

## Physics of Quantum dot junction transistor



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## Applications of QD devices



- Single-electron transistor (SET)
- Single-photon generator (SPG)
- Light-emitting transistor
- Optical signal control (slow light)
- QD lasers & detectors
- Non-linear optical devices
- High-density memory
- Biosensors
- Spintronics devices

## Theoretical approaches for one-particle states



- Envelope Function approximation
- K.p theory --  
Effective-mass model (for spherical conduction Band)  
Luttinger model (for degenerate valence bands)
- Effective bond-orbital model
- VFF model (for strain distribution)
- Tight-binding model
- Density-functional theory (DFT)

## Physical effects



Quantum confinement (size effect)

Selection rule (light polarization)

Discrete energy levels (0-d density of states)

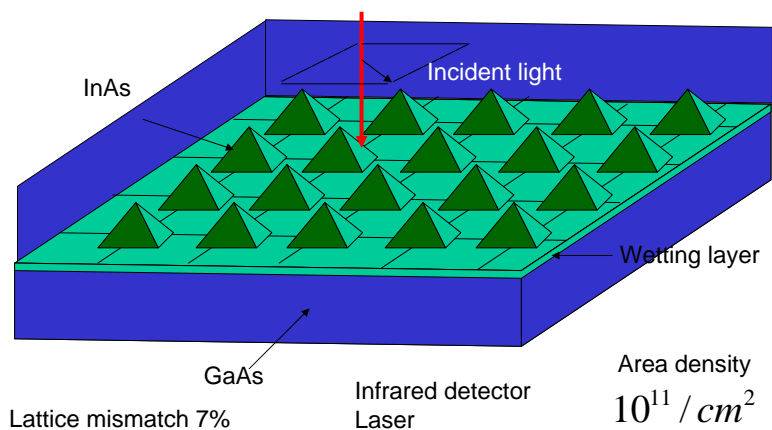
Strong Coulomb interaction (Coulomb blockade)

Strong electron correlation (many-body effect)



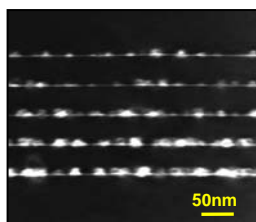
Artificial atom

## Self assembled quantum dots

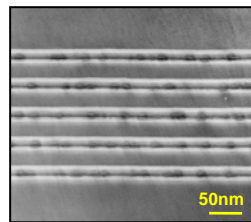


## Optical and Structural Characterization of QDIPs

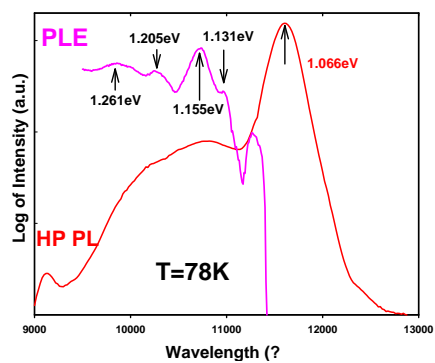
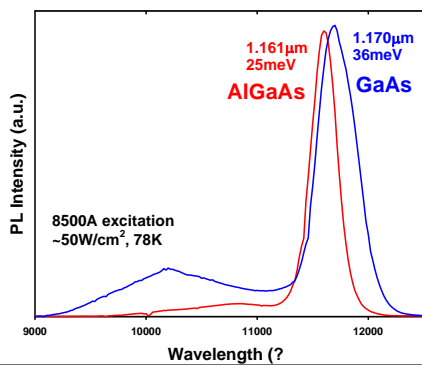
TEM GaAs



TEM AlGaAs



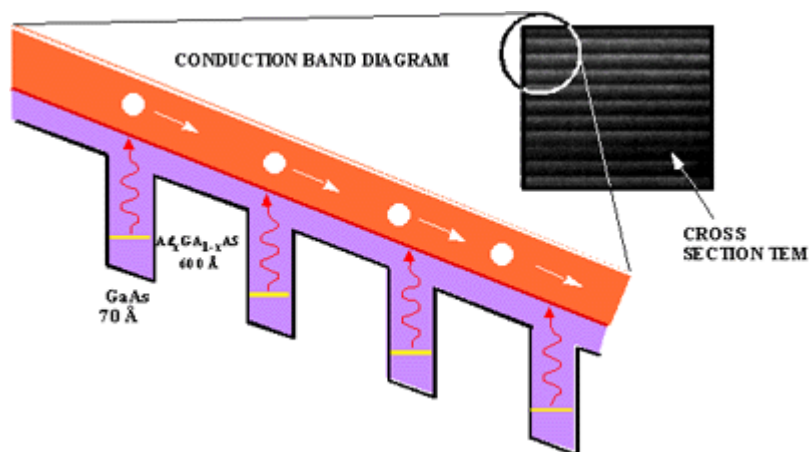
Energies of transitions  
(from PL/PLE data):  
65meV, 89meV  
139meV, 195meV



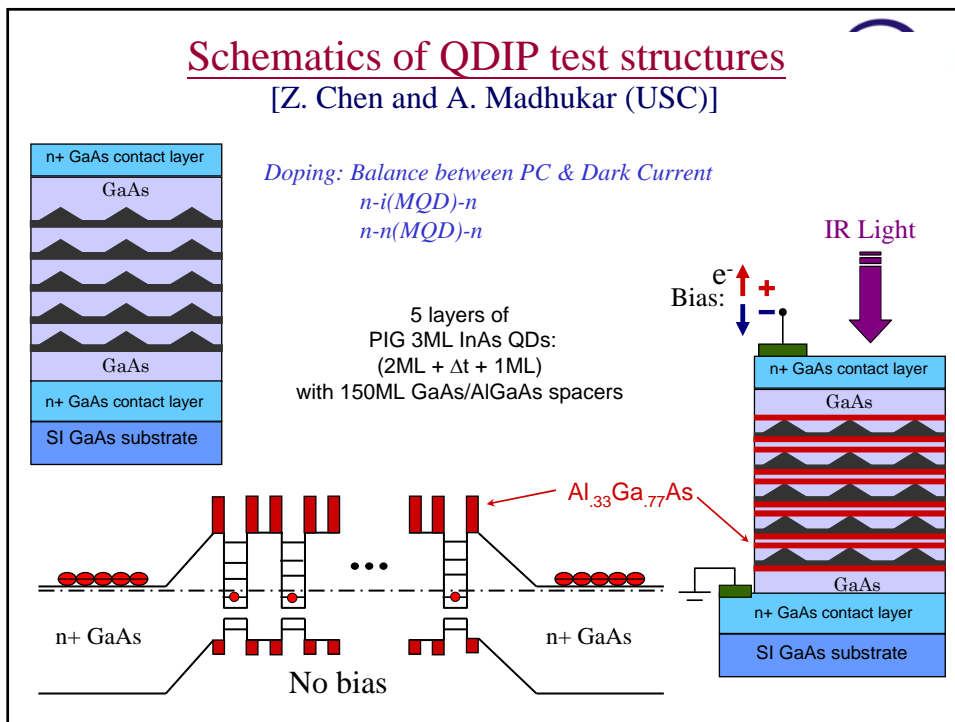
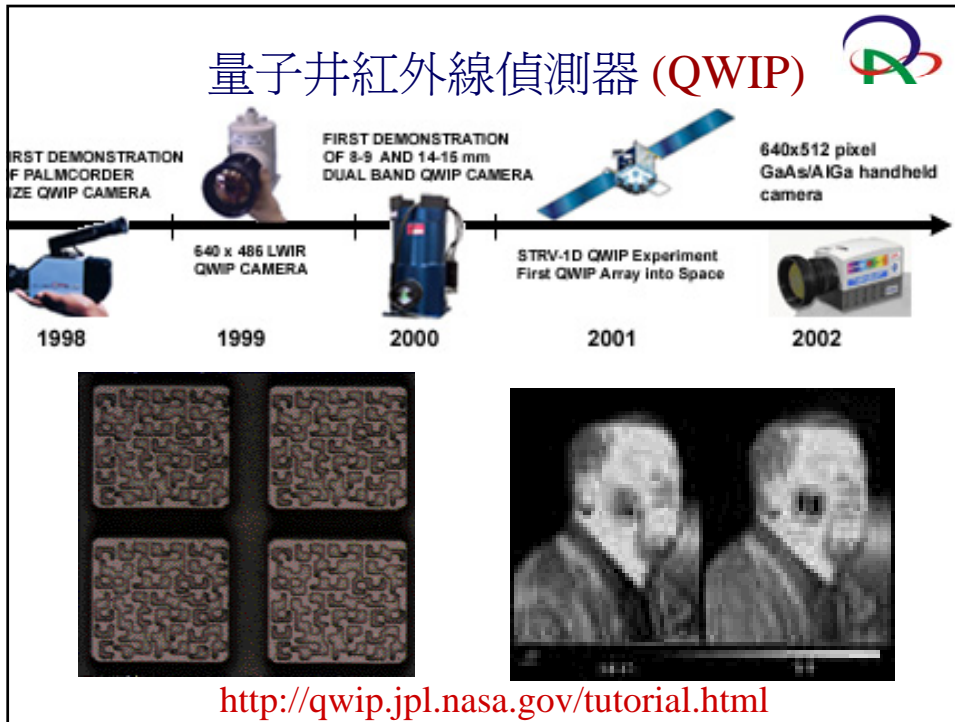
## 紅外線偵測器 (Infrared detector)



## 量子井紅外線偵測器(QWIP)



<http://qwip.jpl.nasa.gov/tutorial.html>

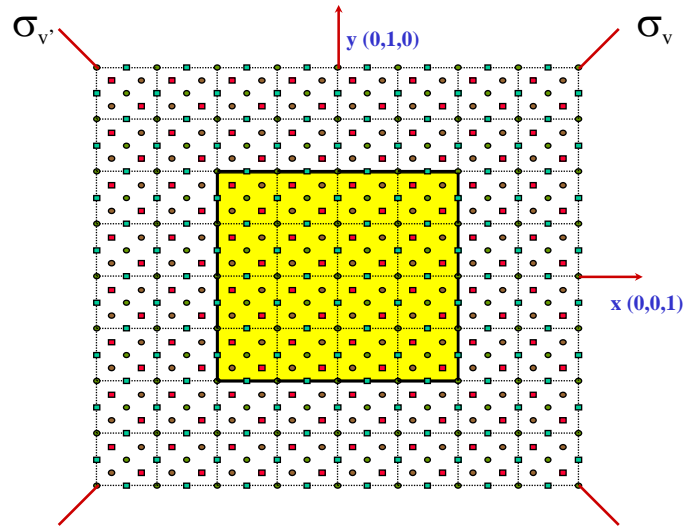




## Microscopic modeling



[S. Sun, Y. C. Chang, PRB 62, 13631 (2000)]



