

CHAPTER III

THE SUPERCONDUCTING PROXIMITY EFFECT

McMillan (McMillan, 1968) studied a theoretical model of the proximity effect between superposed normal (N) and superconducting (S) metal films. He imagined a potential barrier between the films and treated an electron transmission through the barrier as a tunneling process. In the McMillan model :

(a) The tunneling Hamiltonian is used to describe the penetration of electron through the barrier. This restricts the transmission probability of the barrier to be much less than one.

(b) Both N and S films are thin compared with the characteristic superconducting length so that the properties of each film are uniform across its thickness.

(c) Tunneling matrix elements $V_{kk'}$ is taken to be of equal magnitude between energy states in S and N.

Mohabir and Nagi (Mohabir and Nagi, 1979) extended the McMillan tunneling model (MTM) of the superconducting proximity effect by treating the tunneling Hamiltonian to all order of self-consistent perturbation theory. They considered the normal metal in MTM as weak superconductor and treated their system as superconductor-superconductor sandwiches. In this review, we will consider the model of Mohabir and Nagi for S_1 - S_2 sandwiches.

The Hamiltonian for the sandwich is the sum of the Hamiltonians for S_1 and S_2 slabs and the tunneling Hamiltonian.

$$H = H_{S_1} + H_{S_2} + H_T \quad (3.1)$$

where $H_{S_1}(H_{S_2})$ is the Hamiltonian of the 1st (2nd) superconducting metal and H_T is the tunneling Hamiltonian.

$$H_{S_1} = \sum_{k\sigma} \epsilon_{k1} a_{k\sigma}^+ a_{k\sigma} - \Delta_1 \sum_k (a_{k\uparrow}^+ a_{-k\downarrow}^+ + \text{h.c.}) \quad (3.2.1)$$

$$H_{S_2} = \sum_{k\sigma} \epsilon_{k2} d_{k\sigma}^+ d_{k\sigma} - \Delta_2 \sum_k (d_{k\uparrow}^+ d_{-k\downarrow}^+ + \text{h.c.}) \quad (3.2.2)$$

$$H_T = \sum_{k k' \sigma} [V_{kk'} a_{k\sigma}^+ d_{k'\sigma} + \text{h.c.}] \quad (3.2.3)$$

where $a_{k\sigma}^+$ ($d_{k\sigma}^+$) is the creation operator for electron in the $S_1(S_2)$ side, σ is the spin index, and $V_{kk'}$ is the tunneling matrix element, we take $V_{kk'}$ to be independent of k and k' . ϵ_{k1} (ϵ_{k2}) is the band energy of the conduction electron in the $S_1(S_2)$ side measured from the Fermi energy. Δ_1 (Δ_2) is the BCS superconducting order parameter in the $S_1(S_2)$ side.

In order to calculate Δ , we introduce the finite-temperature Green's function.

$$G(k, \omega_n) = - \langle T_\tau \psi_k(\tau) \psi_k^+(0) \rangle \quad (3.3)$$

where $\Psi_{\mathbf{k}}^{\dagger} = (a_{\mathbf{k}\uparrow}^{\dagger} \ a_{-\mathbf{k}\downarrow})$ for S_1 side and T_{τ} is the time ordering operator for the imaginary time $\tau = it$.

We find a single particle Green's function for pure S_1 superconductor $G_1^{\circ}(\mathbf{k}, \omega_n)$ as

$$G_1^{\circ}(\mathbf{k}, \omega_n) = [i\omega_n - \epsilon_{\mathbf{k}}\tau_3 + \Delta_1\tau_1]^{-1} \quad (3.4)$$

where τ_i ($i=1,2,3$) are Pauli matrices.

Treating the electron-phonon interaction and the tunneling Hamiltonian in second order self-consistent perturbation, we find for the Nambu matrix-self-energy of the S_1 side as

$$\Sigma_1(i\omega_n) = V^2 \tau_3 \sum_{\mathbf{k}} G_1(\mathbf{k}, \omega_n) \tau_3 \quad (3.5)$$

where V is the tunneling magnitude.

The self-energy for S_2 side $\Sigma_2(i\omega_n)$ can be obtained from Eq.(3.5) by interchanging the labels 1 and 2.

