

Comparison of energy dependence of transverse momentum of dimuons produced in pN and π^-N interactions with quantum-chromodynamic predictions

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We have analyzed the dependence of the average transverse momentum squared $\langle p_T^2 \rangle$ of dimuons produced in pN and π^-N interactions on s , the square of the center-of-mass energy. The presently available data indicate a linear increase of $\langle p_T^2 \rangle$ with s in both reactions, with the rate of increase being approximately twice as large in the π^-N reaction for $M/\sqrt{s}=0.28$ as in the pN reaction for $M/\sqrt{s}=0.22$. The value of $\langle p_T^2 \rangle$ extrapolated to $s=0$, which is interpreted as the intrinsic $\langle k_T^2 \rangle$ of the π^- and p constituents, is the same within errors for both reactions. First-order QCD consistently predicts a lower $\langle p_T^2 \rangle$ at all s than is observed in the data.

A well known prediction of perturbative quantum chromodynamics (QCD) is that the average transverse momentum squared $\langle p_T^2 \rangle$ of dimuons produced in hadronic collisions is given to order α_s by the relationship¹⁻³

$$\langle p_T^2 \rangle = \langle k_T^2 \rangle + \alpha_s(Q^2) f(\tau, x_F, \ln Q^2) s, \quad (1)$$

where α_s is the running coupling constant, M is the mass of the dimuon, s is the center-of-mass energy squared, $\tau=M^2/s$, and $\langle k_T^2 \rangle$ is the sum of the intrinsic transverse momentum squared of the constituents that take part in

the interaction. Equation (1) implies that $\langle p_T^2 \rangle$ should grow linearly with s at a given τ , Q^2 , and x_F . A linear relationship has been reported.⁴⁻⁷ The aim of this paper is to extract quantitative information from presently existing data on the s variation of $\langle p_T^2 \rangle$ in π^-N and pN reactions and to compare the observed data with the predictions of first-order QCD.

The perturbative expression for the differential cross section to first order in α_s for dimuon production can be written as⁸

$$\begin{aligned} \frac{d\sigma_p}{dM dy dp_T^2} &= \frac{d\sigma_A}{dM dy dp_T^2} + \frac{d\sigma_C}{dM dy dp_T^2} \\ &= \frac{16\alpha_s^2}{27} \frac{s}{M} \int_{x_1^{\min}}^1 dx_1 \frac{x_1 x_2 \alpha_s}{x_1 s + u - M^2} \frac{(\hat{t} - M^2)^2 + (\hat{u} - M^2)^2}{s^2 \hat{u}} \sum_i e_i^2 [q_i^{h_1}(x_1, Q^2) \bar{q}_i^{h_2}(x_2, Q^2) + (1 \leftrightarrow 2)] \\ &\quad + \frac{2\alpha_s^2}{9} \frac{s}{M} \int_{x_1^{\min}}^1 dx_1 \frac{x_1 x_2 \alpha_s}{x_1 s + u - M^2} \frac{\hat{s}^2 + \hat{u}^2 + 2M^2 \hat{t}}{-\hat{s}^3 \hat{u}} \sum_i e_i^2 [q_i^{h_1}(x_1, Q^2) G_i^{h_2}(x_2, Q^2) + (1 \leftrightarrow 2)]. \end{aligned} \quad (2)$$

Here σ_A and σ_C represent the contributions of the order- α_s quark-antiquark-annihilation and Compton-scattering diagrams shown in Figs. 1(b) and 1(c), respectively. The q_i 's and G_i 's in Eq. (2) are the quark and gluon distribution functions, and the various kinematic variables are given by

$$\begin{aligned} x_1^{\min} &= -u/(s+t-M^2), \\ x_2 &= [-x_1 t - (1-x_1)m^2]/(x_1 s + u - m^2), \\ t &= M^2 - \sqrt{s} M_T e^{-y}, \quad u = M^2 - \sqrt{2} M_T e^y, \\ \hat{s} &= x_1 x_2 s, \quad \hat{t} = x_1 t + (1-x_1)M^2, \\ \hat{u} &= x_2 u + (1-x_2)M^2, \\ M_T^2 &= p_T^2 + M^2, \end{aligned}$$

and

$$\alpha_s = 12\pi/[25 \ln(Q^2/\Lambda^2)]$$

with $Q^2=M^2$ in expression (2).

