

Longitudinal-Transverse Separations of Deep-Inelastic Structure Functions at Low Q^2 for Hydrogen and Deuterium

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We report on a study of the longitudinal to transverse cross section ratio, $R = \sigma_L/\sigma_T$, at low values of x and Q^2 , as determined from inclusive inelastic electron-hydrogen and electron-deuterium scattering data from Jefferson Laboratory Hall C spanning the four-momentum transfer range $0.06 < Q^2 < 2.8 \text{ GeV}^2$. Even at the lowest values of Q^2 , R remains nearly constant and does not disappear with decreasing Q^2 , as might be expected. We find a nearly identical behavior for hydrogen and deuterium.

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Since the early experiments at the Stanford Linear Accelerator Center (SLAC), which discovered the substructure of the nucleon and led to the development of the quark parton model [1], deep-inelastic scattering (DIS) has been a powerful tool in the investigation of the partonic substructure of the nucleon. After decades of experiments with electron and muon beams, the nucleon structure function $F_2(x, Q^2)$ is known with high precision over many orders of magnitude in x and Q^2 [2]. Here, Q^2 is the negative square of the four-momentum transfer of the exchanged virtual photon in the scattering process. The Bjorken scaling variable $x = Q^2/2M\nu$, with M the nucleon mass and ν the energy transfer, can be interpreted as the fraction of the nucleon momentum carried by the struck parton.

In the region of large Q^2 and ν , the results of DIS measurements are typically interpreted in terms of partons (quarks and gluons). In this case, the theoretical framework is provided by perturbative quantum chromodynamics

(pQCD), which includes logarithmic scaling violations. This description starts to fail when nonperturbative effects such as higher twist interactions between the struck quark in the scattering process and other quarks or gluons in the nucleon become important. The sensitivity for higher twist effects increases with decreasing Q^2 , since they are proportional to powers of $1/Q^2$.

There is great interest in the behavior of the nucleon structure functions in the low Q^2 region where the transition from perturbative to nonperturbative QCD takes place. However, little is known about this behavior, since at large invariant mass W of the hadronic system there are few data points in this region, except for the (transverse) cross section σ_T at exactly $Q^2 = 0$, which is accessible through real photon absorption experiments. The more plentiful data at low W are typically interpreted in terms of nucleon resonance excitations.

The differential cross section for inclusive electron scattering, after dividing by the virtual-photon flux factor (Γ),

can be written as

$$\frac{1}{\Gamma} \frac{d^2\sigma}{d\Omega dE'} = \sigma_T + \varepsilon\sigma_L, \quad (1)$$

where ε is the virtual-photon polarization and σ_L (σ_T) is the longitudinal (transverse) virtual-photon absorption cross section, which depends on x and Q^2 . Current conservation determines the behavior of the structure functions for $Q^2 \rightarrow 0$, leading to

$$R(x, Q^2) \equiv \frac{\sigma_L}{\sigma_T} = \frac{F_L}{2xF_1} = \mathcal{O}(Q^2), \quad (2)$$

where F_L and F_1 are the longitudinal and transverse nucleon structure functions. The value of Q^2 at which this behavior becomes manifest is, however, neither predictable nor yet observed.

While there is a wealth of data for $F_2 = (2xF_1 + F_L)/(1 + Q^2/\nu^2)$, relatively few data exist for F_L , or equivalently R . Data on $R(x, Q^2)$ on hydrogen and deuterium are available in a limited x and Q^2 ($Q^2 > 1 \text{ GeV}^2$) range, with a typical uncertainty of 0.1–0.2 [3,4], comparable to the size of R itself. For scattering from pointlike spin-1/2 particles, R should vanish at large values of Q^2 because of helicity conservation. At low values of Q^2 , however, R is not small, and typical values are about 0.3.

Precision data on R are necessary for several reasons. Most importantly, determinations of the structure function F_2 from cross section measurements, and the parton distributions derived therefrom, need numerical values for R . If the former are not based on longitudinally and transversely separated measurements, the uncertainties in F_2 introduced by assumptions for R can be as large as 20%. Furthermore, especially in the low Q^2 region, data are needed to study higher twist effects and to search for the onset of the current conservation behavior of the structure functions at low Q^2 described above.

The determination of R is typically accomplished via a Rosenbluth-type separation technique, which requires high precision measurements of the cross section at fixed values of x and Q^2 , but at different values of ε . This technique requires the use of at least two beam energies and correspondingly different scattered-electron angles θ . Only in some experiments have such measurements actually been performed [3–10].

