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### Publication Date

1966-09-21

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Submitted to Physical Review

UCRL-16648 Rev.  
Preprint

UNIVERSITY OF CALIFORNIA  
Lawrence Radiation Laboratory  
Berkeley, California

AEC Contract No. W-7405-eng-48

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THE REACTION  $K^-p \rightarrow \Lambda\omega$  FROM 1.2 TO 2.7 BeV/c:  
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ABSTRACT

We have analyzed over 9000  $K^-p \rightarrow \Lambda\omega \rightarrow (p\pi^-) (\pi^+\pi^-\pi^0)$  events in four momentum regions between 1.2 and 2.7 BeV/c. We have systematically determined the differential cross section and the eleven independent decay-correlation parameters as a function of production angle for each of the four momentum regions. A striking forward peak in the differential cross section at our highest momentum, 2.6 BeV/c, suggests the appearance of strange-meson exchange. Using a new formalism for the absorption model, we show that the behavior of the differential cross section and the decay-correlation parameters at 2.6 BeV/c as a function of production angle is qualitatively explained by the absorption model with K and  $K^*$  exchange. Using available data on  $K^-p \rightarrow \Lambda\phi$  at 2.6 BeV/c, we show that the absorption model also explains the behavior of  $K^-p \rightarrow \Lambda\phi$ , and that the comparison between the couplings of  $K^-p \rightarrow \Lambda\omega$  and  $K^-p \rightarrow \Lambda\phi$  is in reasonable agreement with SU(3) predictions.

## I. INTRODUCTION

### A. Historical Note

In quantum electrodynamics both photons and charged particles provide the forces of interaction; in strong interactions, therefore, it is thought that the forces are created by the exchange of mesons and baryons, the known strongly interacting particles.

The attempt to put this philosophy into practice has led physicists down many paths; simple field theory, a direct translation from q. e. ; S-matrix theory and bootstrap dynamics,<sup>1</sup> based on the analyticity requirements of amplitudes; Regge poles,<sup>2</sup> involving analytic continuation in angular momentum variables; and lately, the absorption model,<sup>3</sup> a more specialized approach with limited application. In all these approaches, exchanges may take place in either the direct channel (resonances) or in crossed channels (meson or baryon exchanges). The absorption model, which is our concern in the body of this paper, deals most successfully with meson exchanges.

The absorption model has its motivation in the simple Feynman diagram of Fig. 1a. The basic contribution of the model is the addition of the diagrams of Fig. 1b involving elastic scattering in the initial and final state. Although the foregoing explanation appears to put the absorption model squarely under field theory, Ball and Frazer<sup>4</sup> have used S-matrix language to justify the basic equations of the model, at least for pseudoscalar exchange. The model has been applied with reasonable success<sup>5</sup> to  $\pi^- p \rightarrow \rho^- p$ ,  $Kp \rightarrow K^* p$ , and many other reactions involving pion and nonstrange vector-meson exchange, as well as  $K^- p \rightarrow \pi^- Y_1^*$ , involving  $K^*$  exchange.<sup>5</sup> Here we apply the absorption model to  $K^- p \rightarrow \Lambda\omega$  and  $K^- p \rightarrow \Lambda\phi$ , which involve  $K$  and  $K^*$  exchange.

### B. Summary of Results

We have analyzed over 9000  $K^-p \rightarrow \Lambda\omega \rightarrow (p\pi^-) (\pi^+\pi^-\pi^0)$  events in four momentum regions between 1.2 and 2.7 BeV/c. We have systematically determined the differential cross section and the eleven decay-correlation parameters as a function of production angle for each of the four momentum regions. In this section of the Introduction, we will indicate our line of thinking as to the implications of our results.