## Valence force field (VFF) Model



$$V = \frac{1}{4} \sum_{ij} \frac{3}{4} \alpha_{ij} (d_{ij}^2 - d_{0,ij}^2)^2 / d_{0,ij}^2$$

$$+ \frac{1}{4} \sum_i \sum_{j \neq k} \frac{3}{4} \beta_{ijk} (\bar{d}_{ij} \cdot \bar{d}_{ik} + d_{0,ij} d_{0,ik} / 3)^2 / d_{0,ij} d_{0,ik}$$

$i$  labels atom positions

$j, k$  label nearest-neighbors of  $i$

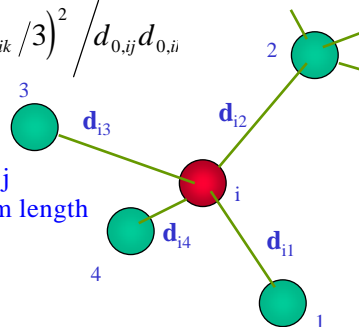
$d_{ij}$  = bond length joining sites  $i$  and  $j$

$d_{0,ij}$  is the corresponding equilibrium length

$\alpha_{ij}$  = bond stretching constants

$\beta_{ijk}$  = bond bending constants

We take  $d_{ijk}^2 = d_{ij} d_{ik}$





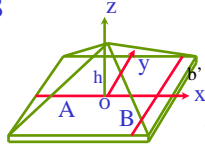
## Strain Potential Along Line A and B of Dot 1



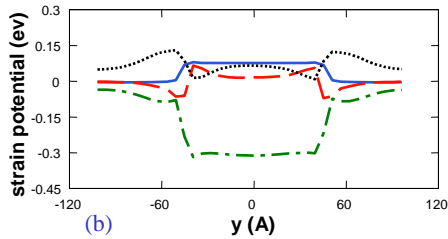
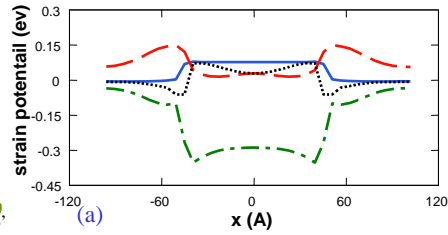
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- **Fig.(a):**  
strain potential  
along line A
- **Fig.(b):**  
strain potential  
along line B

—  $V_{ss}$   
- - -  $V_{xx}$   
⋯  $V_{yy}$   
- · -  $V_{zz}$



- **Band offset:**  
CB: 0.833eV  
VB: 0.260eV



## Strain Distribution Along [001] of Dot 1 and Dot 2



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- **Hydrostatic Strain :**

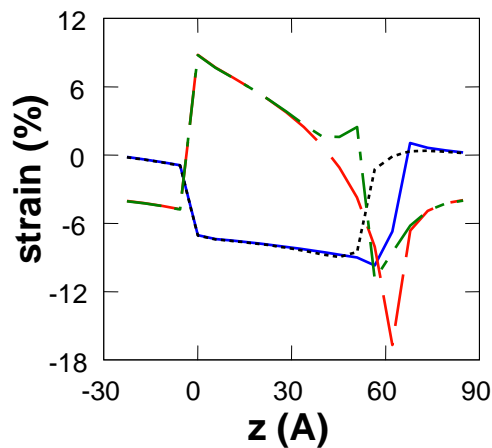
Dot 1: ———  
Dot 2: ⋯⋯⋯

- **Biaxial Strain:**

Dot 1: - - - -  
Dot 2: - · - -

- **[001] is growth direction**

- **Dot 1: 0-62.5 Å**  
**Dot 2: 0-50 Å**



# Comparison with continuum Model



J. Appl. Phys. 104, 083524 (2008)

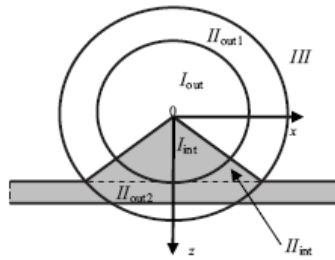


FIG. 1. The schematic representation of the QD-WL system (in gray) in a plan containing OZ axes. The specific regions  $I_{int}$ ,  $I_{ext}$ ,  $II_{int}$ ,  $II_{ext}$ , and III are shown.

using Green's function method one obtains (see Refs. 8–10)

$$\mathbf{u}(\mathbf{r}) = \frac{C}{\pi} \oint \frac{1}{|\mathbf{r} - \mathbf{r}_0|} dS(\mathbf{r}_0), \quad (1)$$

where  $C = \varepsilon_d(1 + \nu)/[4(1 - \nu)]$ ,  $\nu$  is Poisson's ratio,<sup>13</sup>  $\varepsilon_d$  is the interface dilation defined as  $\varepsilon_d = (a_{int} - a_{ext})/a_{ext}$  ( $a_{ext}$ ,  $a_{int}$  are the lattice constant of the external, internal (inclusion) material, respectively) and it coincides with the *compression*

083524-4 T. O. Checha and Y.-C. Chang

J. Appl. Phys. 104, 083524 (2008)

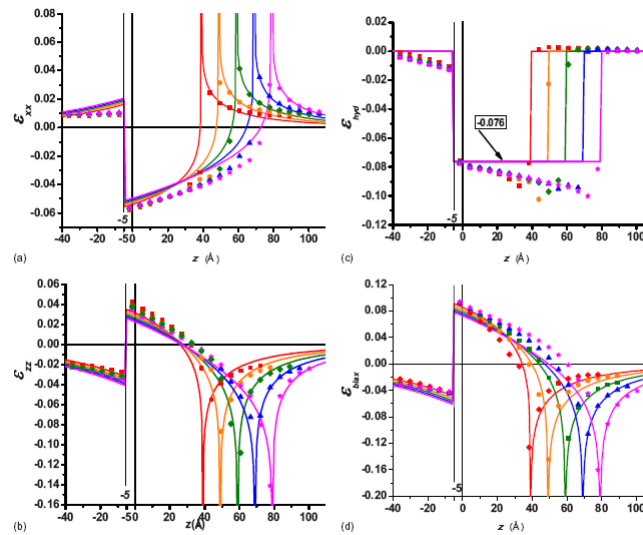


FIG. 3. (Color online) The strains on OZ axes: (a) the radial strain  $\varepsilon_{xx}$ , (b) the strain  $\varepsilon_{zz}$ , (c) the hydrostatic  $\varepsilon_{hyd}$ , and (d) the biaxial strain  $\varepsilon_{biax}$  on the OZ axis, for the ratios  $R/H=10/3.9, 10/4.9, 10/5.9, 10/6.9, 10/7.9$ , with  $R=100$  Å. The dots are for the VFF model and the continuum lines for CM.



## Piezoelectric Potential

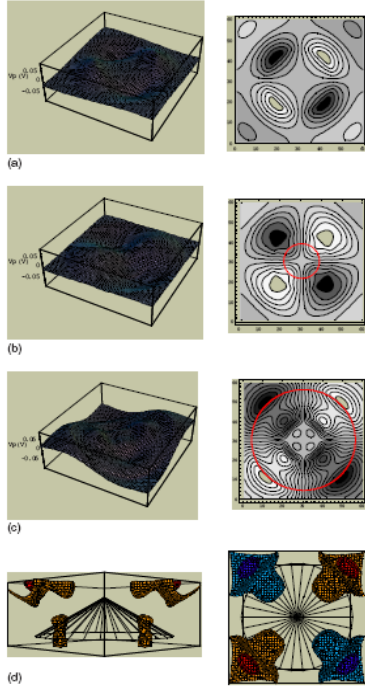


FIG. 5. (Color online) The piezoelectric potential  $V_p$ : (a) in the horizontal plane,  $z=0$  containing the top of the cone; (b) in the horizontal plane  $z=H/3$ ; (c) in the horizontal plane  $z=H$ . The left side figures are the three-dimensional plots and the right side figures are the contour plots. The red circles in these plots represent the cross sections of the QD with the horizontal planes. (d) equipotential surfaces of piezoelectric potential  $V_p$  for 0.065 V (red), -0.065 V (dark blue), 0.04 V (orange), and -0.04 V (blue). In the left side figure only the positive equipotential surfaces are plotted. In the right side figure are plots with a top view.



## Effective bond-orbital Model [Y.C.Chang, PRB 37, 8215 (1988)]



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$$H_{\alpha, \alpha'}(\vec{k}) = E_p \delta_{\alpha, \alpha'} +$$

$$\sum_{\tau} e^{i\vec{k} \cdot \vec{\tau}} \left\{ E_{xy} \tau_{\alpha} \tau_{\alpha'} + \left[ (E_{xx} - E_{xy}) \tau_{\alpha}^2 + E_{zz} (1 - \tau_{\alpha}^2) \right] \delta_{\alpha, \alpha} \right\}$$

### Strain Hamiltonian

$$H_{st} = \begin{pmatrix} -\Delta V_H + D_1 & \sqrt{3}d e_{xy} & \sqrt{3}d e_{xz} \\ \sqrt{3}d e_{xy} & -\Delta V_H + D_2 & \sqrt{3}d e_{yz} \\ \sqrt{3}d e_{xz} & \sqrt{3}d e_{yz} & -\Delta V_H + D_3 \end{pmatrix}$$

$$e_{ij} = (\epsilon_{ij} + \epsilon_{ji})/2$$

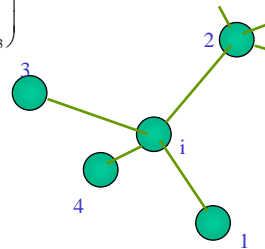
$$\Delta V_H = (a_1 + a_2)(\epsilon_{xx} + \epsilon_{yy} + \epsilon_{zz})$$

$$D_1 = b(2\epsilon_{xx} - \epsilon_{yy} - \epsilon_{zz})$$

$$D_2 = b(2\epsilon_{yy} - \epsilon_{xx} - \epsilon_{zz})$$

$$D_3 = b(2\epsilon_{zz} - \epsilon_{xx} - \epsilon_{yy})$$

$a_1, a_2, b, d$  = deformation potentials.