At sufficiently high Q^2 , pQCD predicts R to be the same for protons and neutrons. Indeed, previous results have shown $R^D - R^H$, the difference in R from hydrogen and deuterium targets, to be consistent with zero [3,9,10]. A low values of Q^2 , however, nonperturbative effects might result into R^H being different from R^D .

In this Letter we present results from a study of R for both hydrogen and deuterium at low values of Q^2 . The experiment (E99-118) was carried out at the Thomas Jefferson National Accelerator Facility (Jlab). Data were obtained at $0.007 < x < 0.55$ and $0.06 < Q^2 < 2.8 \text{ GeV}^2$,

by utilizing 2.301, 3.419, and 5.648 GeV electron beams at a current of $I = 25 \mu\text{A}$. The minimum scattered-electron energy was $E' \approx 0.4 \text{ GeV}$ and the range of the invariant mass squared of the hadronic system W^2 was between 3.5 and 10 GeV^2 . Electrons scattered from 4 cm long liquid hydrogen (H) and deuterium (D) targets were detected in the high-momentum spectrometer in Hall C at various angles between 10° and 60° .

The inclusive double differential cross section for each energy and angle bin within the spectrometer acceptance was determined from

$$\frac{d\sigma}{d\Omega dE'} = \frac{Y_{\text{corr}}}{L\Delta\Omega\Delta E'}, \quad (3)$$

where $\Delta\Omega$ ($\Delta E'$) is the bin width in solid angle (scattered energy), L is the total integrated luminosity, and Y_{corr} is the measured electron yield after correcting for detector inefficiencies, background events, and radiative effects.

To account for backgrounds from π^0 production and its subsequent decay into two photons followed by pair production of electron-positron pairs, positron data were also taken by reversing the polarity of the spectrometer. Other background contributions include electron scattering from the aluminum walls of the cryogenic target cells and electroproduced negatively charged pions. Events from the former were subtracted by performing substitute empty target runs, while events from the latter were identified and removed by use of both a gas Cherenkov counter and an electromagnetic calorimeter. Additional details regarding the analysis and the standard Hall C apparatus employed in this experiment can be found in Ref. [11].

Radiative effects including bremsstrahlung, vertex corrections, and loop diagrams are calculated using the approach by Bardin *et al.* [12]. Additional radiative effects in the target and its exit windows were determined using the formalism of Mo and Tsai [13]. The calculation of such effects includes the emission of one hard photon. There is, however, the possibility that the electron could emit two hard photons. The calculation of this process is unfortunately not fully established and the corresponding effect was therefore treated in the present analysis as a separate uncertainty.

For each energy bin, a weighted average cross section over θ within the spectrometer acceptance was obtained by using a model to correct for the angular variation of the cross section from the central angle value. To minimize the dependence on the model used to compute both this correction and the radiative effects, an iterative procedure was employed.

The total systematic uncertainty in the differential cross section was taken as the quadratic sum of all the systematic uncertainties contributing to the cross section measurement. In a Rosenbluth separation one needs to distinguish between uncertainties that are correlated between measurements at different ε , such as uncertainties in target thickness and integrated charge, and uncorrelated ones. Not

including the contributions from radiative corrections, the uncorrelated systematic uncertainties on the cross section measurements in this experiment amounted to 0.9%, while the total systematic uncertainty was 1.35%.

The size, and consequently the uncertainty, of the radiative effects strongly depends on the kinematics and is largest at low values of E' where the measured cross section is dominated by events from elastic or quasielastic scattering with the emission of one or more photons in the initial or final state. The estimate of these uncertainties was determined by varying all relevant input cross sections within their uncertainties, and amounted to as much as 1.5% for hydrogen and 8.5% for deuterium in the most extreme cases considered. It rapidly decreased for higher values of E' , i.e., higher values of x and Q^2 . The much larger uncertainty in the deuterium cross section is due to the contribution from quasielastic scattering which can only be modeled approximately.

The extractions of purely longitudinal and transverse cross sections and structure functions were accomplished via the Rosenbluth technique, where the reduced cross section is fit linearly as a function of ε , as in Eq. (1). The intercept of such a fit gives the transverse cross section σ_T (and therefore the structure function $F_1(x, Q^2)$), while the slope gives the longitudinal cross section σ_L , from which the structure functions $R(x, Q^2)$ or $F_L(x, Q^2)$ can be extracted.