Figure 2 shows the total cross section for  $K^-p \rightarrow \Lambda\omega$  as a function of beam momentum. We note that there are no striking resonance phenomena; the cross section rises from threshold and falls smoothly in the usual manner for inelastic reactions, at least within our statistics. There is a known resonance,  $Y_0^*(2100)$ , with correct quantum numbers for decay into  $\Lambda\omega$ . A laboratory  $K^-$  momentum of 1.7 BeV/c corresponds to a center-of-mass energy of 2100 MeV. We can set an upper limit for the branching ratio into  $\Lambda\omega$ :

$$\frac{Y_0^*(2100) \rightarrow \Lambda\omega}{Y_0^*(2100) \rightarrow \text{all}} < 0.1 .$$

Now it is perfectly possible that there are other resonances in this region, and it is even possible that an extremely careful analysis of the data could give some indications of them, but the separation of the data into smaller energy intervals would reduce the accuracy of the measurements, because of poor statistics, to such an extent that conclusions could be drawn only with difficulty.

If we focus our attention on the production-angle distributions, Figures 3 through 6, we see that only very low partial waves are needed to explain the 1.5-, 1.7-, and 2.1-BeV/c data, but a very striking forward peak appears at 2.6 BeV/c. This peak could be caused by the interference of high-partial-wave amplitudes coming from some direct-channel activity (see Fig. 7a); we prefer to interpret it as most forward

peaks in interactions around this energy have been interpreted--as the effect of poles in the crossed-channel (see Fig. 7b); in our case, strange-meson exchange.

A broad peak in the differential cross section also appears near the backward direction at 2.6 BeV/c (Fig. 6). Fried and Taylor have interpreted similar data at 3 BeV/c as a manifestation of nucleon exchange.<sup>6</sup> While this explanation is possible, the smooth variation of the production-angle distribution from 1.5 through 2.1 BeV/c makes it seem similarly plausible that the hump at 2.6 BeV/c (and, presumably, the one at 3 BeV/c) is simply a continuation of low partial-wave behavior associated with threshold and resonance effects.

Hence we systematically present our data at all momenta, believing that the data represent the effects of threshold and perhaps some resonance behavior, except for the striking forward peak at 2.6 BeV/c, which we associate with strange-meson exchange.

Before we consider the absorption model, we should discuss why we did not apply any Regge-pole analysis to our data. Briefly, our data are at too low an energy. The requirement that a Regge-pole approximation be valid is usually expressed in terms of  $\cos\theta_t$ , where  $\theta_t$  is the "production angle" in the crossed channel. Since  $\theta_t$  is an unphysical angle,  $|\cos\theta_t|$  is greater than 1; the validity criterion is  $|\cos\theta_t| \gg 1$ . (At least, plead the advocates of Regge poles, have  $|\cos\theta_t| \gtrsim 5$ .) At 1.5 BeV/c in  $K^-p \rightarrow \Lambda\omega$ , we have  $|\cos\theta_t|$  between 1.0 and 1.5; at 2.6 BeV/c, we almost, but not quite, reach  $|\cos\theta_t| = 3$ . Hence it would have very little meaning to apply Regge poles at our energies. It is the absorption model that has had success at these energies.

We use a new formalism for the absorption model developed by R. Huff,<sup>7</sup> in which a linear-momentum representation is used instead of the usual angular-momentum representation involving partial-wave decomposition. We show that;

1. The absorption model has excellent success in fitting the differential cross section and qualitative success in fitting the



decay parameters of  $K^-p \rightarrow \Lambda\omega$  at 2.6 BeV/c in the forward direction.

2. Where it is not applicable, namely the lower momentum regions, the absorption model fails to give reasonable fits.

3. The K-meson-exchange coupling determined in an unconstrained variation of parameters is in remarkable agreement with the SU(3) prediction, and the  $K^*$ -exchange couplings are of a reasonable order of magnitude.

4. The reaction  $K^-p \rightarrow \Lambda\phi$  at 2.6 BeV/c in the forward direction is also reasonably well explained by the absorption model, and the comparison of the K- and  $K^*$ -exchange couplings determined for  $K^-p \rightarrow \Lambda\phi$  with those determined for  $K^-p \rightarrow \Lambda\omega$  is in agreement with SU(3).