We make the ansatz

$$G(\mathbf{k}, \omega_n) = [i\tilde{\omega}_n - \epsilon_{\mathbf{k}}\tau_3 + \tilde{\Delta}\tau_1]^{-1} \quad (3.6.1)$$

here $G(\mathbf{k}, \omega_n)$ is the renormalized Green's function which includes the proximity effect. $\tilde{\omega}_n$ and $\tilde{\Delta}$ are renormalized frequency and renormalized order parameter, respectively.

We find the usual form for matrix Green's function

$$G(\mathbf{k}, \omega_n) = -[i\tilde{\omega}_n + \varepsilon_k \tau_3 - \tilde{\Delta} \tau_1] / [\tilde{\omega}_n^2 + \varepsilon_k^2 + \tilde{\Delta}^2] \quad (3.6.2)$$

Using the Dyson's equation, the self-energy is related to the Green's functions, we have

$$\Sigma_1(i\omega_n) = i(\omega_{1n} - \tilde{\omega}_{1n}) + (\Delta_1 - \tilde{\Delta}_1) \tau_1 \quad (3.7)$$

Substitution of Eq.(3.6.2) into Eq.(3.5) and performing the sum over states, we find

$$\Sigma_1(i\omega_n) = -V^2 A d_2 N_2(0) \pi [(i\tilde{\omega}_{2n} + \tilde{\Delta}_2 \tau_1) / \sqrt{\tilde{\omega}_{2n}^2 + \tilde{\Delta}_2^2}] \quad (3.8)$$

where A, d_2 , and $N_2(0)$ are the area, thickness and bulk density of states (per unit volume) of the second superconductor.

By comparing Eq.(3.7) and (3.8), we obtain the following equations.

$$\tilde{\omega}_{1n} = \omega_{1n} + \Gamma_1 \tilde{\omega}_{2n} [\sqrt{\tilde{\omega}_{2n}^2 + \tilde{\Delta}_2^2}]^{-1} \quad (3.9.1)$$

$$\tilde{\Delta}_1 = \Delta_1 + \Gamma_1 \tilde{\Delta}_2 [\sqrt{\tilde{\omega}_{2n}^2 + \tilde{\Delta}_2^2}]^{-1} \quad (3.9.2)$$

In these equations $\Gamma_1 = V^2 A d_2 N_2(0) \pi$.

Similarly for the second superconductor, we can also obtain the equations for $\tilde{\omega}_{2n}$ and $\tilde{\Delta}_2$ by changing the labels 1 and 2, and define $\Gamma_2 = V^2 A d_1 N_1(0) \pi$.

The solution of the self-energy can be solved by defining a new parameter.

$$U_n = \tilde{\omega}_n / \tilde{\Delta} \quad (3.10)$$

Then, Eqs.(3.9.1) and (3.9.2) can be combined as

$$U_1 = [1/\Delta_1] [\omega_{1n} + \Gamma_1 (U_2 - U_1) [\sqrt{1+U_2^2}]^{-1}] \quad (3.11.1)$$

Similarly, we can show that

$$U_2 = [1/\Delta_2] [\omega_{2n} + \Gamma_2 (U_1 - U_2) [\sqrt{1+U_1^2}]^{-1}] \quad (3.11.2)$$

The self-consistent equation for the order parameter is

$$\Delta_1 = \frac{\omega_{D1}/2\pi T}{2\pi \lambda_1 T} \sum_{n \geq 0} \tilde{\Delta}_1 [\sqrt{\tilde{\omega}_{1n}^2 + \tilde{\Delta}_1^2}]^{-1} \quad (3.12)$$

where $\lambda_1 = g_1 A N_1(0) d_1$.

Substitution of Eq.(3.10) into (3.12) and taking the limit that near T_c , Δ_1 and Δ_2 are small and $U_1, U_2 \gg 1$, we get

$$\Delta_1 = \frac{\omega_{D1}/2\pi T}{2\pi\lambda_1 T} \sum_{n \geq 0} [U_1]^{-1} \quad (3.13)$$

Further, Eqs.(3.11.1) and (3.11.2) give

$$U_1 = [1/\Delta_1] \{ \omega_n (\Gamma_1 + \Gamma_2 + \omega_n) [\Gamma_1 (\Delta_2/\Delta_1) + \Gamma_2 + \omega_n]^{-1} \} \quad (3.14)$$