The results for $R(x, Q^2)$ for hydrogen are shown in Fig. 1 as a function of Q^2 for fixed values of x (open squares). The inner error bars represent the combined statistical and uncorrelated systematic uncertainties. The total error bars represent the statistical and systematic uncertainties added in quadrature. The data are compared to the results of previous measurements at Jlab (E94-110) [7], SLAC [3], and by the EMC [8], NMC [9], and BCDMS [4] collaborations at CERN.

The structure function F_2 determined through the Rosenbluth separation technique was found to agree to better than 2% with a Regge-motivated parametrization of all previously available deep-inelastic scattering data [14], even at very low values of Q^2 [11]. This facilitated the utilization of an alternative approach where R was calculated using this parametrization (including a 2% uncertainty) for the structure function $F_2(x, Q^2)$ from

$$\frac{d^2\sigma}{d\Omega dE'} = \sigma_{\text{Mott}} \frac{2MxF_2}{Q^2\varepsilon} \left(\frac{1 + \varepsilon R}{1 + R} \right). \quad (4)$$

The results (shown by the solid circles in Fig. 1) cover a larger kinematic range than the results from the Rosenbluth separation, since the measurement of this experiment is combined via the model with results of previous experiments at different energies. The inner error bars represent the statistical uncertainty and the total error bars represent the quadratic sum of the statistical uncertainty and the uncertainty in the radiative corrections due to the possible emission of two hard photons. The shaded bands represent

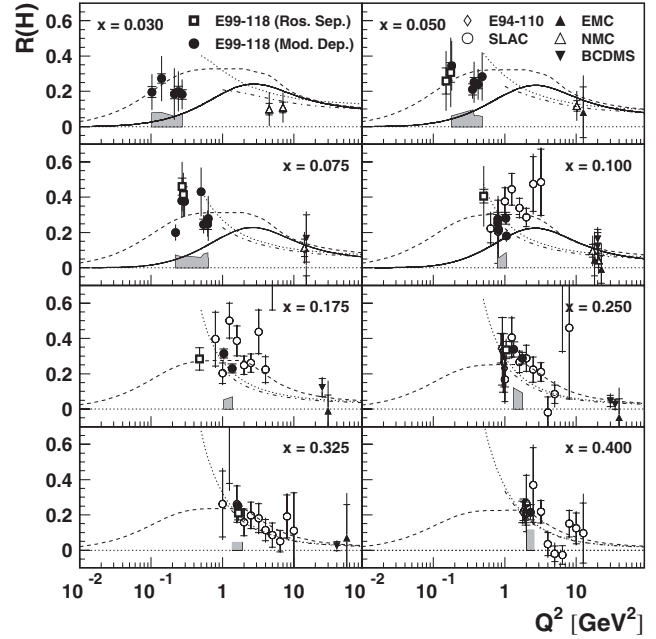


FIG. 1. Comparison of the values of $R(x, Q^2)$ for hydrogen from the present experiment (E99-118) to the results of other experiments. The dashed curves represent the parametrization $R_{e99118}^H(x, Q^2)$ and the solid curves the model developed in [15] (see text for details). The dot-dashed curves are a next-to-next-to-leading order calculation based on the recent MRST parton distributions [16], and the dotted curves show the next-to-leading order result of the GRV model [17]. In both cases target mass effects have been included according to [18].

the uncertainties in the radiative effects due to uncertainties in the input cross sections. Especially in the model dependent extraction, the uncertainties are dominated by the uncertainties in the radiative corrections which are correlated between data points. The results from the second method agree very well with those obtained from the Rosenbluth separation method. Good agreement is also found with previous experiments in the regions of x and Q^2 where the data overlap.

The dashed curve in Fig. 1 represents a new parametrization of R ($R_{e99118}^H(x, Q^2)$) [11] based on all available data including those from this experiment. The functional form of the parametrization has been chosen to satisfy the condition that R vanishes as Q^2 goes to zero. At $Q^2 = 2 \text{ GeV}^2$, it was connected to a previously obtained parametrization from SLAC [3] that is based on measurements at higher values of Q^2 . The solid curves in the upper four panels of Fig. 1 show, within its range of applicability ($x \leq 0.1$), the model developed in [15], which is based on the photon-gluon fusion mechanism suitably extrapolated into the region of low Q^2 .