## II. ANALYSIS OF THE DATA

### A. Introduction

Approximately 9400 events of the reaction  $K^- p \rightarrow \Lambda \omega \rightarrow (p\pi^-)\pi^+\pi^-\pi^0$  have been identified in a  $K^-$  exposure of the 72-in. hydrogen bubble chamber. The momentum settings ranged from 1.2 to 2.7 BeV/c. Figure 8 shows the beam-momentum spectrum for 32 000 events of the type  $K^- p \rightarrow \Lambda \pi^+\pi^-\pi^0$ . Since the cross section for the reaction  $K^- p \rightarrow \Lambda \pi^+\pi^-\pi^0$  is changing in this energy region, Fig. 8 does not reflect the relative amount of film taken at the various momenta. Table I summarizes the data taken at each momentum setting in terms of the number of events per millibarn of cross section.

The bubble chamber was exposed in two different runs, with the use of two entirely different beam configurations.<sup>8,9</sup> The method for identifying the desired events in the first run, designated K72 and with beam momenta from 1.2 to 1.7 BeV/c, has already been given in a previous publication.<sup>10</sup> The analysis of the second run, designated K63 and with beam momenta from 1.7 to 2.7 BeV/c, is given in detail here. For the reader's convenience we include the important parameters of the first run where they are of interest.

### B. Scanning and Measuring

The film was scanned once and the events found were measured. All V + two-prong events were fit to the following hypotheses:

$$K^- p \rightarrow \Lambda \pi^+ \pi^- \quad (1)$$

$$\Lambda \pi^+ \pi^- \pi^0 \quad (2)$$

$$\Lambda K^+ K^- \quad (3)$$

$$\Sigma^0 \pi^+ \pi^- \quad (4)$$

$$\Sigma^0 K^+ K^- \quad (5)$$

$$p \pi^- \bar{K}^0 \quad (6)$$

$$p \pi^- \pi^0 \bar{K}^0 \quad (7)$$

$$n \pi^+ \pi^- \bar{K}^0, \quad (8)$$

where the  $\Sigma^0$  always decays into  $\Lambda\gamma$ ,  $\Lambda$  decays into  $p\pi^-$ , and  $\bar{K}^0$  decays into  $\pi^+\pi^-$ .

V + two-prongs are fitted to hypotheses (1) through (8) in two steps. First, the neutral V direction is taken to be the line connecting the primary vertex to the vertex of the V, and the V is fit to two hypotheses,  $\Lambda \rightarrow p\pi^-$  and  $\bar{K}^0 \rightarrow \pi^+\pi^-$ . These are three-constraint fits. For  $\chi^2(\Lambda) < 32$ , reaction hypotheses (1) through (5) are tried; for  $\chi^2(\bar{K}^0) < 32$ , reaction hypotheses (6) through (8) are tried. For  $\chi^2(\Lambda)$  and  $\chi^2(\bar{K}^0)$  each less than 32, all production hypotheses are tried; in this case if an acceptable  $\chi^2$  is obtained from some production process for both interpretations of the V, the event is classified as ambiguous between  $\Lambda$  and  $\bar{K}^0$  production. The percentage of ambiguous events varied from 2.2% at 1.7 BeV/c to 6.7% at 2.6 BeV/c. (In K72 the percentage varied from 1.2% at 1.2 BeV/c to 2.5% at 1.7 BeV/c. The two independently analyzed samples at 1.7 BeV/c thus agree.) Most of the ambiguous events are  $\Lambda$  events.<sup>10</sup>