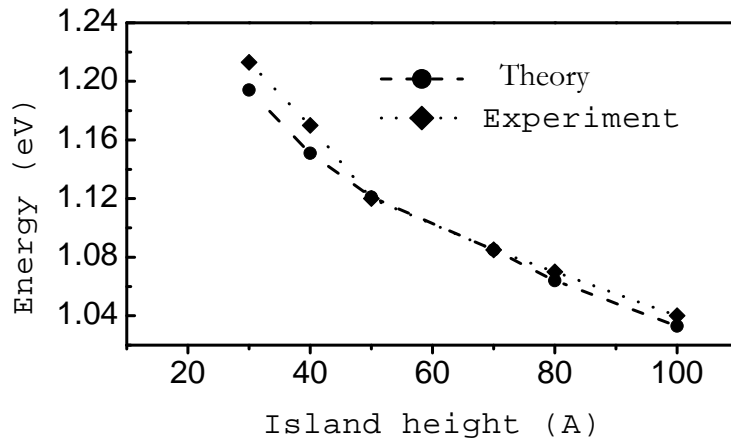


## Ground Transition Energy Varying With Dot Height (comparing to Experiment)

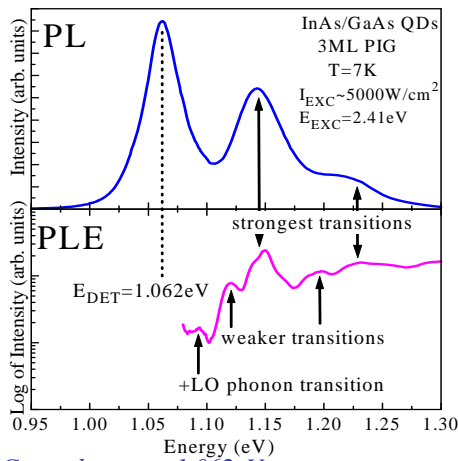


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Dot base length 200Å



## PL/PLE Characterization: Electronic Structure

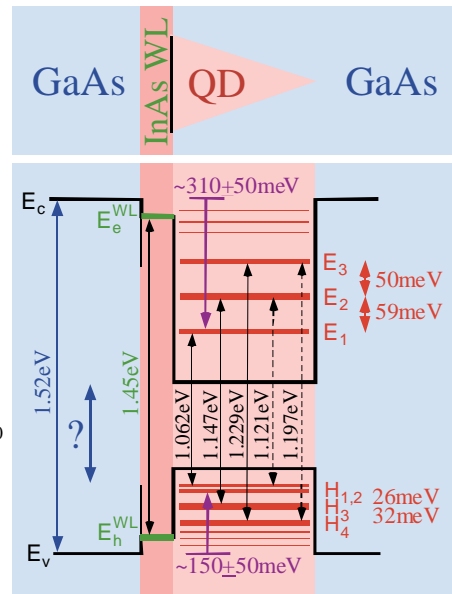


Ground state at 1.062eV

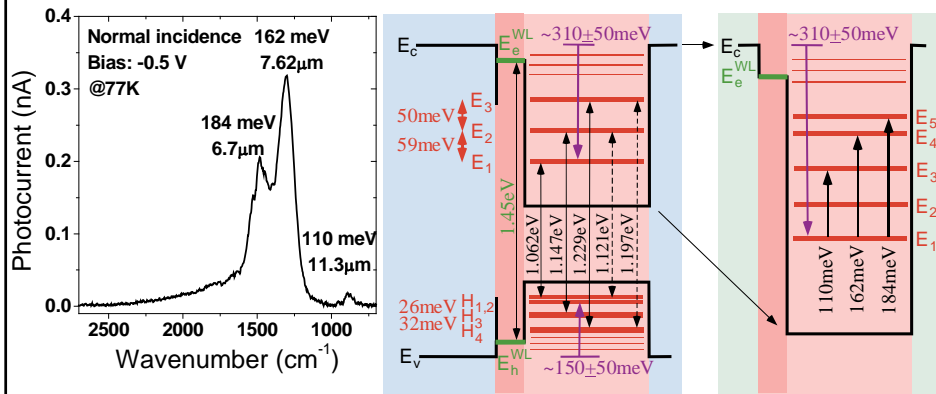
Excited states:

Strongest at 1.147eV and 1.229eV

Weaker at 1.121eV and 1.197eV



## Intra-band Transitions



A. Madhkar (USC)



## Intra-band Transitions



Table 4 Inter-sub band transition matrix elements of ground electron state to upper three electron states,  $|\langle \phi_{i,c} | r | \phi_{i,c} \rangle|^2$ . B=200A, h=80A.

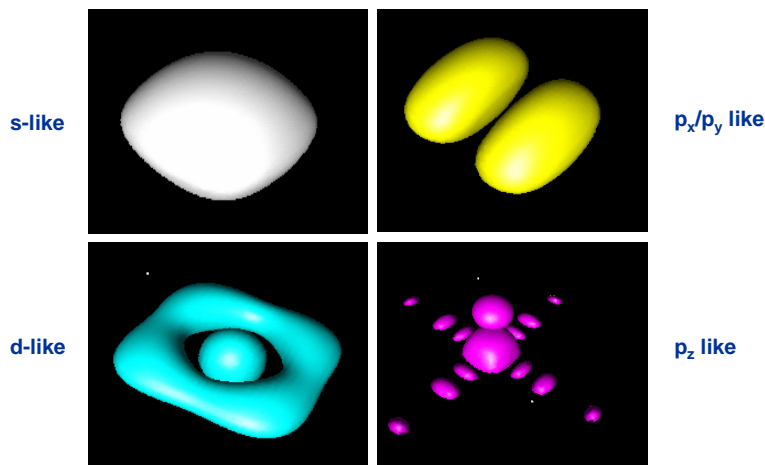
| Symmetry | state i    | x    | y    | z    |
|----------|------------|------|------|------|
| A1       | #2 (0.111) | 0    | 0    | 0.2  |
|          | #3 (0.123) | 0    | 0    | 57   |
|          | #4 (0.197) | 0    | 0    | 201  |
| A2       | #2 (0.106) | 0    | 0    | 28.5 |
|          | #3 (0.114) | 0    | 0    | 0    |
| B1,B2    | #2 (0.109) | 0    | 0    | 15   |
|          | #3 (0.138) | 0    | 0    | 42   |
|          | #4 (0.201) | 0    | 0    | 14   |
| A1-B1n   | #1 (0.062) | 446  | 446  | 0    |
|          | #2 (0.162) | 0.2  | 0.2  | 0    |
|          | #3 (0.218) | 0.4  | 0.4  | 0    |
| B1-A1n   | #2(0.049)  | 536  | 536  | 0    |
|          | #3(0.061)  | 659  | 659  | 0    |
|          | #4(0.135)  | 376  | 376  | 0    |
|          | #5(0.161)  | 10.2 | 10.2 | 0    |

## Theoretical Analysis



- Energy and wave functions computed using a stabilized transfer matrix technique by dividing the system into many slices along growth direction.
- Envelope function approximation with energy-dependent effective mass is used.
- Effective-mass Hamiltonian in k-space:  
$$[(k_x^2+k_y^2)/m_l(E)+\partial_z^2/m_l(E)-E]F(\mathbf{k}) + \sum_{\mathbf{k}'} [V(\mathbf{k},\mathbf{k}') + V_{\text{imp}}(\mathbf{k},\mathbf{k}')]F(\mathbf{k}') = 0$$
is solved via plane-wave expansion in each slice.
- 14-band k-p effects included perturbatively in optical matrix elements calculation
- Dopant effects incorporated as screened Coulomb potential
- The technique applies to quantum wells and quantum dots (or any 2D periodic nanostructures)

## Charge densities of low-lying states in lens-shaped QD



## Quantum well intrasubband photodetector (QWISP) for far infrared and terahertz radiation detection

[Ting et al., APPLIED PHYSICS LETTERS 91, 073510 (2007)]

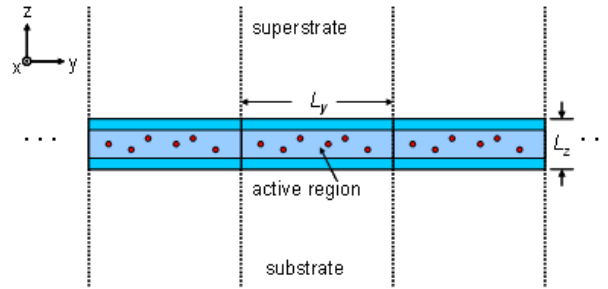


Fig. 4 A schematic illustration of the laterally repeating supercell geometry used in our simulations. Note that the supercell also repeats in the  $x$  direction with periodicity of  $L_x$  (not drawn). The example shown in this figure illustrates a quantum well containing several randomly placed dopant impurities.

## Quantum well intrasubband photodetector (QWISP) for far infrared and terahertz radiation detection

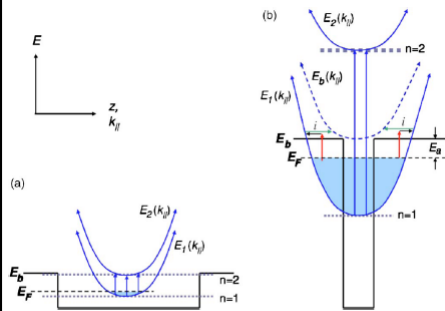


FIG. 1. (Color online) Schematic illustrations of the energy dispersions and energy band diagrams of (a) FIR/terahertz QWIP; (b) QWISP showing intersubband and impurity scattering assisted intrasubband optical absorption and carrier extraction mechanisms.

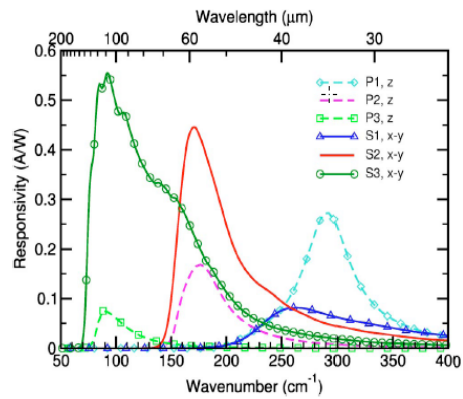


FIG. 3. (Color online) Low-temperature FIR/terahertz QWIP (set P) side-incidence responsivity and QWISP (set S) normal-incidence responsivity. Modeling parameters are listed in Table I.

## Quantum well intrasubband photodetector (QWISP) for far infrared and terahertz radiation detection

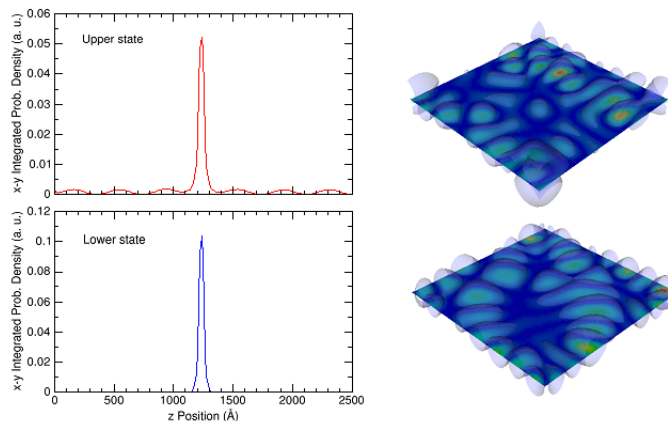


Fig. 5 The zero-bias wave functions of a pair of states involved in an intra-subband transition in a QWISP structure. The energy of the lower state is slightly below the Fermi level, and the energy of the upper state is just above the barrier band edge. The left panel shows the in-plane averaged probability densities as functions of  $z$  (growth direction). The random dopant distribution induced in-plane fluctuations of the wave functions are displayed in the right panel, which shows the translucent probability density isosurfaces, overlaid on top of probability density color contour plane located at the middle of the quantum well.