Eqs.(3.13) and (3.14) give

$$\Delta_1 = \lambda_1 \{ \Delta_1 f_1(T) - \Gamma_1 (\Gamma_1 + \Gamma_2)^{-1} [\Delta_1 - \Delta_2] K_1(\rho) \} \quad (3.15.1)$$

where

$$f_1(T) = \frac{\omega_{D1}/2\pi T}{2\pi T} \sum_{n \geq 0} [\omega_n]^{-1} = \ln(2\gamma \omega_{D1}/\pi T) \quad (3.15.2)$$

and

$$\begin{aligned} K_1(\rho) &= \frac{\omega_{D1}/2\pi T}{2\pi T} \sum_{n \geq 0} \{ (\omega_n)^{-1} - (\omega_n + \Gamma_1 + \Gamma_2)^{-1} \} \\ &= \psi(-1/2 + \rho/2) - \psi(-1/2) - \ln(1 - \Gamma/\omega_{D1}) \end{aligned} \quad (3.15.3)$$

where $\Gamma = \Gamma_{N^+} \Gamma_S$, $\gamma = 1.781$, $\rho = \Gamma/\pi T$ and ψ is the digamma function (Davis,1965).

The equation for Δ_2 and U_2 are obtained from the above equations by interchanging the labels 1 and 2.

From Eq.(3.15.1) we set up an equation for Δ_2 , Eliminating Δ_1 / Δ_2 from these two equations, we finally get

$$\lambda_1 \lambda_2 f_1(T) - f_1(T) \{ \lambda_1 + \lambda_2 - \lambda_1 \lambda_2 \chi + \lambda_1 \lambda_2 [\gamma_1 K_1(\rho) + \gamma_2 K_2(\rho)] \} \\ + [1 - \chi \lambda_2] [1 + \lambda_1 \gamma_1 K_1(\rho)] + \lambda_2 \gamma_2 K_2(\rho) = 0 \quad (3.16)$$

where $\chi = \ln(\omega_{D2}/\omega_{D1})$, $\gamma_1 = \Gamma_1/\Gamma$, $\gamma_2 = \Gamma_2/\Gamma$.

For a superconductor-normal metal sandwich, we can make the assumption that $\lambda_2 = 0$. Eq.(3.16) gives the transition temperature of an S-N sandwich as

$$\ln(T_{cs}^*/T_c) = \Gamma_N/\Gamma [\psi(-1/2 + \Gamma/2\pi T_c) - \psi(-1/2)] \quad (3.17)$$

where T_{cs}^* is the transition temperature of the single S slab.

Eq.(3.17) is the McMillan's transition temperature of S-N sandwich (McMillan,1968).