We must now consider how to separate type-2 events from those of types 1, 3, 4, 5, and 6. Events that simultaneously fit reactions (2) and (1) or (3) constitute less than 2% of the sample which fits (2). Consequently the  $\Lambda\pi^+\pi^-\pi^0$  events are free from contamination by  $\Lambda\pi^+\pi^-$  or  $\Lambda K^+K^-$  events. The task of separating the other reactions is not so simple. Since the  $\gamma$  ray and the  $\Lambda$  of (4) and (5) are constrained to have the  $\Sigma^0$  mass, reactions (4) and (5) are two-constraint fits, while reaction (2) is a one-constraint fit. If our measurement errors were properly estimated and were free from systematic errors, the mean value of  $\chi^2$  (production), for events that are truly of the type being fitted, would be equal to the number of the constraint class. Actually our errors are underestimated, so that this equality does not hold in general. Nevertheless, a confidence level is calculated for each hypothesis. Events are accepted as being a particular reaction if the confidence level for that reaction is greater than the confidence levels of all other hypotheses and the confidence level is greater than 0.005. If all confidence levels are less than 0.005, the event is classified as a failure.

The failing rate for first measurements is not small (between 30 and 40%) and therefore events that have failed are measured a second time, and sometimes even a third time. Both second and third measurements have about a 50% failing rate.

### C. Scanning and Measuring Biases

We must now consider the possibility that the loss of events due to scanning and measuring errors has biased the angular distributions in which we are interested.

#### 1. Scanning Biases

We have checked for two possible scanning biases.

a. Opening angle of the  $\Lambda$ . The direction of the pion in the  $\Lambda$  rest frame makes an angle  $\psi$  with respect to the direction of the  $\Lambda$  in the laboratory. (The  $\Lambda$  laboratory direction remains the same when transformed to the  $\Lambda$  rest frame.) If the scanning contains no biases against certain opening angles, then the distribution of  $\cos\psi$  should be flat. Figure 9 illustrates the distribution of  $\cos\psi$  for 1.7 BeV/c. The other momenta have similarly flat distributions.

b.  $\Lambda$  length cutoff. An event in which the  $\Lambda$  has decayed within a few millimeters of the production vertex is difficult to distinguish from a four-prong event. The distribution of the length of the  $\Lambda$ , shown in Fig. 10 for 1.7 BeV/c, deviates from the expected approximate exponential at 2 or 3 mm and less. To check whether this causes a bias in angular distributions for  $K^-p \rightarrow \Lambda\omega$  we have compared the center-of-mass (c. m.) production-angle distribution for events with  $750 \text{ MeV} < M(\pi^+\pi^-\pi^0) < 810 \text{ MeV}$  at 2.6 BeV/c with a production-angle distribution obtained from the same events in the following way: All events whose  $\Lambda$  went less than 2 mm in the laboratory were discarded, and each remaining event was weighted by the factor  $\exp[x/\eta c\tau]$ , where  $x$  is the  $\Lambda$  length cutoff (2 mm),  $\eta$  is the laboratory momentum of the  $\Lambda$  in BeV/c divided by the mass of the  $\Lambda$  in  $\text{BeV}/c^2$  and  $\tau$  is the mean life of the  $\Lambda$ . It should be noted that at 2.6 BeV/c and below, the  $\Lambda$  is constrained to the forward 45-deg cone

in the laboratory. We have chosen 2.6 BeV/c as our sample because two reasons indicate that the bias should be worst at the highest momentum; first, the  $\Lambda$  can go slowest in the laboratory, and second, the correction is largest for  $\Lambda$ 's that go backwards in the center of mass and the 2.6-BeV/c production-angle distribution is sharply peaked in the backward direction (forward direction for the three-pion system). Figure 11 shows the unweighted distribution with the weighted points shown as boxes. The corrections are within the error bars, and it should be remembered that when background is subtracted, the estimated errors will increase. Since the decay correlations will be much less affected by this bias than the production-angle distributions, we have not weighted events in any of our analyses of angular distributions.

No scanning biases relating to the two prongs in the V + two-prong events have been discovered.