## Submonolayer QD infrared photodetector

[Ting et al., APPLIED PHYSICS LETTERS 94, 1 (2009)]

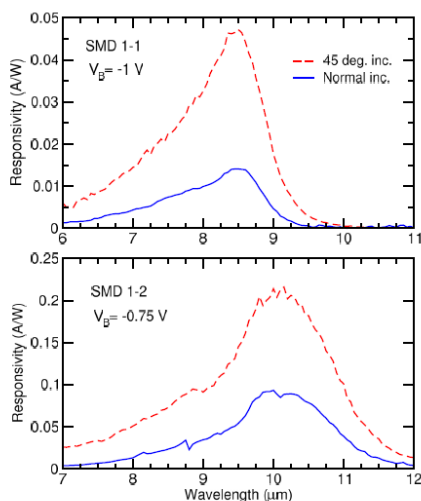


FIG. 2. (Color online) Measured normal and 45° incidence spectral responsivity curves for two SML QDIP samples.

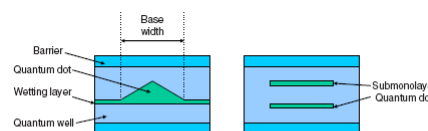


FIG. 1. (Color online) The left panel illustrates a conventional DWELL structure where a SK QD, consisting of pyramidal shape QD resting on a wetting layer, is embedded in a QW structure. The right panel shows two stacks of SML QDs embedded in a QW.

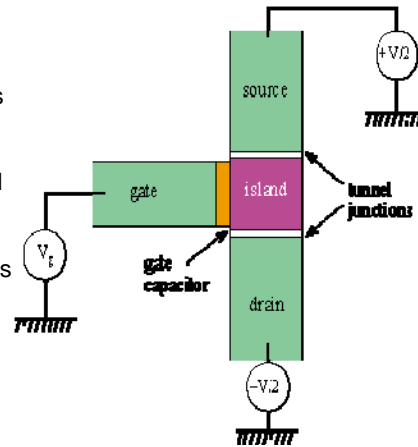


FIG. 4. (Color online) 80 K images taken with a 1024×1024 pixel SML QDIP FPA with  $f/2$  optics.

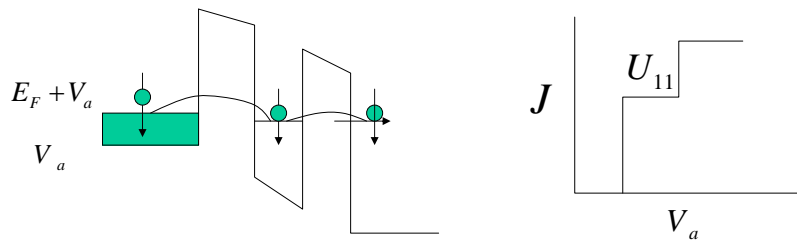
## Principle of the SET transistor



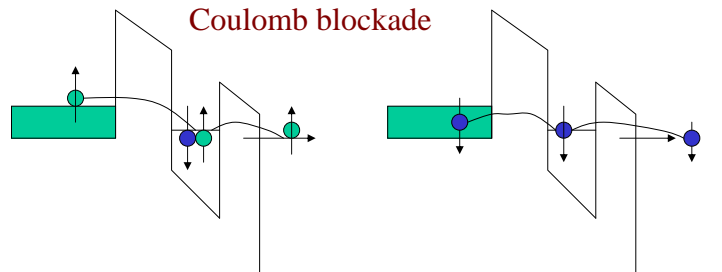
Like a MOSFET, the single-electron tunnelling (SET) transistor consists of a gate electrode that electrostatically influences electrons travelling between the source and drain electrodes. However, the electrons in the SET transistor need to cross two tunnel junctions that form an isolated conducting electrode called the island. Electrons passing through the island charge and discharge it, and the relative energies of systems containing 0 or 1 extra electrons depends on the gate voltage.



## Single-electron transistor (SET)



Coulomb blockade



## Tunneling current with Coulomb blockade



$$J = \frac{2e}{h} \sum_{\alpha} \int dw \Gamma [f_L(w - \mu_L) - f_R(w - \mu_R)] \text{Im} G_{d\alpha}(w)$$

$$G_{d\alpha}(w) = \left[ \frac{1 - \langle n_{d,\alpha} \rangle}{w - \varepsilon_0 + i\Gamma} + \frac{\langle n_{d,\alpha} \rangle}{w - \varepsilon_0 - U + i\Gamma} \right]$$

$$\langle n_{d,-\alpha} \rangle = \langle n_{d,\alpha} \rangle = \frac{1}{\pi} \int_{-\infty}^{\infty} dw \frac{f_L(w - \mu_L) + f_R(w - \mu_R)}{2} \text{Im} G_{d\alpha}(w)$$

## CdSe nanoparticles

(as fluorescent tags for biochemical analysis)



average diameter of particles (nm)

2.4    2.8    3.4    3.8    4.2

**Fluorescence from Nanoparticles  
of Various Diameters**

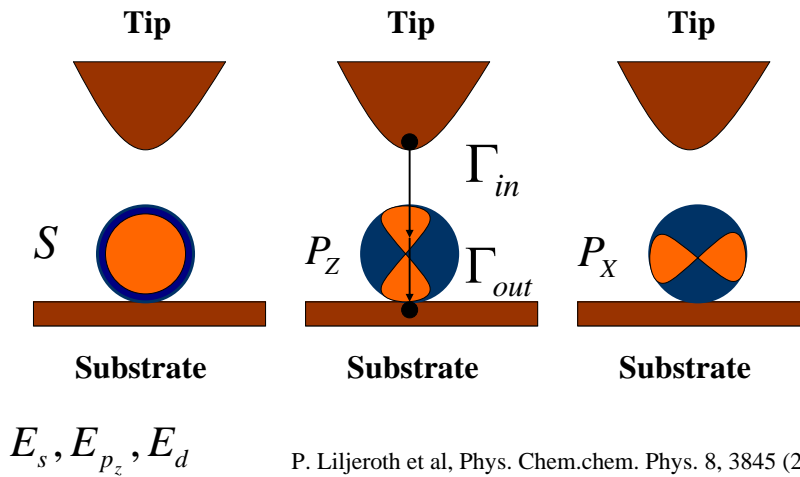
<http://unit.aist.go.jp/kyushu/mischel/eng/theme/theme2.htm>



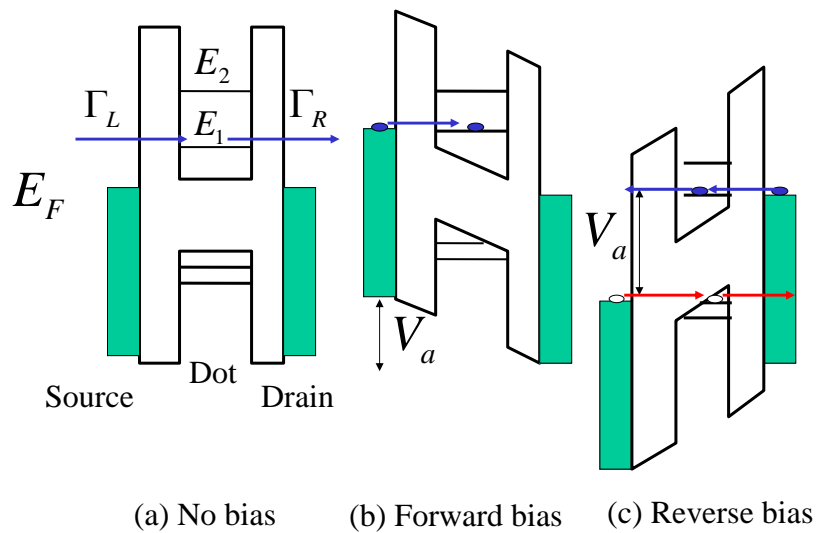
Tunneling current spectroscopy of a nanostructure junction involving multiple energy levels



David M T Kuo & Y. C. Chang



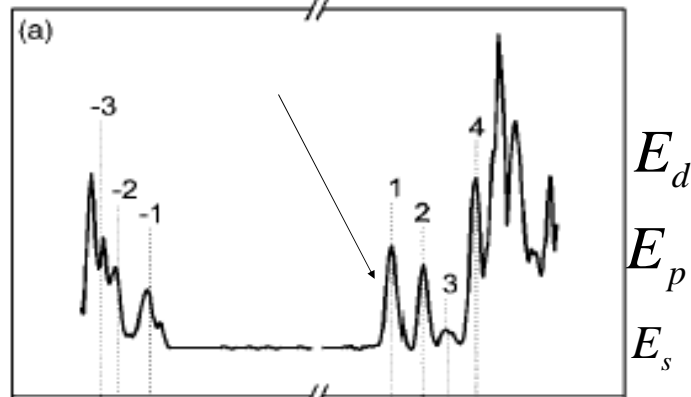
Energy diagram for STM-QD junction



## Experimental results



- Differential conductance oscillation



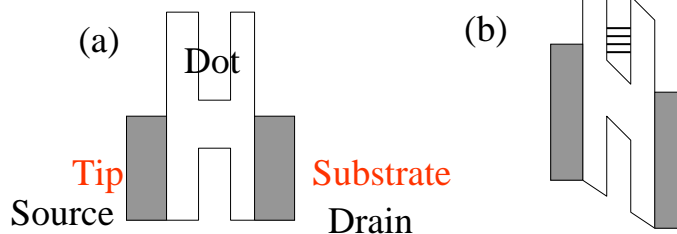
Identify current spectra.

