega}(k)} t\right)$$

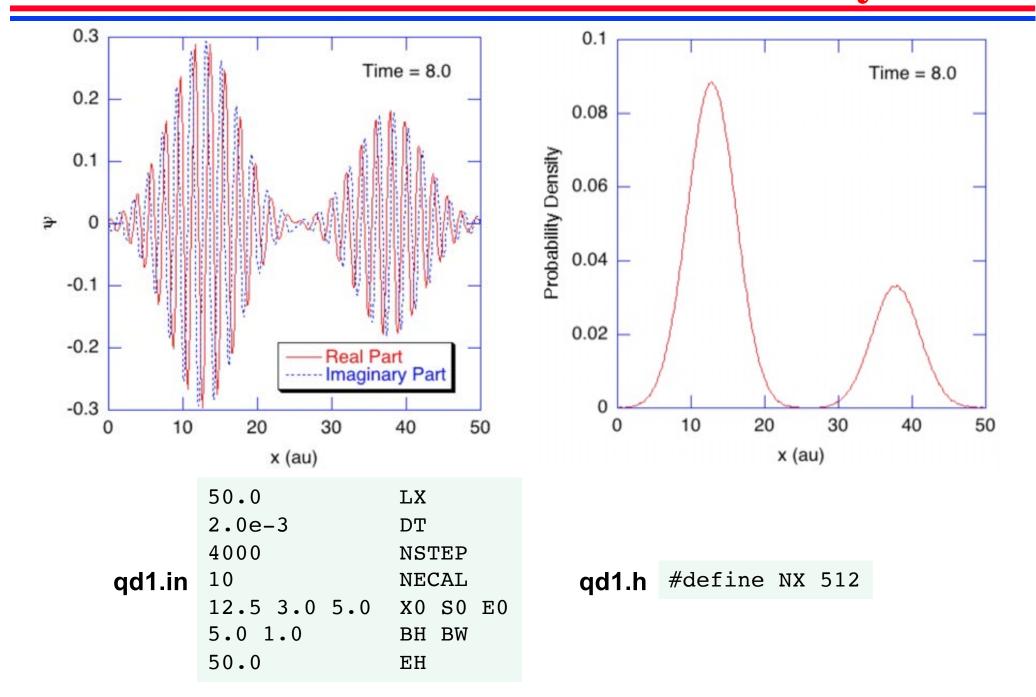
$$= \exp\left(ik_0 \left(x - \frac{\omega(k_0)}{k_0} t\right)\right) \int dk \tilde{\psi}(k) \exp\left(i(k - k_0) \left(x - \frac{d\omega}{dk_0} t\right)\right)$$
Thus evelocity