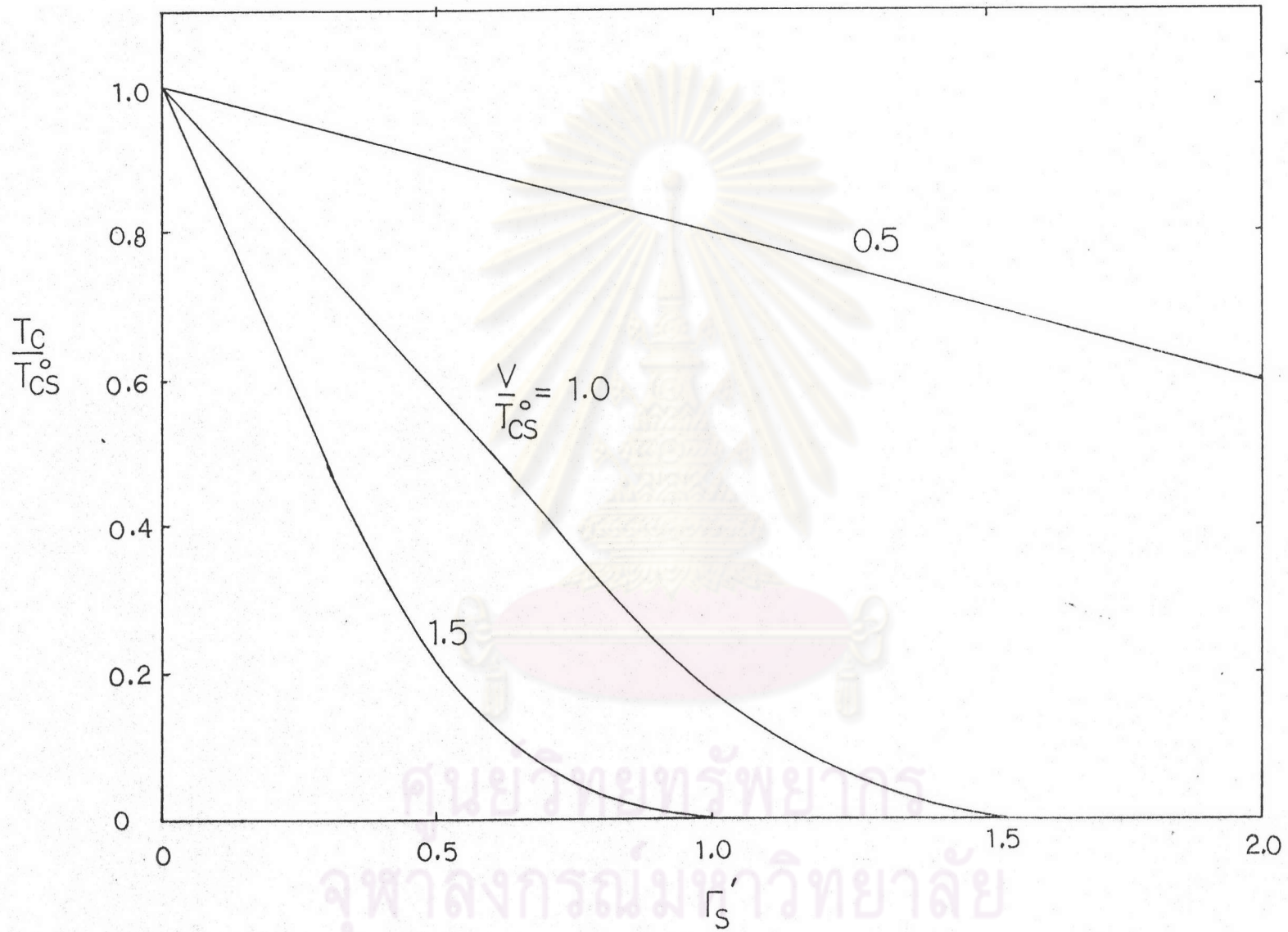


Fig. 3.1 Critical temperature T_c of S-N sandwich vs. Γ'_s ($\Gamma'_s = \pi A d_N N_N(0) T_{c0}$) for various values of V/T_{c0} .
The parameter is $\Gamma_N = 0.2$ (Mohabir and Nagi, 1979).

In Fig. 3.1, we have introduced dimensionless parameter Γ'_s ($\Gamma'_s = \pi A d_N N_N(0) T_{cs}^*$). When $\Gamma'_s = 0$, the ratio of T_c/T_{cs}^* is equal 1 for various values of V/T_{cs}^* . For fixed values of V/T_{cs}^* , the ratio of T_c/T_{cs}^* decreases when V/T_{cs}^* increases.

The transition temperature of an S_1 - S_2 sandwich is given by solving the quadratic Eq.(3.16), we finally get

$$\ln(T_{c1}^*/T_c) = (1/2) [A + H - \sqrt{A^2 + H^2 + 2AG}] \quad (3.18)$$

where

$$A = (\lambda_2)^{-1} - (\lambda_1)^{-1} - \chi$$

$$H(\rho_{c1}) = \gamma_1 K_1(\rho_{c1}) + \gamma_2 K_2(\rho_{c1})$$

$$G(\rho_{c1}) = \gamma_2 K_2(\rho_{c1}) - \gamma_1 K_1(\rho_{c1})$$

Eq.(3.18) is the transition temperature of an S_1 - S_2 sandwich as shown by Mohabir and Nagi (Mohabir and Nagi,1979).

In Figs. (3.2) and (3.3) , we show calculated values of the transition temperature of an S_1 - S_2 sandwich.

