

INELASTIC ELECTRON-POSITRON SCATTERING

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Abstract : The two-photon exchange in the inelastic Bhabha scattering $e^+e^- \rightarrow e^+e^- + \text{anything}$ is studied without the quasi real photon approximation.

Résumé : L'échange de deux photons dans la diffusion inélastique $e^+e^- \rightarrow e^+e^- + \Gamma$ est étudié sans l'approximation des photons quasi réels.



I wish to report very briefly on a recent work done in Paris on the two-photon contributions to inelastic Bhabha scattering

$$e^+ + e^- \Rightarrow e^+ + e^- + \Gamma \quad (1)$$

where Γ is a one or multiparticle system not detected.

1°) More precisely we have computed the dominant diagram at high energy

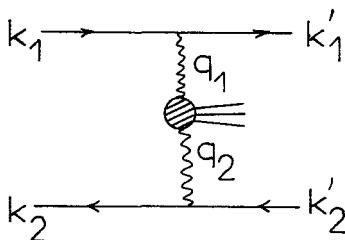


Fig. 1

The knowledge of the cross section for the process (1) will give information on the two virtual photon annihilation reaction

$$\gamma + \gamma \Rightarrow \Gamma \quad (2)$$

When the particles entering the system Γ of effective mass W are not detected, we can define 10 structure functions depending on q_1^2 , q_2^2 and W^2 only. If time reversal holds, this number is reduced to 8.

An experiment with unpolarized electrons and positrons is able to measure only 6 structure functions, and we get

$$\frac{E'_1}{d_3 k'_1} \frac{E'_2}{d_3 k'_2} d\Gamma = \frac{\alpha^2}{2\pi^4} \frac{1}{3} \frac{KW}{q_1^2 q_2^2} \left\{ \Gamma_{LL} K_{LL} + \Gamma_{LT} K_{LT} + \Gamma_{RL} K_{RL} + \Gamma_{RR} K_{RR} + (\Gamma_{II} - \Gamma_{\perp\perp}) K_{TT}^{\text{ex}} + (\Gamma_{\perp\perp} + \Gamma_{II}) K_{LT}^{\text{ex}} \right\}$$

where K is the c.m. momentum of the $\gamma\gamma$, S the square of the total incident energy $S = (E_1 + E_2)^2$.

The function $K_{\lambda\lambda'}$ are known functions of \vec{k}_1^{λ} and $\vec{k}_2^{\lambda'}$ but obviously they depend only on 5 variables because of an overall symmetry around the incident axis of the beams. The structure functions depending only on 3 variables, the dependence of the cross section on the remaining 2 variables is known as a consequence of the two-photon exchange approximation and is completely contained in the functions $K_{\lambda\lambda'}$.

The two main features of this process is the smallness of the electron mass and the presence of the photon propagators. The poles existing at $q_1^2 = 0$ and $q_2^2 = 0$ dominate the cross section and suggests the use of a quasi-real photon approximation. Our calculations are free of such an approximation.

2°) The calculations have been made for a state Γ free of any e^+e^- pair.

(a) The differential cross section $\frac{d^3\sigma}{dq_1^2 dq_2^2 dW^2}$ has been computed exactly, integrating analytically over the two kinematical variables [1].

(b) The dominant contribution of $\frac{d\sigma}{dW^2}$ is associated to the total cross section $\bar{\sigma}_{\gamma\gamma}(0,0,W^2)$ for process (2) with real photons

$$\frac{d\sigma_{\gamma\gamma}}{dW^2} = \left(\frac{a}{\pi}\right)^2 \frac{\bar{\sigma}_{\gamma\gamma}(0,0,W^2)}{W^2} F_{\gamma\gamma}(W^2, s) \quad (3)$$

The function $F_{\gamma\gamma}(W^2, s)$ can be exactly computed and is independent of any model. It exhibits a logarithmic dependence in the energy we can order when W^2 is small as compared to s , according to the number of logarithms involved as

$$3\text{-Log} + 2\text{-Log} + 1\text{-Log} + C = O\left(\frac{W^2}{s}\right)$$

For instance we get, for the 3-Log terms [1]

$$(\log \frac{s}{m^2})^2 \log \frac{s}{w^2} - \frac{1}{3} (\log \frac{s}{w^2})^3 \quad (4)$$

(c) The other contributions of $\frac{d\mathcal{G}}{dw^2}$ which are model dependent and where enter the other helicity states and the dependence in q_1^2 and q_2^2 of the structure functions [2].