Group velocity

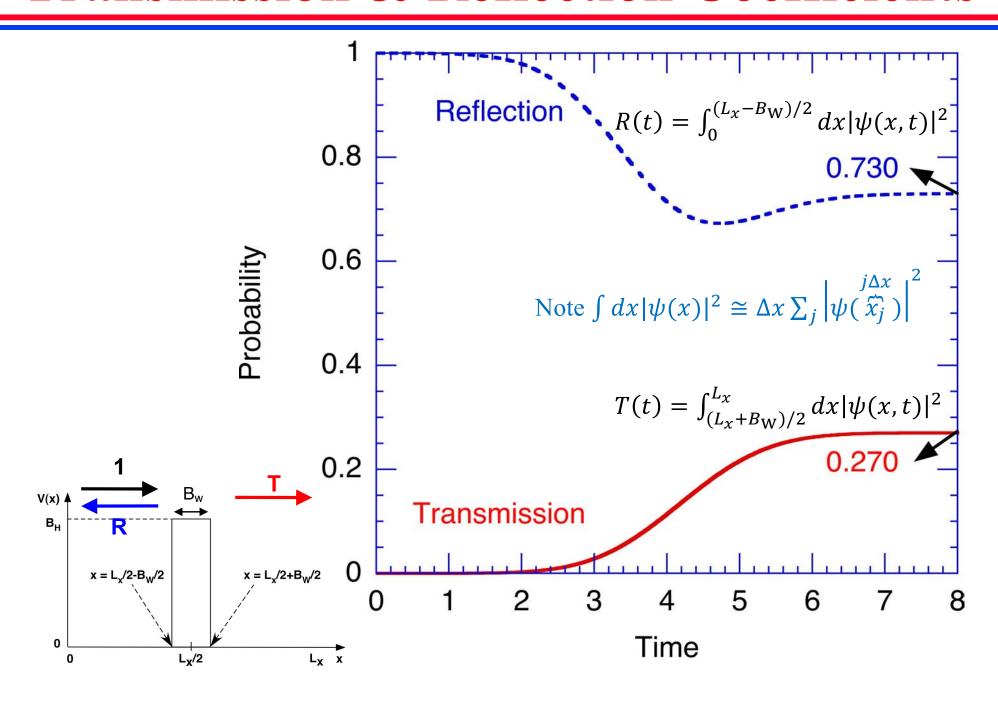
Numerical Example



Wave Function & Probability



Transmission & Reflection Coefficients



Top 10 Algorithms in History

In putting together this issue of *Computing in Science & Engineering*, we knew three things: it would be difficult to list just 10 algorithms; it would be fun to assemble the authors and read their papers; and, whatever we came up with in the end, it would be controversial. We tried to assemble the 10 algorithms with the greatest influence on the development and practice of science and engineering in the 20th century. Following is our list (here, the list is in chronological order; however, the articles appear in no particular order):

- Metropolis Algorithm for Monte Carlo
- Simplex Method for Linear Programming
- Krylov Subspace Iteration Methods
- The Decompositional Approach to Matrix Computations
- The Fortran Optimizing Compiler
- QR Algorithm for Computing Eigenvalues
- Quicksort Algorithm for Sorting
- Fast Fourier Transform
- Integer Relation Detection
- Fast Multipole Method

IEEE CiSE, Jan/Feb (2000)

PHYS 516 CSCI 596 CSCI 653

Fast Fourier Transform

• Danielson-Lanczos algorithm:

$$\psi_{j} = \sum_{m=0}^{N-1} \tilde{\psi}_{m} \exp(ik_{m}x_{j}) = \sum_{m=0}^{N-1} \tilde{\psi}_{m} \exp(i2\pi mj/N) \quad O(N^{2})!$$

$$= \sum_{m=0}^{N/2-1} \tilde{\psi}_{2m} \exp(i2\pi (2m)j/N) + \sum_{m=0}^{N/2-1} \tilde{\psi}_{2m+1} \exp(i2\pi (2m+1)j/N)$$

$$= \sum_{m=0}^{N/2-1} \tilde{\psi}_{2m} \exp(i2\pi mj/(N/2)) + \exp(i2\pi j/N) \sum_{m=0}^{N/2-1} \tilde{\psi}_{2m+1} \exp(i2\pi mj/(N/2))$$

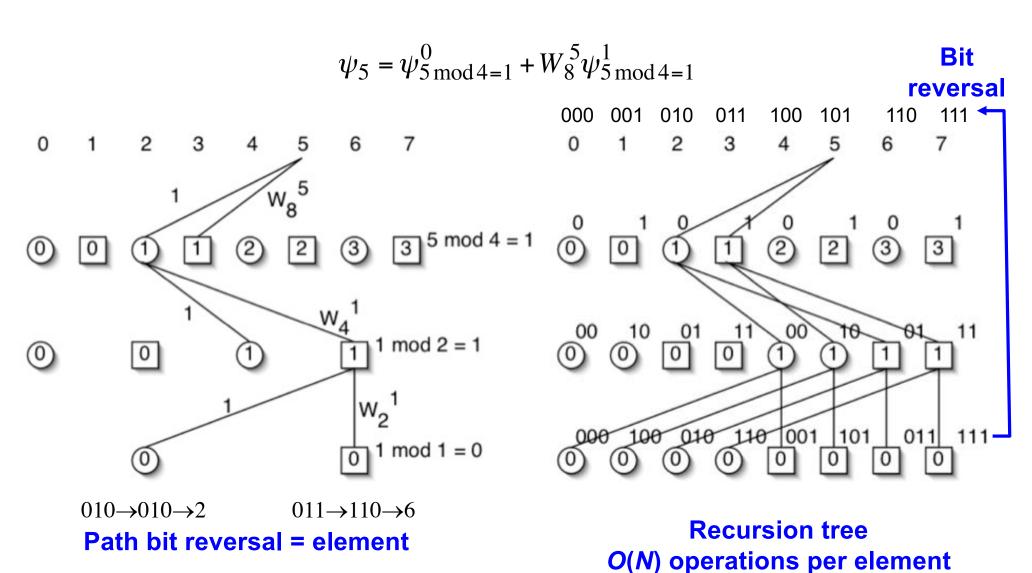
$$= \sum_{m=0}^{N/2-1} \tilde{\psi}_{2m} \exp(i2\pi mj/(N/2)) + \exp(i2\pi j/N) \sum_{m=0}^{N/2-1} \tilde{\psi}_{2m+1} \exp(i2\pi mj/(N/2))$$