L. Jdira et al, Phys. Rev. B 73, 115305 (2006)

## Multi-level Anderson Hamiltonian



$$\begin{aligned}
 H = & \sum_{k,\sigma,\beta} \epsilon_k a_{k,\sigma,\beta}^\dagger a_{k,\sigma,\beta} + \sum_{\ell,\sigma} E_\ell d_{\ell,\sigma}^\dagger d_{\ell,\sigma} \quad (1) \\
 & + \sum_{\ell,j,\sigma,\sigma'} U_{\ell,j} d_{\ell,\sigma}^\dagger d_{\ell,\sigma} d_{j,\sigma'}^\dagger d_{j,\sigma'} + \sum_{k,\sigma,\beta,\ell} V_{k,\beta,\ell} a_{k,\sigma,\beta}^\dagger d_{\ell,\sigma} \\
 & + \sum_{k,\sigma,\beta,\ell} V_{k,\beta,\ell}^* d_{\ell,\sigma}^\dagger a_{k,\sigma,\beta}
 \end{aligned}$$



H. Haug and A. P. Jauho, *Quantum Kinetics in Transport and Optics of Semiconductor* (Springer, Heidelberg, 1996).



## Tunneling Current formula

$$J = \frac{-2e}{\hbar} \sum_{\ell} \int \frac{d\epsilon}{2\pi} [f_{in}(\epsilon - \mu_{in}) - f_{out}(\epsilon - \mu_{out})] \quad (2)$$

$$\frac{\Gamma_{\ell,in}(\epsilon)\Gamma_{\ell,out}(\epsilon)}{\Gamma_{\ell,in}(\epsilon) + \Gamma_{\ell,out}(\epsilon)} \text{Im}G_{\ell,\sigma}^r(\epsilon),$$

**Is it possible to have a simple, efficient and reliable spectral function for a QD with arbitrary number of energy levels to determine the electronic structure of a single QD?**

H. Haug and A. P. Jauho, *Quantum Kinetics in Transport and Optics of Semiconductor* (Springer, Heidelberg, 1996).



## Mean-field approximation

$$\langle d_{i\sigma}^+ d_{i\sigma} d_{j\sigma}^+ d_{j\sigma} \rangle = \langle d_{i\sigma}^+ d_{i\sigma} \rangle \langle d_{j\sigma}^+ d_{j\sigma} \rangle \text{ for } i \neq j$$

- **One level case**

$$G_{1,\sigma}^r = (1 - N_{1,-\sigma}) \left( \frac{1}{\epsilon - E_1 + i \frac{\Gamma_1}{2}} \right) + N_{1,-\sigma} \left( \frac{1}{\epsilon - E_1 - U_1 + i \frac{\Gamma_1}{2}} \right)$$

The mean-field approximation is adequate for describing the Coulomb blockade, but not the Kondo effect

$N_{1\sigma} = \langle d^+ d \rangle$  is the average one-particle occupation number, which determines the probability of each resonant level.

How about the expression of retarded Green's functions for two-level, three-level and n-level case ?

H. Haug and A. P. Jauho, *Quantum Kinetics in Transport and Optics of Semiconductor* (Springer, Heidelberg, 1996).

## Two-level case (n=2)



$$G_{\ell,\sigma}^r(\varepsilon) = (1 - N_{\ell,-\sigma}) \sum_{m=1}^3 \frac{p_m}{\varepsilon - E_\ell - \Pi_m + i \frac{\Gamma_\ell}{2}}$$

$$+ N_{\ell,-\sigma} \sum_{m=1}^3 \frac{p_m}{\varepsilon - E_\ell - U_\ell - \Pi_m + i \frac{\Gamma_\ell}{2}}$$

Probability factors

Interlevel Coulomb factors

$$p_1 = a^j \equiv 1 - (N_{j,\sigma} + N_{j,-\sigma}) + \langle n_{j\sigma} n_{j-\sigma} \rangle \quad \Pi_1 = 0$$

$$p_2 = b^j \equiv N_{j,\sigma} + N_{j,-\sigma} - 2 \langle n_{j\sigma} n_{j-\sigma} \rangle \quad \Pi_2 = U_{\ell j}$$

$$p_3 = c^j \equiv \langle n_{j\sigma} n_{j-\sigma} \rangle \quad \Pi_3 = 2U_{\ell j}$$

$\ell \neq j$

## M-level case

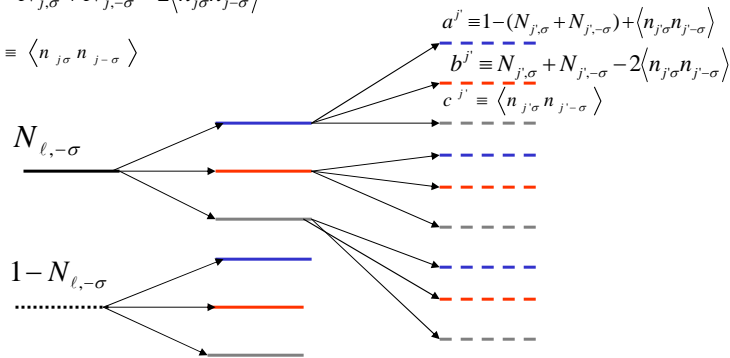


$$a^j \equiv 1 - (N_{j,\sigma} + N_{j,-\sigma}) + \langle n_{j\sigma} n_{j-\sigma} \rangle$$

$$b^j \equiv N_{j,\sigma} + N_{j,-\sigma} - 2 \langle n_{j\sigma} n_{j-\sigma} \rangle$$

$$c^j \equiv \langle n_{j\sigma} n_{j-\sigma} \rangle$$

$$\ell \neq j \neq j'$$



$$p_1 = a^j a^{j'}, p_2 = b^j a^{j'}, p_3 = a^j b^{j'}, p_4 = c^j a^{j'}, p_5 = c^j a^{j'}, \dots$$

$$\Pi_1 = 0, \Pi_2 = U_{\ell j}, \Pi_3 = U_{\ell j'}, \Pi_4 = 2U_{\ell j}, \Pi_5 = 2U_{\ell j'},$$

We define the  $(2+n)$ - and  $(1+n)$ -particle ( $n \leq 2M-2$ ) Green's functions as

$$G_{\ell, j_1, \dots, j_n}^r = \langle n_{\ell, -\sigma} n_{j_1} \dots n_{j_n} d_{\ell, \sigma} d_{\ell, \sigma}^\dagger \rangle$$

and

$$G_{\ell, j_1, \dots, j_n}^r = \langle n_{j_1} \dots n_{j_n} d_{\ell, \sigma} d_{\ell, \sigma}^\dagger \rangle,$$

where  $j_1, \dots, j_n$  label any  $n$  states out of the  $2(M-1)$  states (excluding spin up and down states in level  $\ell$ ). Here  $j_\alpha$  is a composite index for energy level and spin. In general, the  $(2+n)$ - and  $(1+n)$ -particle ( $n \leq 2M-2$ ) Green's functions satisfy

$$\begin{aligned} (\mu_l - U_\ell - \sum_{\alpha=1}^n U_{\ell, j_\alpha}) G_{\ell, j_1, \dots, j_n}^r &= N_{\ell, -\sigma} \langle n_{j_1} \dots n_{j_n} \rangle \\ &+ \sum_{j'}^n U_{\ell, j'} G_{\ell, j_1, \dots, j_n, j'}^r, \end{aligned} \quad (32)$$

and

$$\begin{aligned} (\mu_l - \sum_{\alpha=1}^n U_{\ell, j_\alpha}) G_{\ell, j_1, \dots, j_n}^r &= \langle n_{j_1} \dots n_{j_n} \rangle + U_\ell G_{\ell, j_1, \dots, j_n}^r \\ &+ \sum_{j'}^n U_{\ell, j'} G_{\ell, j_1, \dots, j_n, j'}^r, \end{aligned} \quad (33)$$

where the "double prime" on the summation indicates that  $j'$  is not among the  $n$  states and not in the level  $\ell$ . In the  $n$ -particle state labeled by  $j_1, \dots, j_n$  if there are  $m$  levels occupied with two particles and the rest singly occupied, we can label them by  $j_1, j_1, \dots, j_m, j_m, j_{2m+1}, \dots, j_n$ . We then assume that the  $n$ -particle correlation function can



$$G_{2M}^r = \frac{N_{\ell, -\sigma} \prod_j' c_j}{(\mu_l - U_\ell - 2 \sum_j' U_{\ell, j})}, \quad (4)$$

where  $\mu_\ell \equiv \epsilon - E_\ell + i(\Gamma_{\ell, L} + \Gamma_{\ell, R})/2$  and  $U_{\ell, j}$  denotes the on-site Coulomb energy in level  $\ell$ .  $N_{j, -\sigma} = \langle n_{j, -\sigma} \rangle$  and  $c_j = \langle n_{j, -\sigma} n_{j, \sigma} \rangle$  denote the one-particle and two-particle average occupancy in the level  $j$ . Here and henceforth,  $\prod_j'$  means taking the product of terms labeled by  $j$  with  $j = 1, \dots, M$ , excluding  $\ell$ .

$$\begin{aligned} G_{\ell, \ell}^r(\epsilon) &= \prod_j' (\hat{a}_j + \hat{b}_j + c_j) G_{2M}^r / \prod_j' c_j \\ &\equiv N_{\ell, -\sigma} \sum_{m=1}^{3^M-1} \frac{P_m}{\mu_\ell - U_\ell - \Pi_m} \end{aligned} \quad (5)$$

and

$$\begin{aligned} G_{\ell, \sigma}^r(\epsilon) &= (\hat{b}_\ell N_{\ell, -\sigma}^{-1} + 1) G_{\ell, \ell}^r(\epsilon) \\ &= (1 - N_{\ell, -\sigma}) \sum_{m=1}^{3^M-1} \frac{P_m}{\mu_\ell - \Pi_m} + N_{\ell, -\sigma} \sum_{m=1}^{3^M-1} \frac{P_m}{\mu_\ell - U_\ell - \Pi_m}, \end{aligned} \quad (6)$$

where  $\hat{b}_\ell$ ,  $\hat{a}_j$ , and  $\hat{b}_j$  are operators that put a factor  $b_\ell = 1 - N_{\ell, -\sigma}$ ,  $a_j \equiv 1 - (N_{j, \sigma} + N_{j, -\sigma}) + c_j$ , and  $b_j \equiv N_{j, \sigma} + N_{j, -\sigma} - 2c_j$  in the numerator and increases the value of the denominator by  $U_\ell$ ,  $2U_{\ell, j}$  and  $U_{\ell, j}$ , respectively when acting on a fractional function. For ex-



Differential conductance  $dJ/dV$  is determined by:



$$G_{\ell,\sigma}^r(\epsilon) = (1 - N_{\ell,-\sigma}) \sum_{m=1}^{3^n-1} \frac{p_m}{\epsilon - E_\ell - \Pi_m + i\frac{\Gamma_\ell}{2}} + N_{\ell,-\sigma} \sum_{m=1}^{3^n-1} \frac{p_m}{\epsilon - E_\ell - U_\ell - \Pi_m + i\frac{\Gamma_\ell}{2}},$$

$\Pi_m$  = sum of Coulomb interactions in configuration  $m$ .

$p_m$  = probability of finding the system in configuration  $m$ .

Level  $j$  ( $j \neq \ell$ ) can be occupied by 0, 1 or 2 particles.