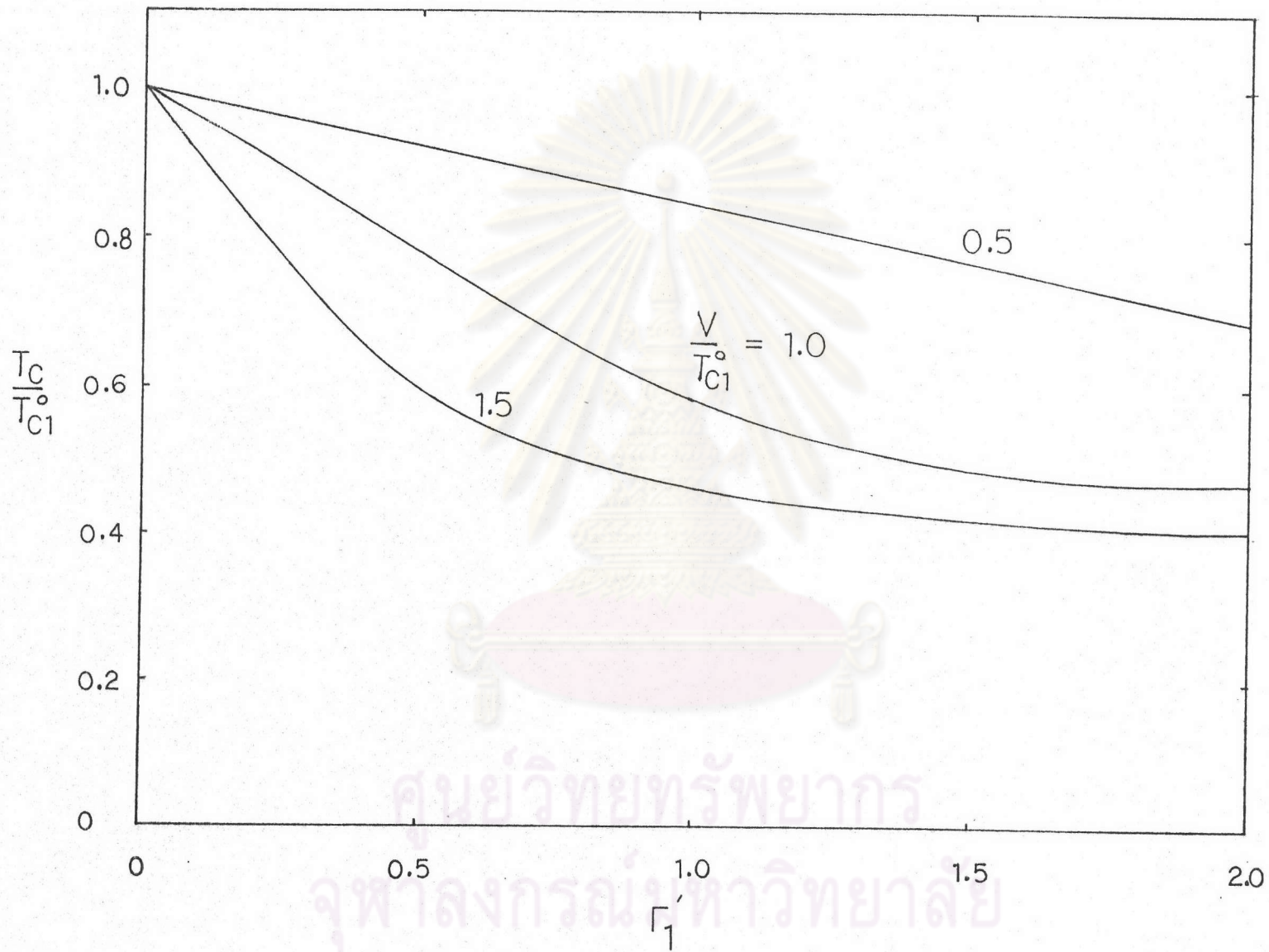


Fig. 3.2 Critical temperature T_c of S_1 - S_2 sandwich vs. Γ_1 for various values of V/T_{c1} . The parameters used are $\lambda_1=0.2447$, $\lambda_2=0.1669$, $T_{c1}^*=3.72$ K (Mohabir and Nagi, 1979).