## 2. Measuring Biases

Possible measuring biases due to the large failure rates in first and second measurements have been investigated in the following way: Angular distributions for events which passed the first measurement are compared with those that failed the first measurement but passed the second. Figure 12 shows the production-angle distributions for two such samples. No significant differences are noted. Twice-failing events have been scanned and no obvious biases were detected.

### D. Ambiguities

Among the 6300 events in the K63 run which fit  $K^-p \rightarrow \Lambda \pi^+ \pi^- \pi^0$  and have a  $M(\pi^+ \pi^- \pi^0)$  in the  $\omega$  region, there are undoubtedly a small number of events that are really of other reaction types. However, there is no reason to suppose that these events create a peak in the mass spectrum near the  $\omega$  mass, which might be confused with the  $\omega$ . Since ambiguities are known to constitute less than 10% of the  $\Lambda \pi^+ \pi^- \pi^0$  sample, the contamination of other reactions in the  $\omega$  region is less than 10% of background, and therefore is negligible. (Of course, we believe that the contamination from other reactions is much less than

this 10% ambiguity percentage because we think we have estimated confidence levels reasonably well. The upper limit considered here is nevertheless satisfying.) We have further reduced the effect of any background by the subtraction technique outlined in Section III.

### E. Total-Cross-Section Determinations

Total cross sections in the K72 run have been published,<sup>11</sup> The values are listed in Table I.

In the K63 experiment total  $K^-$  path lengths have been determined by Lindsey and Smith<sup>12</sup> at all momenta except 1.7 BeV/c. We determined the path length at 1.7 BeV/c by counting  $\tau$  decays of the  $K^-$  in the same manner as they.

If we divide the total number of good events in a certain fiducial volume that come through the system at a particular momentum setting by the path length at that momentum we will obtain a total cross section for the reaction we are studying. The number we obtain, however, needs several corrections.

#### 1. Scanning Inefficiency

We obtained the scanning efficiencies by scanning the film a second time and comparing the list of events found with the list of good events whose  $\Lambda$  has a length greater than 5 mm. (We correct for  $\Lambda$  length cutoff separately, and we would not want to do it twice.) Good events are those that were found on the first scan and that fit the hypothesis  $K^-p \rightarrow \Lambda\pi^+\pi^-\pi^0$ . Then the scanning efficiency is (number of good events found on the second scan)/(number of good events). The scanning efficiencies varied from 94 to 97%. (In K72, the scanning efficiencies varied from 94 to 98%.) Table II lists scanning efficiencies.

#### 2. Measuring Inefficiency

We calculated an effective measuring efficiency by computing passing rates (number passed/number measured) for first, second, and third measurements, and using these numbers to project the failing and unmeasured events through third measurement. Of 120 000  $V +$  two-prong events in K63, 79 000 have passing measurements, and

21 000 more would pass if we completed the measuring program through third measurements. It might appear that 20 000 events are unaccounted for. However, twice-failing events were scanned, and it was discovered that about 50% of them were not V + two-prong events. Projections show that 23 000 events should fail twice and thus we know that in our sample about 12 000 events are not V + two-prong events. We have therefore accounted for all the V + two-prong events to within 6%.

Table II lists correction factors that must be used to multiply the number of passing events to obtain the true number of events of a particular reaction.

### 3. $\Lambda$ Length Cutoff

The distribution in proper time (length/momentum) for the  $\Lambda$ 's in our sample, which we expect to be an exponential with decay corresponding to the mean life of the  $\Lambda$ , is seen to drop in the region of short times. We account for the missing events at both ends of the time spectrum, and find corrections of  $4 \pm 3\%$  at 1.7 BeV/c,  $5 \pm 3\%$  at 2.1 BeV/c and  $5 \pm 3\%$  at 2.6 BeV/c.