For a two-level ( $n = 2$ ) system ( $\ell \neq j$ ), we have three configurations:

$$p_1 = a^j \equiv 1 - (N_{j,\sigma} + N_{j,-\sigma}) + \langle n_{j,\sigma} n_{j,-\sigma} \rangle$$

(the probability with no particle in level  $j$ ).

$$p_2 = b^j \equiv N_{j,\sigma} + N_{j,-\sigma} - 2\langle n_{j,\sigma} n_{j,-\sigma} \rangle$$

(the probability with one particle in level  $j$ )

$$p_3 = c^j \equiv \langle n_{j,\sigma} n_{j,-\sigma} \rangle$$

(the probability with two particles in level  $j$ )

$$\Pi_1 = 0, \Pi_2 = U_{\ell j} \text{ and } \Pi_3 = 2U_{\ell j}.$$

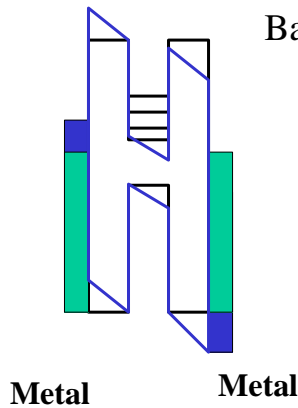
$U_{\ell j}$  = Coulomb interaction between level  $\ell$  and level  $j$ .

$$p_m = \prod_j (\hat{a}_j + \hat{b}_j + c_j)$$

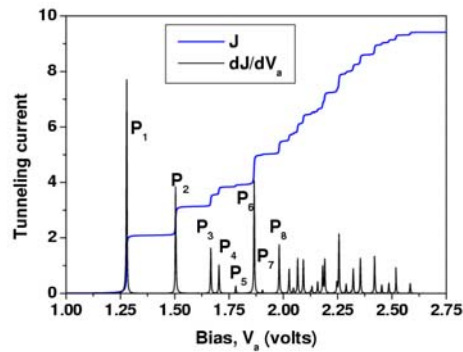
## 2. Three-level system at symmetry case



Based on three-level case



At zero temperature



- d-orbital

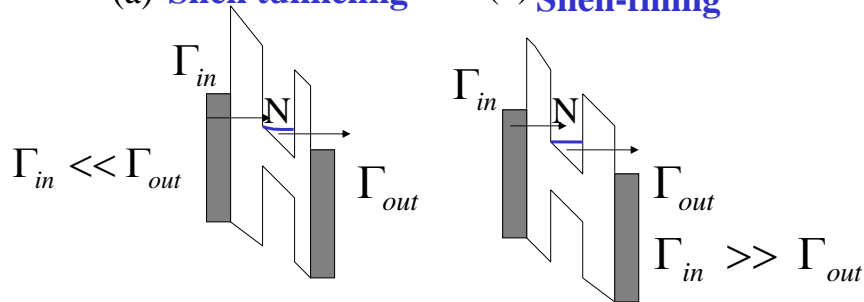
## Asymmetry tunneling rates



- Tunneling rates  $\Gamma_{in}$  and  $\Gamma_{out}$

(a) **Shell tunneling**

(b) **Shell-filling**

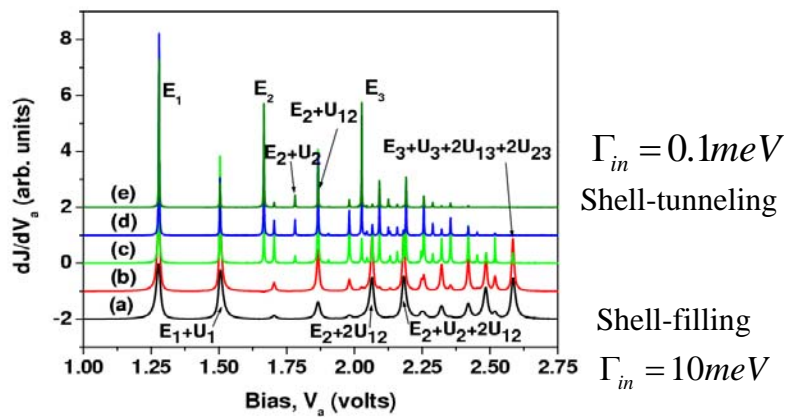


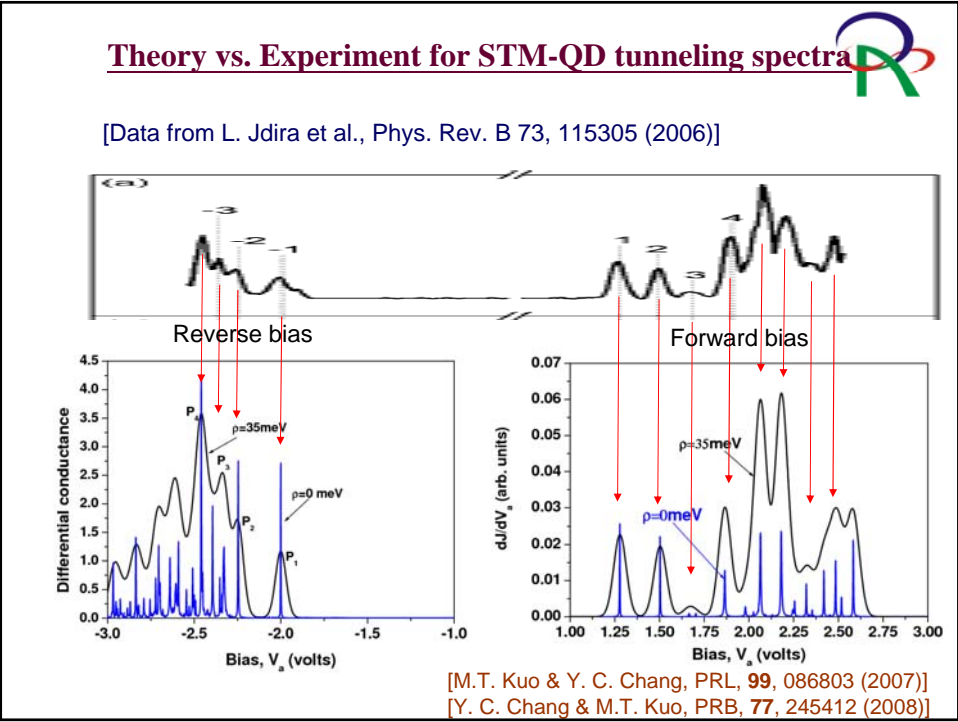
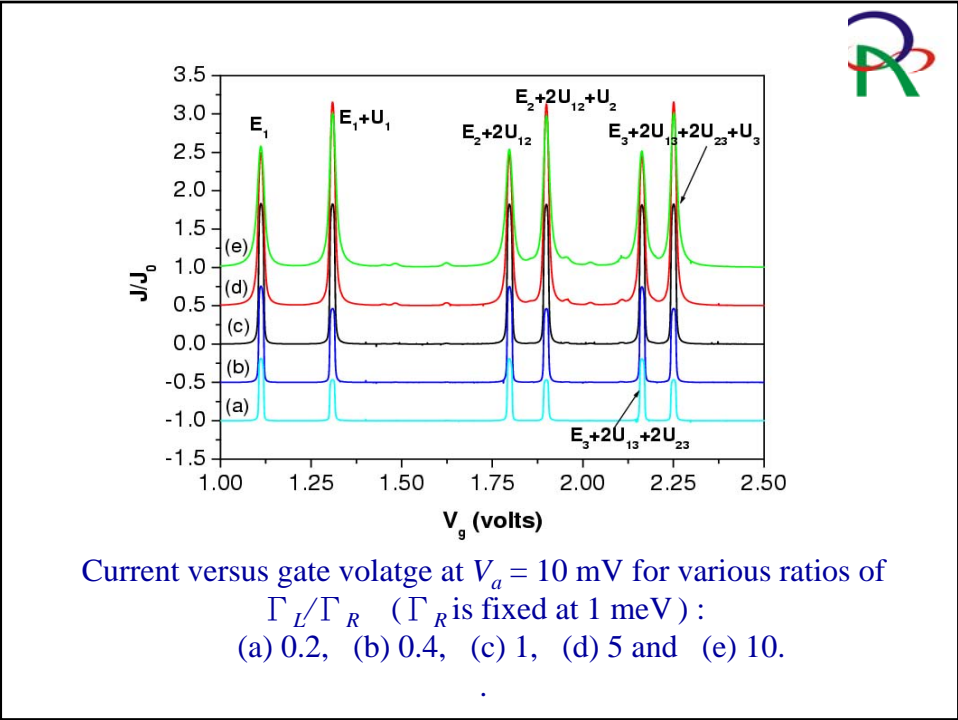
The characteristics of tunneling current of junctions can be classified into either the shell-tunneling case (no charge accumulation) or the shell-filling case (with charge accumulation)

## Effect of asymmetric tunneling



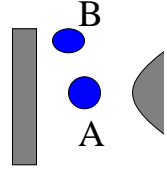
$$\Gamma_{out} = 1meV$$



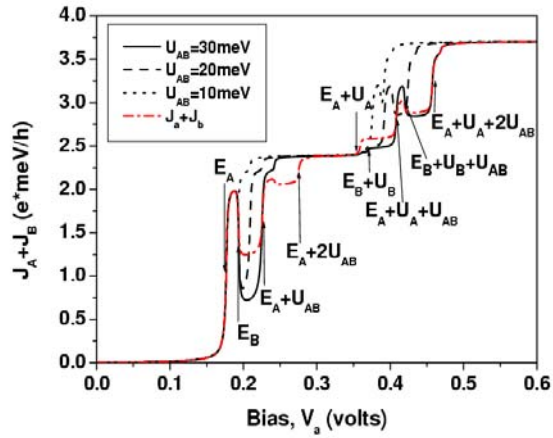




## Proximity effect for coupled QDs



$$\begin{aligned}\Gamma_{in,A} &= 1\text{meV} \\ \Gamma_{out,A} &= 1\text{meV} \\ \Gamma_{in,B} &= 1\text{meV} \\ \Gamma_{out,B} &= 0.1\text{meV}\end{aligned}$$