Equation (2) is useful only for large p_T since it has a singularity at $p_T=0$. Furthermore, it ignores the intrinsic transverse momentum k_T of the constituents. Altarelli, Parisi, and Petronzio¹ have given the following prescription which uses this intrinsic k_T to regularize the $p_T=0$ divergence:

$$\begin{aligned} \frac{d\sigma}{dM dy dp_T^2} &= \int d^2 q_T \frac{d\sigma_p}{dM dy dp_T^2} [f(\vec{p}_T - \vec{q}_T) - f(p_T)] \\ &\quad + \pi f(p_T) \frac{d\sigma_{DY}}{dM dy}. \end{aligned} \quad (3)$$

Here $\vec{q}_T = \vec{p}_T - \vec{k}_T$, σ_{DY} is the Drell-Yan cross section [see Fig. 1(a)], and f is the soft k_T distribution. We have

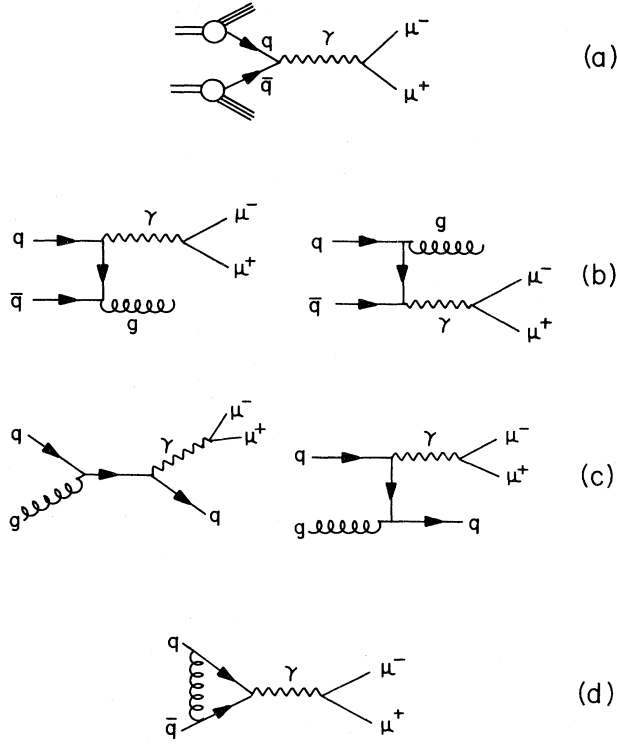


FIG. 1. (a) Drell-Yan process. (b) $O(\alpha_s)$ QCD quark-antiquark-annihilation process. (c) $O(\alpha_s)$ QCD Compton-scattering process. (d) Vertex correction to the Drell-Yan process with virtual gluon exchange.

assumed the k_T distribution to have a Gaussian form with the normalization

$$\int f(p_T) d^2 p_T = \frac{1}{\pi \langle k_T^2 \rangle} \int d^2 p_T \exp[-p_T^2 / \langle k_T^2 \rangle] = 1. \quad (4)$$

The structure functions used in this calculation of $\langle p_T^2 \rangle$ are given by

$$\begin{aligned} xu_p &= u_0 x^{0.52-0.16\bar{s}} (1-x)^{2.79+0.77\bar{s}}, \\ xd_p &= d_0 x^{0.52-0.16\bar{s}} (1-x)^{3.79+0.77\bar{s}}, \\ xS_p &= (0.26+0.18\bar{s})(1-x)^{7.8+0.78\bar{s}}, \\ xG_p &= 3.06(1-x)^{5.0}, \\ xV_\pi &= V_0 x^{0.4}(1-x)^{0.9}, \\ xS_\pi &= 0.24(1-x)^{6.9}, \\ xG_\pi &= 2.0(1-x)^{3.0} \end{aligned} \quad (5)$$

with $\bar{s} = \ln[(Q^2/0.25)/\ln(20/0.25)]$, where Q^2 is in GeV^2 . The valence-quark distributions are normalized to obtain the proper number of valence quarks for π 's and protons. We have assumed that the strange quark-antiquark pairs are suppressed by a factor of 2 compared to $u\bar{u}$ or $d\bar{d}$ pairs. The quark distribution functions for the proton

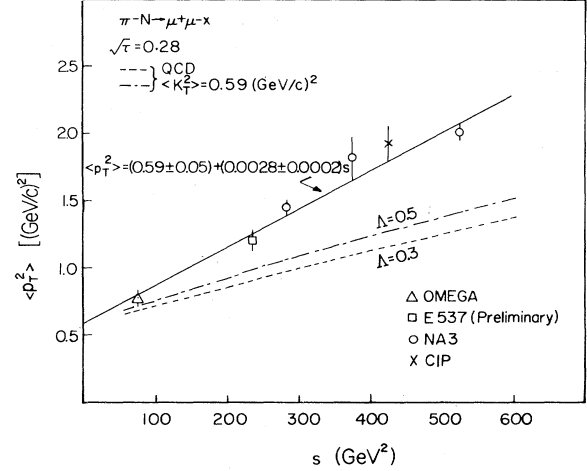


FIG. 2. $\langle p_T^2 \rangle$ versus s for dimuons produced in π^- -nucleon interactions. The solid curve is a linear fit to the data. The dashed and dot-dash curves are the predictions of first-order QCD using the Altarelli *et al.* prescription for different values of Λ .