### 4. Dalitz Decay of the $\pi^0$

The  $\pi^0$  from  $\omega$  decay will give a Dalitz pair at the production vertex 1.25% of the time. The event would then be a V + four-prong event and hence would be lost to our V + two-prong sample. We must increase each cross section by 1.25% to account for this effect.

Total cross sections are given in Table I and in Fig. 2.

After considerable analysis we have obtained an unbiased group of events of the type  $K^- p \rightarrow \Lambda \omega \rightarrow \Lambda \pi^+ \pi^- \pi^0$ , and the normalization needed to obtain total cross sections is well understood.

### III. RESULTS OF THE ANALYSIS

#### A. Introduction

We are considering  $K^- p \rightarrow \Lambda \omega$  events in which the  $\omega$  decays into  $\pi^+ \pi^- \pi^0$  and the  $\Lambda$  decays into  $p \pi^-$ . There is an extremely large amount of data contained in every event of this type; our problem is to present the data in a useful and understandable form.

First we define the variables (vectors, such as the decay pion momentum from the  $\Lambda$ ; and scalars, such as the c. m. energy) which characterize each event and which can vary from event to event. The differential cross section can then be expressed as a function of these variables in a simple way. The parameters of this function express concisely our knowledge of the reaction. (For example, one may express the knowledge of an angular distribution by giving only the coefficients of the Legendre polynomials in the expression for the angular distribution.)

#### B. Definition of Internal Variables

In the c. m. system, illustrated in Fig. 13, we use as variables the c. m. energy  $E$  and the production angle  $\theta$ , defined by

$$\cos \theta = \underline{\underline{K}} \cdot \underline{\underline{\omega}} / |\underline{\underline{K}}| |\underline{\underline{\omega}}|.$$

We obtain all rest-frame quantities by first transforming to the center-of-mass system and then to the rest frame in question.

Unit vectors defined in the  $\omega$  rest frame are:

$\underline{\underline{n}}$  = normal to the plane of the pions from the  $\omega$  decay ( $\underline{\underline{\pi}}^- \times \underline{\underline{\pi}}^+$ ).

$\underline{\underline{N}}$  = normal to the production plane  $\underline{\underline{K}} \times \underline{\underline{\omega}}$  (defined in the c. m. frame and unchanged when shifted to the  $\omega$  rest frame).

$\underline{\underline{X}}$ ,  $\underline{\underline{Y}}$ , and  $\underline{\underline{N}}$  = an orthogonal set of axes defined by the production process (e. g.,  $\underline{\underline{P}}_1, \underline{\underline{N}} \times \underline{\underline{P}}_1, \underline{\underline{N}}$ ).

Unit vectors defined in the  $\Lambda$  rest frame are:

$\underline{\underline{\pi}}$  = direction of the pion from the  $\Lambda$  decay.

$\underline{\underline{N}}$  = normal to the production plane (same as in  $\omega$  rest frame).

$\underline{\underline{X}}'$ ,  $\underline{\underline{Y}}'$ , and  $\underline{\underline{N}}$  = an orthogonal set of axes defined by the production process (e. g.,  $\underline{\underline{q}}_1, \underline{\underline{N}} \times \underline{\underline{q}}_1, \underline{\underline{N}}$ ).



### C. Expressions for Cross Sections

Byers and Yang<sup>13</sup> and Berman and Oakes<sup>14</sup> have exhibited the general dependence of this reaction on the angles formed in the decays of the final-state particles, given an unpolarized target. Huff<sup>15</sup> has also discussed this reaction. Ademollo and Gatto<sup>16</sup> treated the production characteristics of reactions of this type by means of a density-matrix formalism; such a treatment is the connecting link between the correlations in this section and the production amplitudes. Of course, the spins and parities of the  $\omega$  and  $\Lambda$  are taken to be  $1^-$  and  $1/2^+$ , respectively. We may express the entire dependence of the cross section on internal variables as

