

Using Matlab functions

1. Create a text file with extension ".m" in your path
2. In the first line, define
function <output> = <function name>(<input>)
3. In subsequent lines, calculate output.

Example: a text file called 'genH.m':

```
function H=genH(N)
H=diag(-2*ones(N,1))+diag(ones(N-1,1),1)+diag(ones(N-1,1),-1);
```

when called at the Matlab command prompt,
we can now generate arbitrarily large matrices w/
desired structure $\longrightarrow$

```
>> genH(4)
```

```
ans =
```

```
-2  1  0  0
 1 -2  1  0
 0  1 -2  1
 0  0  1 -2
```

```
>> genH(5)
```

```
ans =
```

```
-2  1  0  0  0
 1 -2  1  0  0
 0  1 -2  1  0
 0  0  1 -2  1
 0  0  0  1 -2
```

"Diagonalization"

$\vec{X}_i$'s are orthonormal eigenvectors of Hermitian matrix $\vec{H}$ w/ eigenvalues λ_i .

What is $\vec{X}_i^\dagger \vec{H} \vec{X}_j$ ($i, j = 1, 2, \dots$)?

Since $\vec{H} \vec{X}_j = \lambda_j \vec{X}_j$, $\vec{X}_i^\dagger \vec{H} \vec{X}_j = \vec{X}_i^\dagger \lambda_j \vec{X}_j = \lambda_j \vec{X}_i^\dagger \vec{X}_j = \lambda_j \delta_{ij} = \begin{bmatrix} \lambda_1 & 0 & \dots \\ 0 & \lambda_2 & 0 \dots \\ & & \ddots \end{bmatrix}$

So eigenvectors diagonalize the matrix!

Example

$\vec{H} = \begin{bmatrix} 0 & 1 \\ 1 & 0 \end{bmatrix}$ eigenvalue: $\lambda_1 = 1$ $\lambda_2 = -1$
eigenvector: $\vec{X}_1 = \frac{1}{\sqrt{2}} \begin{bmatrix} 1 \\ 1 \end{bmatrix}$ $\vec{X}_2 = \frac{1}{\sqrt{2}} \begin{bmatrix} 1 \\ -1 \end{bmatrix}$

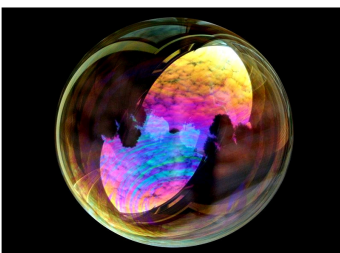
$i=1, j=1$: $\vec{X}_1^\dagger \vec{H} \vec{X}_1 = \frac{1}{2} \begin{bmatrix} 1 & 1 \end{bmatrix} \begin{bmatrix} 0 & 1 \\ 1 & 0 \end{bmatrix} \begin{bmatrix} 1 \\ 1 \end{bmatrix} = \frac{1}{2} \begin{bmatrix} 1 & 1 \end{bmatrix} \begin{bmatrix} 1 \\ 1 \end{bmatrix} = 1$

$i=1, j=2$: $\vec{X}_1^\dagger \vec{H} \vec{X}_2 = \frac{1}{2} \begin{bmatrix} 1 & 1 \end{bmatrix} \begin{bmatrix} 0 & 1 \\ 1 & 0 \end{bmatrix} \begin{bmatrix} 1 \\ -1 \end{bmatrix} = \frac{1}{2} \begin{bmatrix} 1 & 1 \end{bmatrix} \begin{bmatrix} -1 \\ 1 \end{bmatrix} = 0$

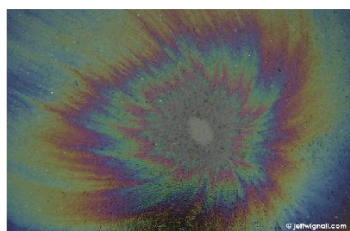
etc. . .

Matter Waves: Interference

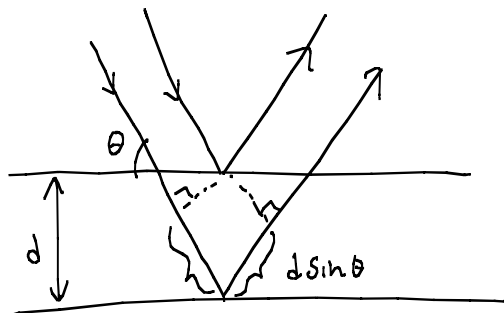
EM waves: "Fabry-Perot"



soap bubble



oil slick



path length difference = $2d \sin \theta = n\lambda$ $n = 1, 2, 3, \dots$ for Constructive interference.