## Proximity effect for tunneling current in nanojunction with coupled QDs

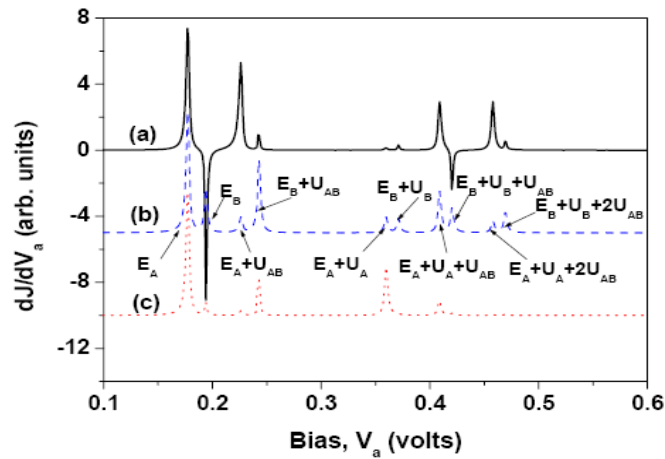
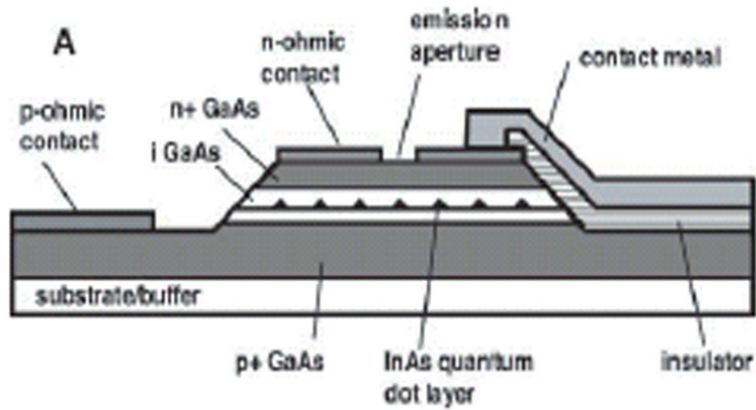


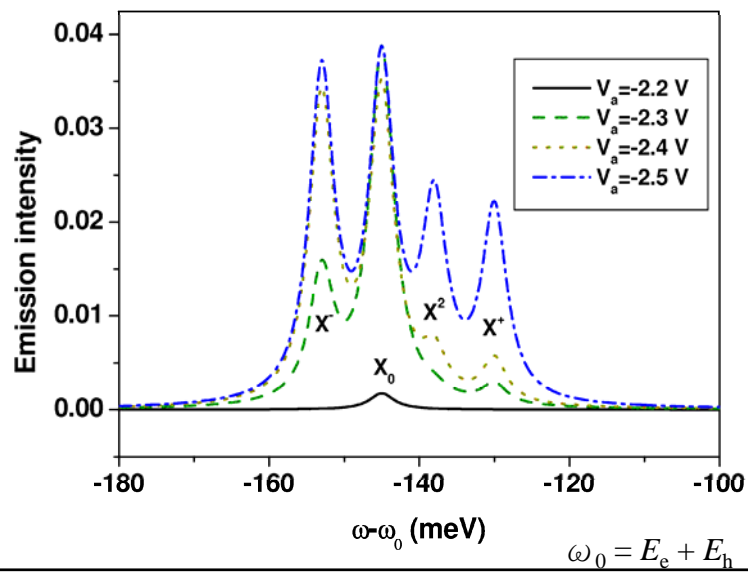
FIG. 4: Differential conductance as a function of applied bias for various tunneling rate ratios of  $\Gamma_{L,B}/\Gamma_{R,B}$ .  $U_{AB} = 30$  meV. (a)  $\Gamma_{L,B} = 1$  meV and  $\Gamma_{R,B} = 0.1$  meV. (b)  $\Gamma_{L,B} = 1$  meV and  $\Gamma_{R,B} = 1$  meV. (c)  $\Gamma_{L,B} = 0.1$  meV and  $\Gamma_{R,B} = 1$  meV.

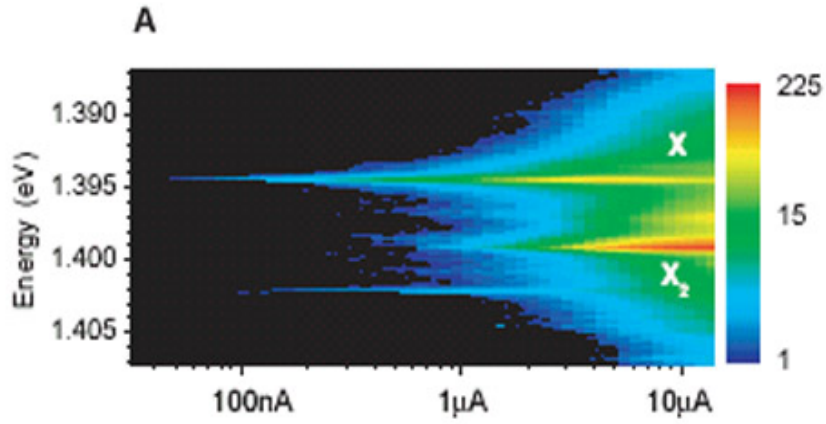
## Single-Photon generator



Z. Yuan et al., Science, 295, 102 (2002)

## Emission spectra of exciton complexes for various applied biases in a reverse bias setup





Z. Yuan et al., Science, 295, 102 (2002)



### Including optical field

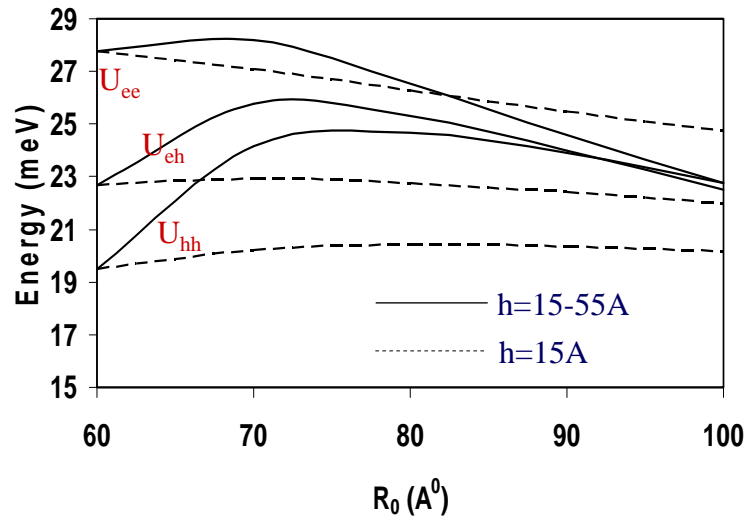
$$\Gamma_1 N_1 = -\text{Imp} - \int \frac{d\varepsilon}{\pi} \Gamma_1^{L/R} f_{L/R}(\varepsilon) \text{Im} G_{11}(\varepsilon + \frac{\omega}{2})$$

$$\Gamma_2 N_2 = \text{Imp} - \int \frac{d\varepsilon}{\pi} \Gamma_2^{L/R} f_{L/R}(\varepsilon) \text{Im} G_{22}(\varepsilon - \frac{\omega}{2})$$

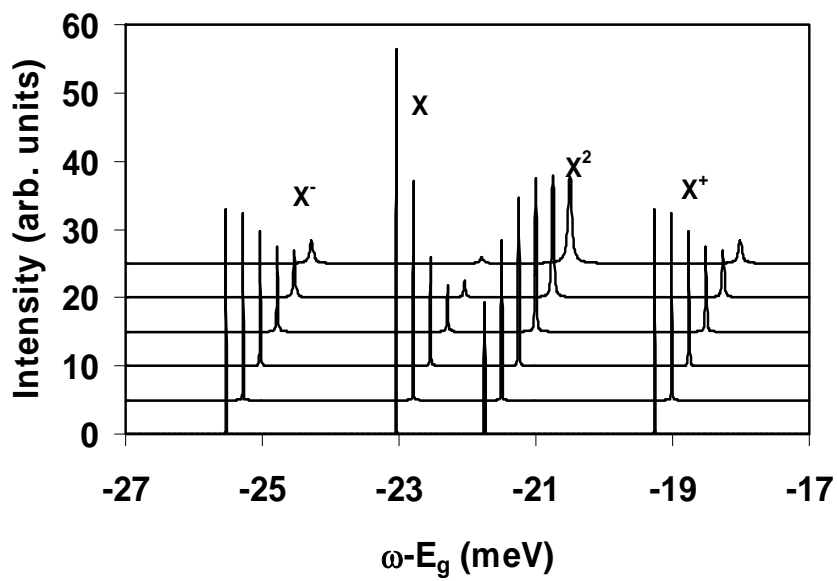
$$\text{Imp} \approx 2\lambda^2 \int \frac{d\varepsilon}{2\pi} \frac{\Gamma_1^L f_L(\varepsilon) + \Gamma_1^R f_R(\varepsilon)}{\Gamma_1} \left\{ \text{Im} \left( \frac{1 - N_1}{\varepsilon - E_1 + i \frac{\Gamma_1}{2}} \right) \text{Im} \left( \frac{1}{\varepsilon - E_2 + \omega - i \frac{\Gamma_2}{2}} \right) \right.$$

$$\left. + \text{Im} \left( \frac{N_1}{\varepsilon - E_1 - U_{11} + i \frac{\Gamma_1}{2}} \right) \text{Im} \left( \frac{1}{\varepsilon - E_2 - U_{12} + \omega - i \frac{\Gamma_2}{2}} \right) \right\}$$