(d) The total cross section $\mathcal{G}(s)$ for the multiparticle states Γ [2].

3°) Some applications of these calculations have been investigated [2].

(a) The first one concerns the $\mu^+ \mu^-$ production where the quantum electrodynamics gives us a model for an exact and complete calculation. In particular, the differential $\frac{d}{dw^2} \mathcal{G}(s, w^2)$ and total $\mathcal{G}(s)$ cross sections have been computed. The comparison of $\mathcal{G}(s)$ with previous calculations is as follows, see fig. 2.

1.- With formula (3) of Bonneau-Gourdin-Martin, exactly computed, the agreement is better than 2%.

2.- With the Baier-Fadin [3] calculation, the 3-Log, 2-Log, and 1-Log terms are identical, but in that paper the $\mathcal{O}(w^2)$ and $\mathcal{O}(\frac{m^2}{s})$ terms have not been evaluated so that we have a nice agreement asymptotically, but at intermediate energy, Baier and Fadin make too small an evaluation, having an incorrect threshold.

3.- With Brodsky-Kinoshita-Terazawa [4] 's calculation where the pole term $(\log \frac{E}{m})^2$ only is taken into account, giving systematically too high a cross section. Of course the threshold is there correct and the disagreement is about 15-20% at intermediate energy where the future machine will operate and of course 50% at infinite energy where the second term of equation (4) plays a role.

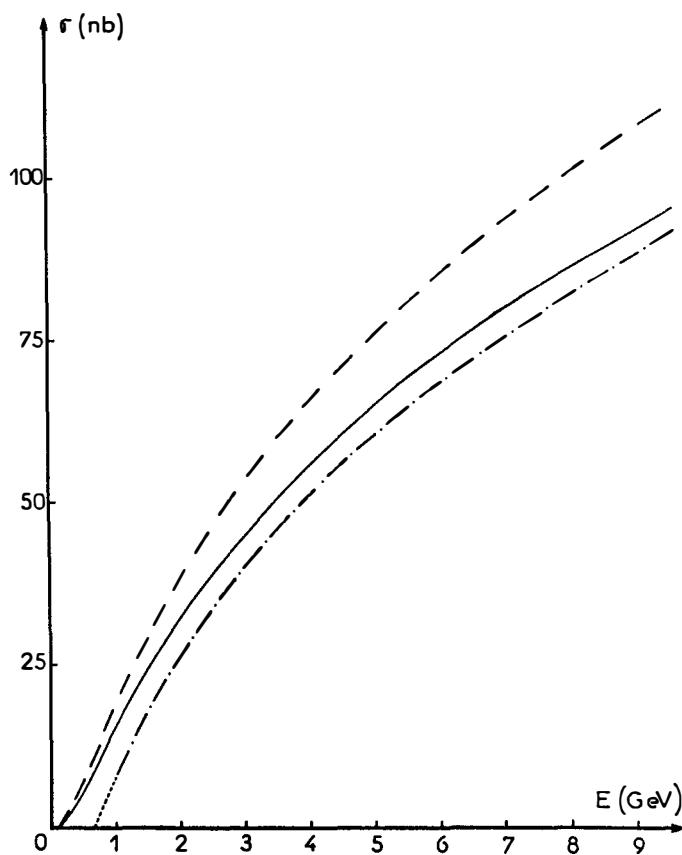


Fig. 2 - The total cross section for the process $ee \rightarrow ee \mu^+ \mu^-$ as a function of E :
 — our complete result
 - - - result of ref. [4]
 - - - result of ref. [3]

(b) The second application of the general calculations concerns the production of pseudoscalar meson $\Gamma = \pi^0, \Sigma, \Lambda^0$. As in reference [4] a vector meson dominance type model has been used with one parameter M in order to describe the variation with q_1^2 and q_2^2 of the amplitude

For π^0 production, the model independent part gives 3/4 of the cross section and the model dependent part 1/4 for M of the order of the ρ meson mass. If M varies between 5 and 1.5 GeV, then the model dependent part can change by 30% so that the computed cross section varies by approximatively 5%-10%.

For Σ and Λ^0 production, the situation is better in the sense that the model independent part is relatively more important so that if the model dependent part is more sensible to the mass parameter M as in the π^0 case, the total prediction is not too sensible to that parameter.

Of course these computations agree with similar ones of Brodsky, Kinoshita, Terazawa [4] corresponding to specific values of M .

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