"Bragg's Law" $\rightarrow$ X-rays for $d \sim$ atomic spacing

Electron waves: Davisson-Germer experiment (Nobel Prize, 1937)

De Broglie: $p = \frac{h}{\lambda} \rightarrow \lambda = \frac{h}{p} = \frac{h}{\sqrt{2mE}}$ $\left(E = \frac{1}{2}mv^2 = \frac{m^2v^2}{2m} = \frac{p^2}{2m} \right)$

graphite
demo!



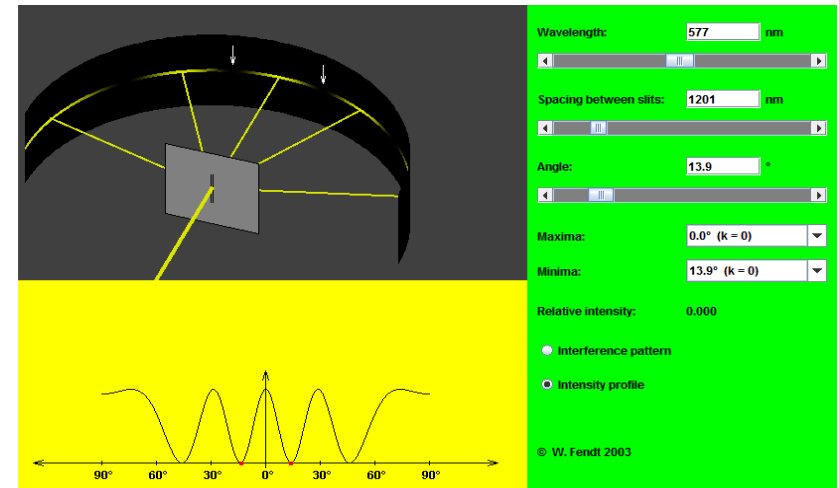
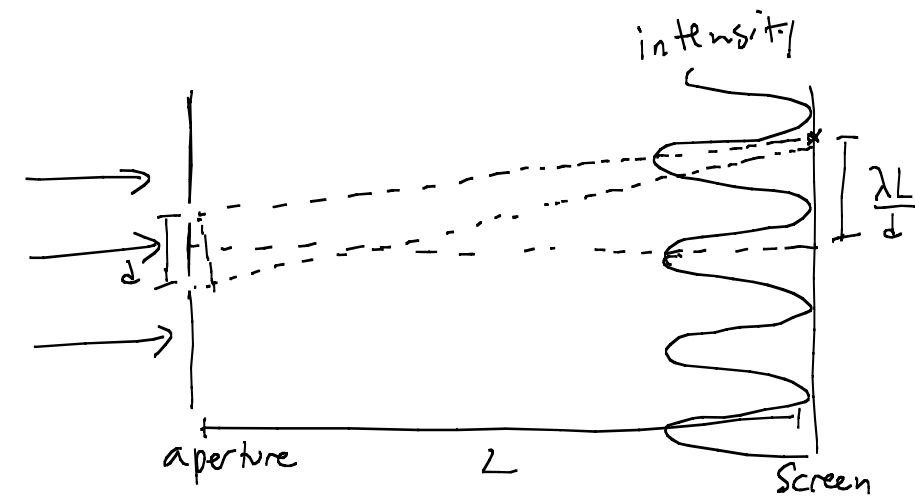
From Bragg's Law,

$$d = \frac{n\lambda}{2 \sin \theta} = \frac{nh}{2 \sin \theta \sqrt{2mE}} \sim \frac{h}{2^{1/10} \sqrt{2m} 10^3 \text{ eV}} = \frac{5.4 \times 10^{-15} \text{ eV} \cdot \text{s}}{\sqrt{2 \cdot 5 \times 10^{-16} \text{ eV} \frac{\text{s}^2}{\text{cm}^2} \cdot 10^3 \text{ eV}}} \sim 2 \text{ \AA}$$

atomic scale spacing!
C.f. actual $\sim 3 \text{ \AA}$! ✓

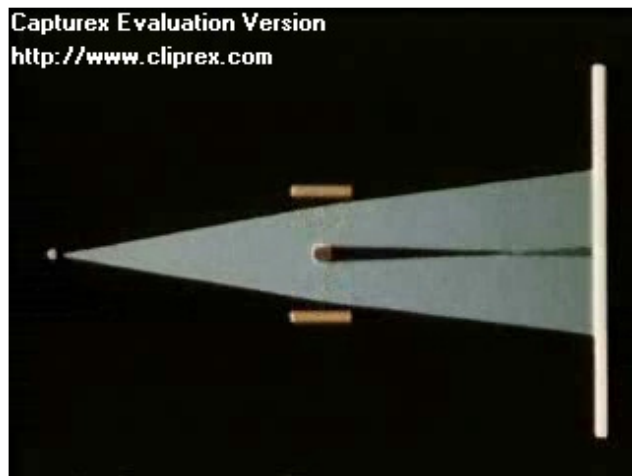
$E = mc^2 \rightarrow m = E/c^2$ (mass of electron is $511 \text{ KeV}/c^2$)