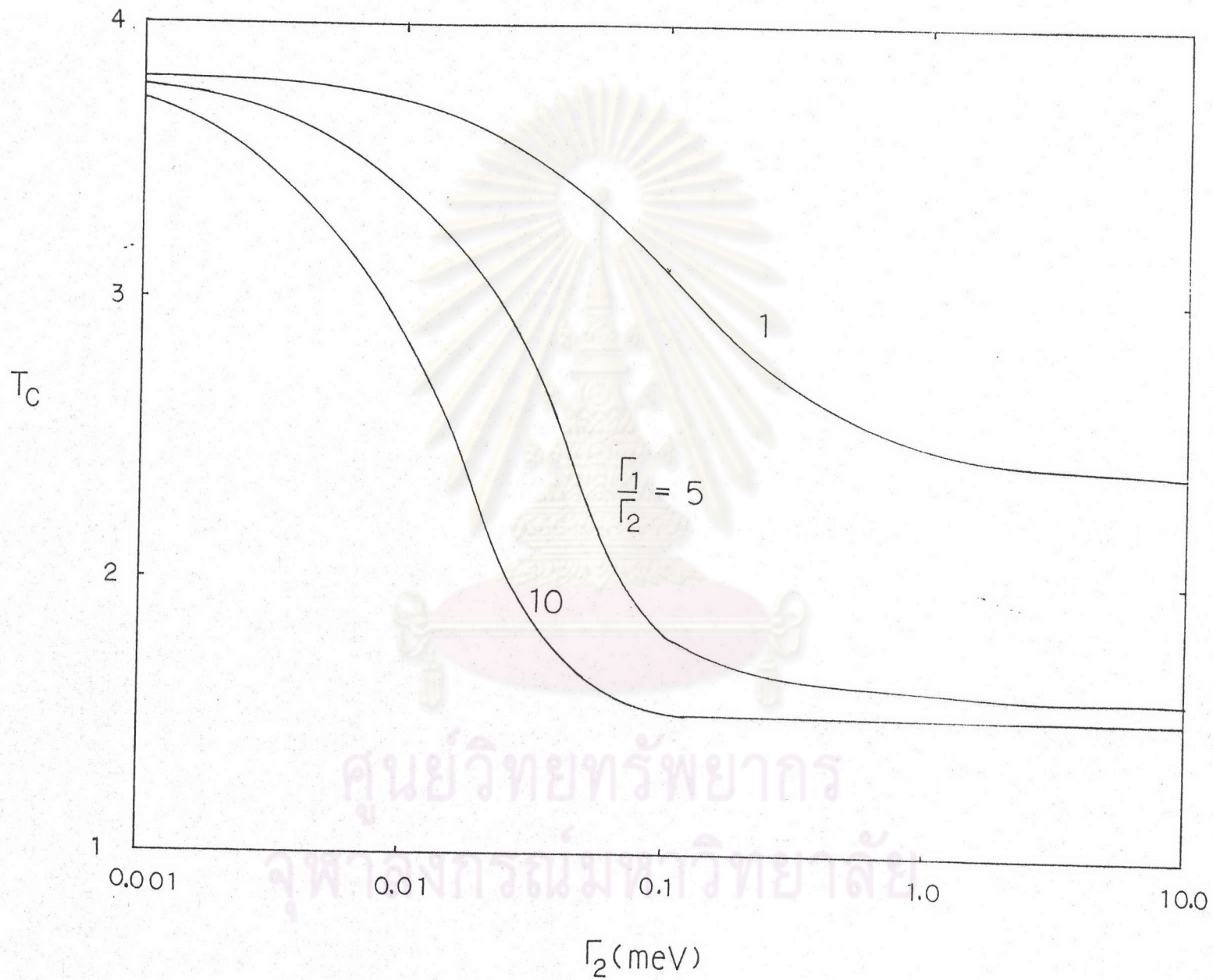


Fig. 3.3 Critical temperature T_c of an S_1 - S_2 sandwich vs. Γ_2 for various values of Γ_1/Γ_2 . The parameters used are $\lambda_1 = 0.246$, $\lambda_2 = 0.171$, $\omega_{D1} = 16.78$ meV, $\omega_{D2} = 32.21$ meV (Mohabir and Nagi, 1979).

In Fig. (3.2), we have introduced dimensionless parameter Γ_1' ($\Gamma_1' = \pi A d_2 N_2(0) T_{c1}^*$). When $\Gamma_1' = 0$, the ratio of T_c/T_{c1}^* is equal to 1. We found that for fixed value of Γ_1' and V/T_{c1}^* , ratio T_c/T_{c1}^* of an S_1 - S_2 sandwich decreases more slowly than the ratio T_c/T_{cs}^* of S-N sandwich.

In Fig. (3.3), we show calculated values of the critical temperature of an S_1 - S_2 sandwich as a function of Γ_2 (for various values of Γ_1/Γ_2). For fixed values of Γ_2 , the values of T_c decreases when Γ_1/Γ_2 increases.



ศูนย์วิทยทรัพยากร
จุฬาลงกรณ์มหาวิทยาลัย