correspond to the CERN-Dortmund-Heidelberg-Saclay (CDHS) results⁹ while the quark distribution functions for the pion correspond to the NA3 results.^{6,10} The Q^2 dependence is quite small over the range of interest. We have used a Q^2 -independent gluon distribution for the proton corresponding to counting rules; the conclusions arrived at here are not affected if the power index is varied in the range 5.0 ± 1.0 .

When we use the structure functions (5) in Eq. (3) and integrate over p_T , we obtain the Q^2 -dependent Drell-Yan cross section which is known to be lower than the experimental cross sections for dimuon production by π 's, p 's, or antiprotons by factors in the range 1.6 to 2.6 [compilations of K factors $\equiv (d\sigma/dM dy)_{\text{data}} / (d\sigma/dM dy)_{\text{DY}}$ for various reactions are given, for example, in Refs. 7 and 11–14]. At the level of the accuracy of the data these K

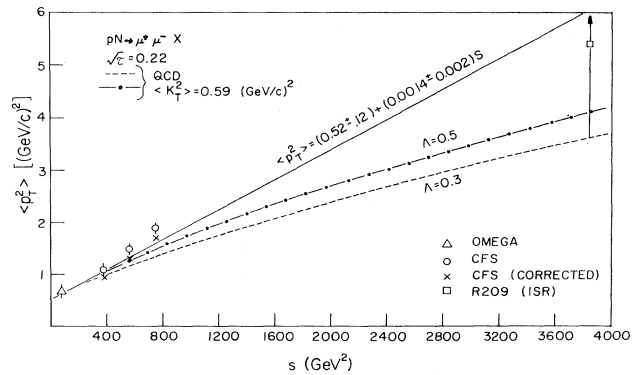


FIG. 3. $\langle p_T^2 \rangle$ vs s for dimuons produced in p -nucleon interactions. The solid curve is the linear fit to the data. The dashed and dot-dash curves are the predictions of first-order QCD using the Altarelli *et al.* prescription for different values of Λ .

TABLE I. Fits of $\langle p_T^2 \rangle = A + B(\tau)s$ to existing dimuon data.

Fit	Reaction	$\sqrt{\tau}$	Intercept (A) [(GeV/c) ²]	Slope (B)	χ^2/DF
$\langle p_T^2 \rangle$	$\pi^- N$	0.28	0.59±0.05	$(2.8 \pm 0.2) \times 10^{-3}$	1.2
$\langle p_T^2 \rangle$	pN	0.22	0.52±0.11	$(1.4 \pm 0.2) \times 10^{-3}$	0.3

factors have been determined to be independent of τ and y . Theoretical justifications of constant K factors have been provided by several authors¹⁵ who have carried out a full $O(\alpha_s)$ calculation and have confirmed the observed independence of K of τ and y for $\tau < 0.5$. In these calculations the largest part of correction to the cross section for $\tau \leq 0.3$ (i.e., for the range of the current experiments) is found to arise from the vertex-correction diagram [Fig. 1(d)]. This diagram, however, contributes no p_T other than the k_T of the constituents to the dimuons. In addition, more complete calculations to order α_s^2 which have been done by Ellis, Martinelli, and Petronzio,¹⁶ indicate that K is only a slowly varying function of p_T . Therefore, we have used empirical K factors assumed to be independent of τ , x_F , and p_T in fitting the dimuon p_T distributions at various energies.

Using Eq. (3), the structure functions given by (5), and the empirical K factors, we have been able to fit¹³ published dimuon transverse-momentum distributions for proton interactions at 400 and 2050 GeV/c, for π^- interactions from 125 to 280 GeV/c, and for \bar{p} interactions at 125 GeV/c. While the agreement of the calculations with the proton data is quite good at a given s , a $\langle k_T^2 \rangle$ varying from 0.9 (GeV/c)² at $\sqrt{s} = 23.4$ GeV to 1.5 (GeV/c)² at $\sqrt{s} = 62$ GeV is required. Similarly $\langle k_T^2 \rangle$ must vary from 0.9 (GeV/c)² at $\sqrt{s} = 15.3$ GeV to 1.2 (GeV/c)² at $\sqrt{s} = 22.8$ GeV for the π^- reaction. Thus, we find that QCD to order α_s with a constant $\langle k_T^2 \rangle$ is unable to describe the existing p_T data at different energies.¹³