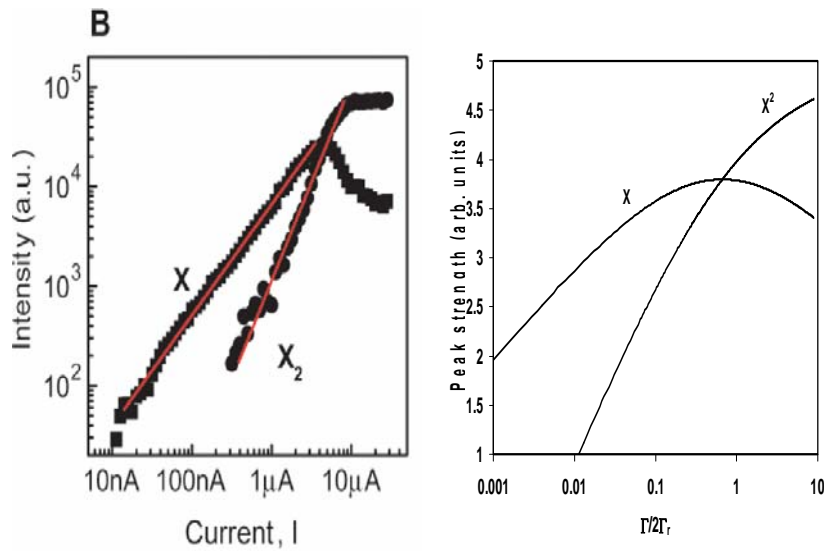
### Coulomb interactions vs. QD size



### Spontaneous emission spectrum



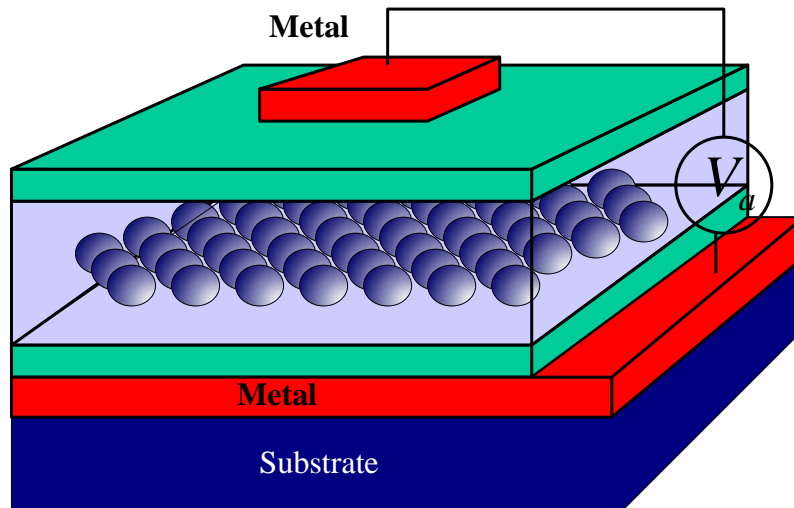
## Intensity vs current



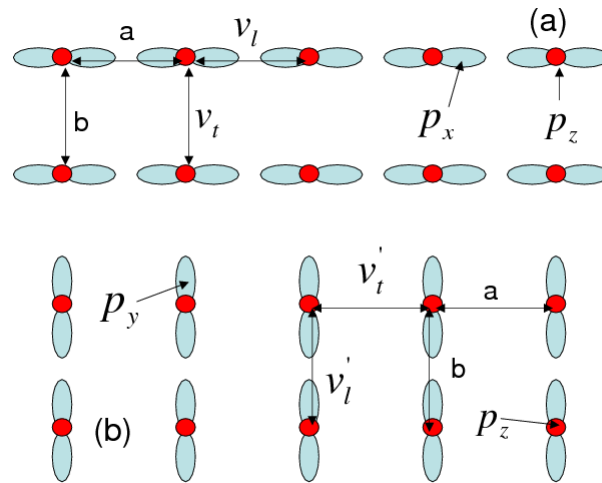
[Z. Yuan et al., Science, 295, 102 (2002)]

[Phys. Rev. B **69**, R41306 (2004)]

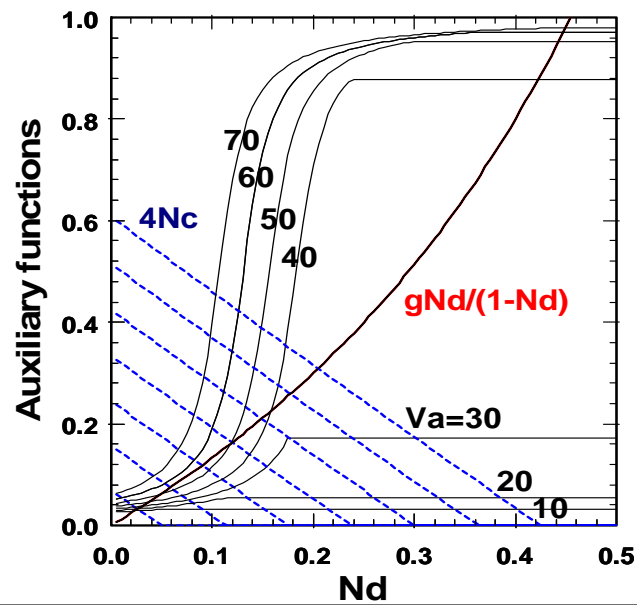
## Quantum dot array(QDA) memory



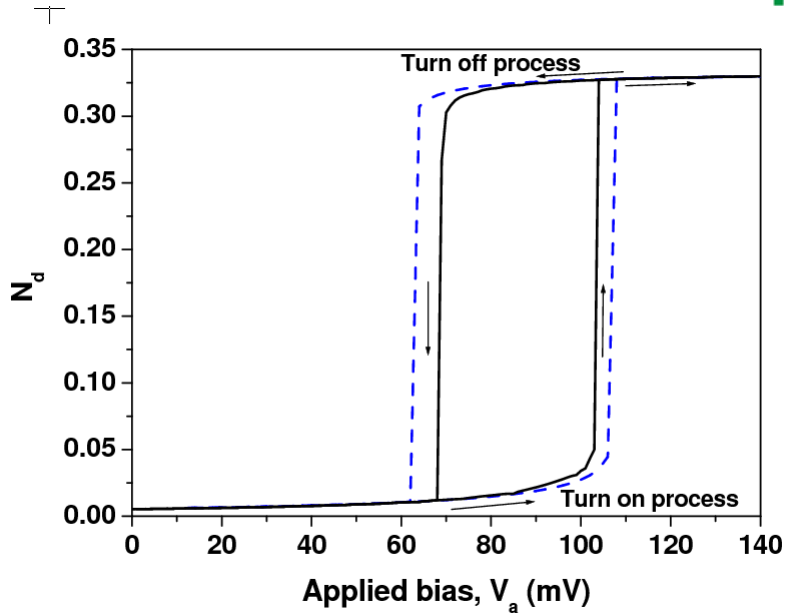
## Coupling of degenerate p orbitals in QDA



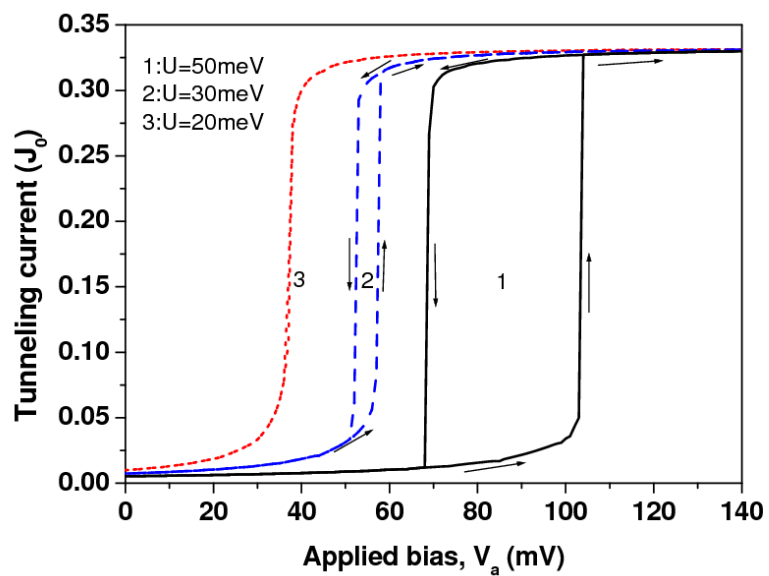
## Bistable solution to occupation number



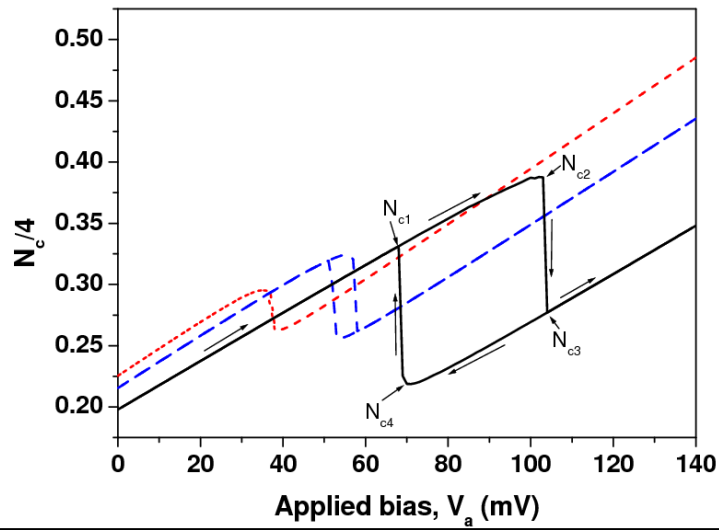
## Bistable behavior of occupation number



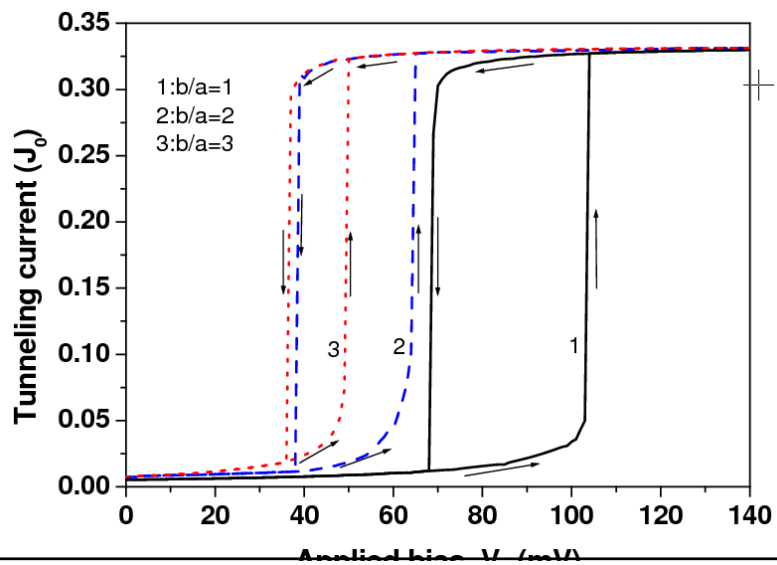
## Bistable tunneling current through QDA



## Bistable behavior of occupation number in the QDA bands

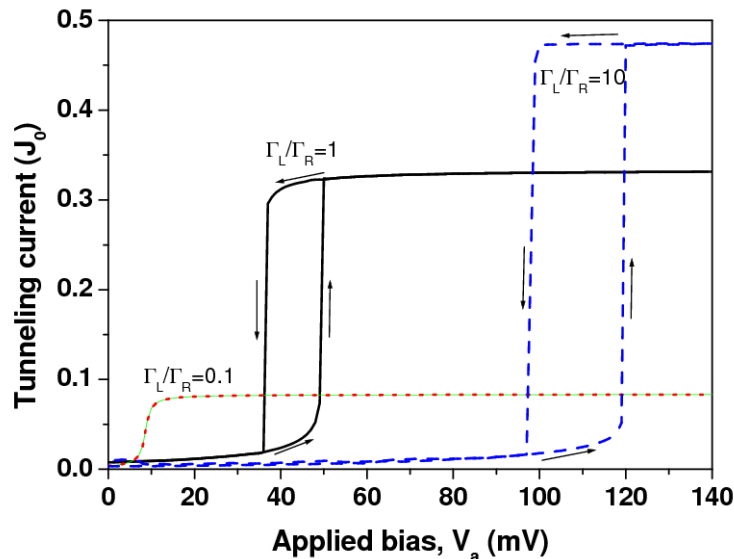


## Bistable tunneling current from 2D to 1D





## Effects of in/out ratio of tunneling rate in 1D QDA



## Summary



- **A simple, efficient and reliable closed-form formula is derived to reveal the tunneling current spectra of a complicated nanostructure junction**
- **Negative differential conductance arises from interdot Coulomb interactions**
- **The formula can also be used to study the transport properties of coupled nanostructures and QDA**
- **Memory effects of QDA with degenerate p orbitals are demonstrated**