$$\psi_j^0 = \sum_{m=0}^{N/2-1} \widetilde{\psi}_{2m} \exp(i2\pi mj/(N/2)) \xrightarrow{\text{subarray Fourier decompositions}} \psi_j = \psi_j^0 + W_N^j \psi_j^1 \qquad \begin{cases} \psi_j^1 = \sum_{m=0}^{N/2-1} \widetilde{\psi}_{2m+1} \exp(i2\pi mj/(N/2)) \\ w_N = \exp(i2\pi/N) \end{cases} \text{ j read as } j \bmod N/2$$

Divide-and-conquer

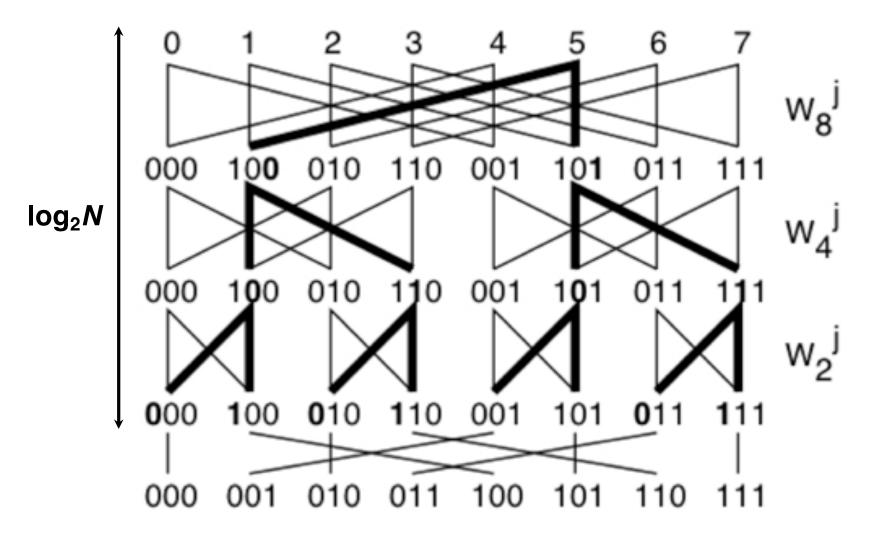
Fast Fourier Transform

• Recursive sub-Fourier transforms: $\psi_j = \psi_j^0 + W_N^j \psi_j^1$



Fast Fourier Transform Algorithm

• Butterfly (hypercube) data exchange after bit-reversal:

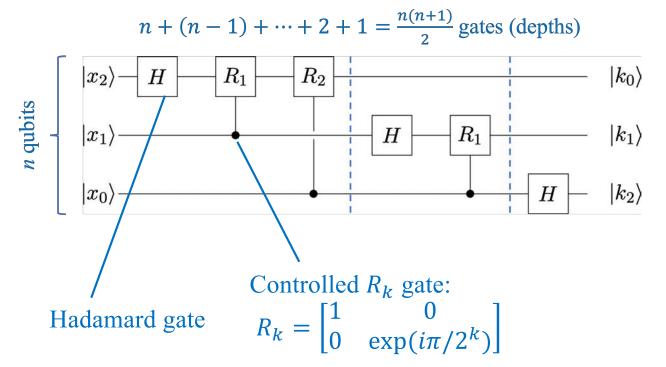


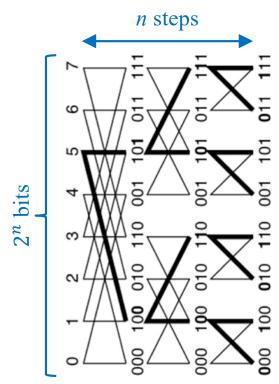
- Many computations are shared among the recursion trees
- 2Nlog₂N arithmetic operations

Quantum Fourier Transform (QFT)

• On a quantum computer, quantum parallelism allows Fourier transform to be performed using only n qubits & $O(n^2)$ gates for $N=2^n$

QFT:
$$|x\rangle \to \frac{1}{\sqrt{N}} \sum_{k=0}^{N-1} e^{i\frac{2\pi kx}{N}} |k\rangle$$





• Compare QFT with $2N\log_2 N = O(2^n n)$ arithmetic operations in classical fast Fourier transform (FFT):