To make explicit the failure of first-order QCD with constant $\langle k_T^2 \rangle$ to explain the transverse-momentum distribution of dimuons at different energies, we have compared the observed second momenta of the p_T distributions of the data to the $\langle p_T^2 \rangle$ predicted by (3). Figure 2 shows the existing measurements^{3,6,17,18} of $\langle p_T^2 \rangle$ for $\sqrt{\tau} = 0.28$ and $x_F > 0$ at different values of s for the $\pi^- N$ reaction. Figure 3 shows $\langle p_T^2 \rangle$ for the pN data^{5,19,20} for $\sqrt{\tau} = 0.22$. The pN data points from the Columbia-Fermilab-Stony Brook (CFS) experiment¹⁹ are for $\langle y \rangle$ ranging from 0.03 for 400 GeV/c to 0.40 for 200 GeV/c. To take into account the expected dependence of $\langle p_T^2 \rangle$ on y we have used the QCD formalism described above to correct the CFS data to correspond to $y \geq 0$. Linear fits of the form $\langle p_T^2 \rangle = A + Bs$ are shown for both sets of data. The details of the fits are given in Table I. As shown in the figures and in Table I, the $s = 0$ intercept is the same within errors for the pN and $\pi^- N$ data but the slope of the rise of $\langle p_T^2 \rangle$ with s is approxi-

mately twice as large for the π^- data as for the pN data.

Since within errors the intercepts $\langle P_T^2 \rangle_{s=0}$ are the same and since we interpret $\langle P_T^2 \rangle_{s=0}$ to be equal to $\langle k_T^2 \rangle$ from Eq. (1), we have set $\langle k_T^2 \rangle = 0.59$ (GeV/c)² in the QCD calculations. The results of the calculation of $\langle P_T^2 \rangle$ with this choice of $\langle k_T^2 \rangle$ are shown in Figs. 2 and 3 for two different values of Λ (0.3 and 0.5 GeV/c). The predictions of QCD fall well below the data in the case of both the pN and the $\pi^- N$ reactions. However, the calculations do show a steeper slope for $\pi^- N$ than for pN in qualitative agreement with the data.

In carrying out the computations we have restricted the calculations to $p_T \leq 10$ GeV/c. In order to check the effect of this truncation we have repeated the calculation for $s = 3850$ GeV² (since the effect is expected to be maximum at large s) but extending p_T to $p_T \leq 15$ GeV/c. (Note that kinematics restrict the maximum value of p_T to 29.5 GeV/c for $\sqrt{\tau} = 0.22$.) For $\Lambda = 0.5$ GeV/c, we obtain $\langle p_T^2 \rangle = 4.20$ (GeV/c)² compared to $\langle p_T^2 \rangle = 4.07$ (GeV/c)² for $p_T \leq 10$ GeV/c. Thus, the effect of truncation of the p_T distribution is fairly small for $s \leq 3850$ GeV². In order to explore the source of the curvature seen in Fig. 3, we have repeated calculations by using a constant value of 0.3 for α_s . This yields a very nearly straight-line behavior for $\langle p_T^2 \rangle$ vs s . Thus, the source of the curvature is almost entirely due to the $\ln(Q^2)$ dependence of α_s .

In conclusion we have examined the behavior of $\langle p_T^2 \rangle$ as a function of s for the high-mass dimuons produced in pN and $\pi^- N$ interactions. We find that the data are consistent within errors with the same $\langle p_T^2 \rangle$ intercept at $s = 0$ for p and π^- , but that the rise of $\langle p_T^2 \rangle$ with s is approximately twice as fast for the $\pi^- N$ reaction as for the pN reaction. It may, however, be noted that the values of τ used for pN ($\sqrt{\tau} = 0.22$) and $\pi^- N$ ($\sqrt{\tau} = 0.28$) are different which, in principle, would lead one to expect different slopes for the two cases. We have interpreted the $s = 0$ intercept as the sum of the $\langle k_T^2 \rangle$ of the constituents of the proton and pion participating in the interaction and have used the observed intercept $\langle k_T^2 \rangle = 0.59$ GeV²/c² in the first-order QCD prescription of Altarelli *et al.* to calculate the expected behavior of $\langle p_T^2 \rangle$ with s . We find that this $O(\alpha_s)$ formalism of QCD to (with the assumption of a p_T -independent K factor and $Q^2 = M^2$) is unable to account for the observed energy dependence of $\langle p_T^2 \rangle$ with the disagreement worsening at larger s . We therefore conclude that if the p_T distribution of the dimuons has to be understood on the basis of QCD, it is necessary to consider diagrams representing higher orders in α_s .

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