As mentioned above, at low values of Q^2 , current conservation requires R to be proportional to Q^2 . However, in the data from the present experiment this behavior is not yet observed, and R remains nearly constant over the measured range in Q^2 . Thus, the transition to $R \propto Q^2$

TABLE I. The values of R_H and $R^D - R^H$ calculated via the Rosenbluth separation. Note that the systematic error in the difference accounts for the correlation between the uncertainties in the hydrogen and deuterium data.

Q^2 (GeV 2)	x	R^H	Stat.	Syst.	$R^D - R^H$	Stat.	Syst.
0.150	0.041	0.259	0.074	0.153	0.036	0.131	0.136
0.175	0.050	0.307	0.056	0.188	-0.100	0.091	0.196
0.273	0.077	0.460	0.049	0.132	-0.162	0.084	0.153
0.283	0.081	0.414	0.045	0.117	-0.141	0.071	0.138
0.476	0.156	0.283	0.063	0.025	-0.115	0.091	0.061
0.508	0.091	0.406	0.038	0.168	-0.164	0.065	0.172
1.045	0.200	0.335	0.048	0.041	-0.155	0.068	0.046
1.670	0.320	0.211	0.038	0.021	-0.040	0.057	0.051

must occur below a Q^2 value of about 0.1 GeV 2 at low x , or below 1 GeV 2 for $x > 0.2$. This result also disagrees with the model calculation shown in Fig. 1 and might indicate a larger higher twist contribution in the longitudinal structure function than assumed in that model.

Both the Rosenbluth separation and the model dependent extraction of R were also carried out for the deuterium data. While the precision of the results from the Rosenbluth separation is comparable to that of the hydrogen data, the systematic uncertainty in the model dependent extraction is much bigger for the deuterium data due to a large uncertainty in the calculation of the quasielastic radiation tail which is significant at low x and Q^2 .

Thus, the difference $R^D - R^H$ was calculated using only the results from the Rosenbluth separation method, compiled in Table I, and compared in Fig. 2 to previous results from the NMC Collaboration [9] and SLAC [10]. In this plot only the most recent and precise data from SLAC [10] are shown, while additional SLAC results [3] are included in the statistical analysis of $R^D - R^H$, below. Previously, the conclusion was drawn that there is no difference between R^D and R^H . However, most of the data from NMC are at rather high Q^2 values, where R itself is small, and the $R^D - R^H$ values extracted from the SLAC measurements [3] were averaged over all Q^2 , including high Q^2 values, and were hence biased toward smaller differences. Including our results, the data are still consistent with R^D being identical to R^H . However, at values of $Q^2 < 1.5$ GeV 2 there is a hint both in the present data and in the highest-precision data from SLAC [10] that R^D is smaller than R^H . The global average (including all data) yields $R^D - R^H = -0.054 \pm 0.029$.

The results presented here are measurements of the longitudinal to transverse cross section ratio below Q^2 of about 2 GeV 2 for hydrogen and deuterium targets. These data appear in a region where R might be expected to disappear as Q^2 gets very small. However, a nearly constant behavior of R^H and R^D is observed down to Q^2 of about 0.1 GeV 2 at low values of x . For $Q^2 < 1.5$ GeV 2 , the data hint at a small difference between R^D and R^H .

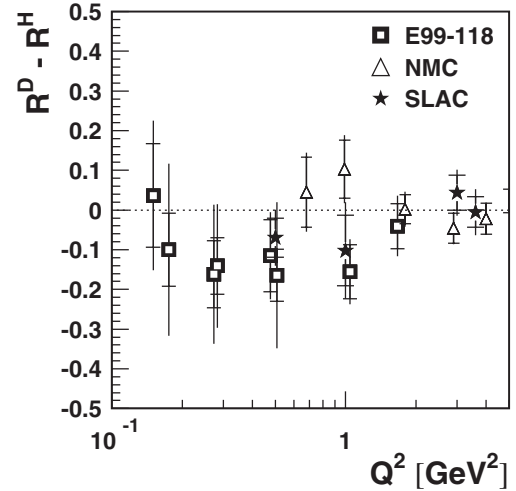


FIG. 2. The difference $R^D - R^H$ as a function of Q^2 from the present experiment calculated via Rosenbluth separation. The data from previous experiments are also shown for $Q^2 < 5$ GeV 2 .

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