Exponential operation reduction: $O(2^n n/n^2) = O(2^n/n)$

Preskill, arXiv:2106.10522 ('21)

Parallelizing Quantum Dynamics

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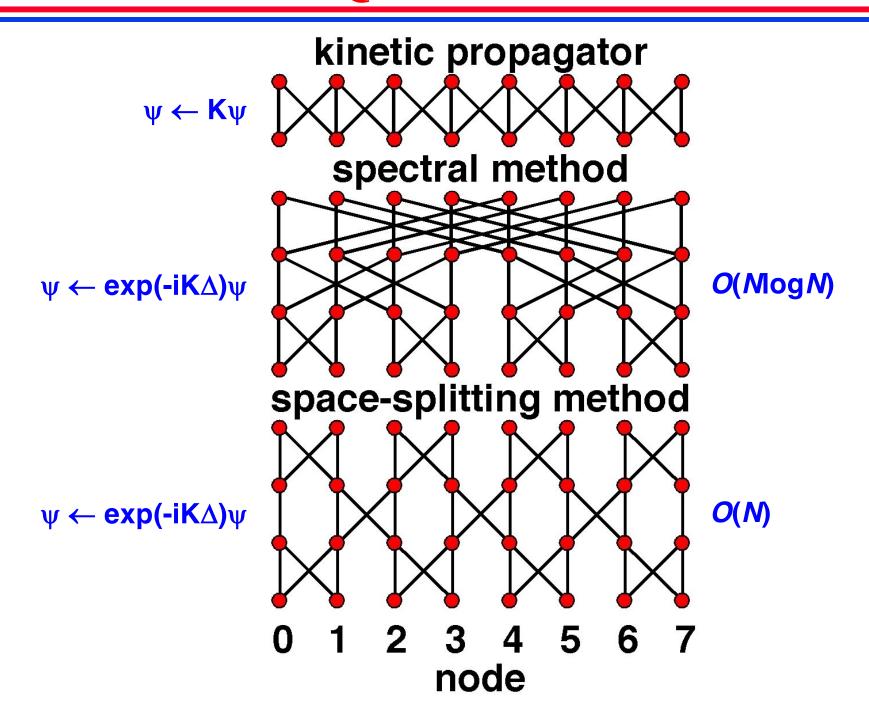
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See https://aiichironakano.github.io/cs596.html



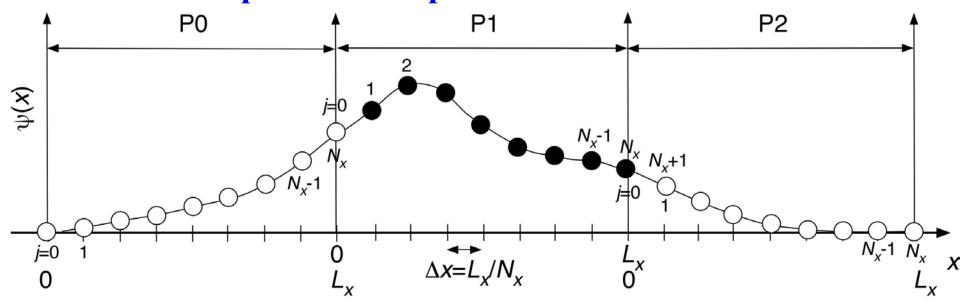


Parallel QD Communications



Parallelization of Space Splitting Method

• Self-centric spatial decomposition



Local & global coordinates

$$\begin{cases} x_j = j\Delta x \\ x_j^{\text{(global)}} = j\Delta x + pL_x \end{cases}$$

Global coordinates only in init_prop() & init_wavefn()

Boundary Wave Function Caching

• Parallelized periodic_bc()

(X) ↑

```
plw = (myid-1+nproc)%nproc; /* Lower partner process */
  pup = (myid+1)
                        )%nproc; /* Upper partner process */
  /* Cache boundary wave function value at the lower end */
  dbuf[0:1] \leftarrow psi[NX][0:1];
  Send dbuf to pup;
  Receive dbufr from plw;
  psi[0][0:1] \leftarrow dbufr[0:1];
  /* Cache boundary wave function value at the upper end */
  dbuf[0:1] \leftarrow psi[1][0:1];
  Send dbuf to plw;
  Receive dbufr from pup;
  psi[NX+1][0:1] \leftarrow dbufr[0:1];
                                                        P2
           P0
                                  P1
j=0
                              \Delta x = L / N_{y}
```