EM waves: Young's two-slit expt



<http://www.walter-fendt.de/ph14e/doubleslit.htm>

electrons: Biprism interference

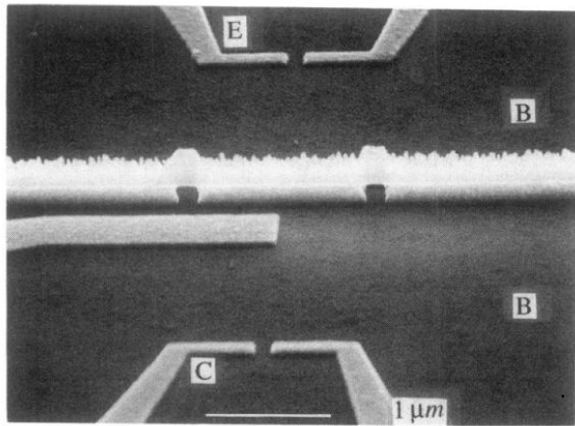


electron
fringes
↘

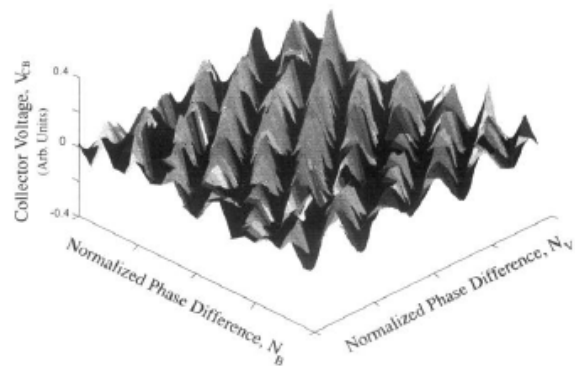


Electron Interferometry: Solid State (semiconductors)

two-slit:



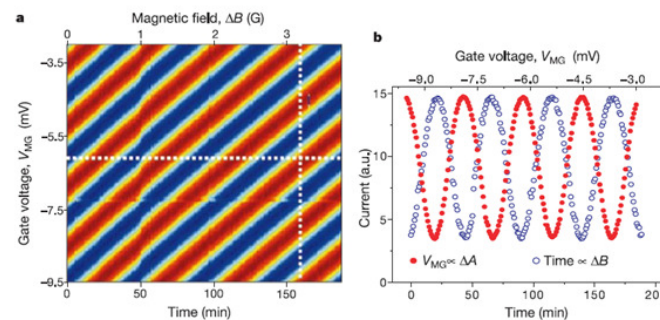
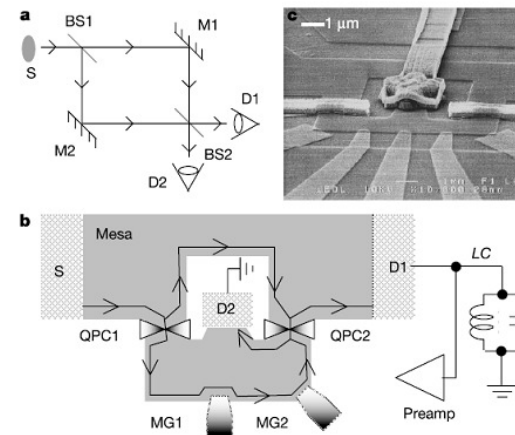
A. Yacoby, M. Heiblum, V. Umansky, H. Shtrikman, and D. Mahalu
• Phys. Rev. Lett. 73, 3149 (1994)



An electronic Mach-Zehnder interferometer

Yang Ji, Yunchul Chung, D. Sprinzak, M. Heiblum, D. Mahalu & Hadas Shtrikman

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Potential Exam Topics

Ordinary diff. Eqs.

Partial diff. Eqs.

Separation of Variables

Complex variables

Taylor series

Fourier Transforms

Orthogonality of fns
and vectors (incl. inner product)

"basis" functions and vectors

delta fns

Gaussians

Uncertainty rel'n

Boundary conditions

Finite differences

Differential eigenvalue/eigenfunction problem

Matrix eigenvalue/eigenvector problem

Hermitian matrices