$$d^5\sigma = \left[ F_1(\underline{n}\cdot\underline{N})^2 + F_2(\underline{n}\cdot\underline{X})^2 + F_3(\underline{n}\cdot\underline{Y})^2 + F_4(\underline{n}\cdot\underline{X})(\underline{n}\cdot\underline{Y}) + F_5(\underline{n}\cdot\underline{N})^2(\underline{\pi}\cdot\underline{N}) \right. \\ \left. + F_6(\underline{n}\cdot\underline{X})^2(\underline{\pi}\cdot\underline{N}) + F_7(\underline{n}\cdot\underline{Y})^2(\underline{\pi}\cdot\underline{N}) + F_8(\underline{n}\cdot\underline{X})(\underline{n}\cdot\underline{Y})(\underline{\pi}\cdot\underline{N}) + F_9(\underline{n}\cdot\underline{N})(\underline{n}\cdot\underline{X})(\underline{\pi}\cdot\underline{X}') \right. \\ \left. + F_{10}(\underline{n}\cdot\underline{N})(\underline{n}\cdot\underline{Y})(\underline{\pi}\cdot\underline{X}') + F_{11}(\underline{n}\cdot\underline{N})(\underline{n}\cdot\underline{X})(\underline{\pi}\cdot\underline{Y}') + F_{12}(\underline{n}\cdot\underline{N})(\underline{n}\cdot\underline{Y})(\underline{\pi}\cdot\underline{Y}') \right] \\ \cdot \left[ \frac{3}{(4\pi)^2} d\Omega_{\underline{\pi}} d\Omega_{\underline{n}} d\cos\theta \right].$$

Each  $F_i$  is an unknown function of  $E$  and  $\cos\theta$ , and depends on the dynamics of the process.

It is convenient to introduce another parameterization of the cross-section formula:

$$d^5\sigma = C(E, \cos\theta) \left[ f_1(\underline{n}\cdot\underline{N})^2 + f_2(\underline{n}\cdot\underline{X})^2 + f_3(\underline{n}\cdot\underline{Y})^2 + f_4(\underline{n}\cdot\underline{X})(\underline{n}\cdot\underline{Y}) \right. \\ \left. + f_5(\underline{n}\cdot\underline{N})^2(\underline{\pi}\cdot\underline{N}) + \dots \right] \left[ \frac{3}{(4\pi)^2} d\Omega_{\underline{\pi}} d\Omega_{\underline{n}} d\cos\theta \right],$$

with the subsidiary condition  $f_1 + f_2 + f_3 = 1$ , which is the normalization condition after integration over the two solid angles involved. By this parameterization we have provided a convenient normalization for the dependence of the cross section on the decay angles of the  $\Lambda$  and  $\omega$ .

That is, the dependence of the cross section on the decay angles (which

means the dependence on the spin alignments of the  $\Lambda$  and  $\omega$ ) is contained in the  $\{f_i\}$  in the form of a probability density whose integral is 1. Thus we have

$$\begin{aligned} d\sigma &= \int_{\Omega_\pi} \int_{\Omega_n} d^5\sigma \\ &= C(E, \cos\theta) d\cos\theta \iiint \left[ f_1 (\underline{n} \cdot \underline{N})^2 + \dots \right] \frac{3}{(4\pi)^2} d\Omega_\pi d\Omega_n \\ &= C(E, \cos\theta) d\cos\theta. \end{aligned}$$

Thus  $C(E, \cos\theta)$  is the differential cross section, integrated over all decay angles, of the reaction taking place at a given  $E$  and  $\cos\theta$ .

The total cross section is given by

$$\sigma_T = \int C(E, \cos\theta) d\cos\theta.$$

#### D. General Model

At this point we might tabulate  $d\sigma/d\cos\theta$  and the set of  $f_i$  as a function of  $E$  and  $\cos\theta$ . However, we still face the problem of choosing the vectors  $\underline{X}'$ ,  $\underline{Y}'$  and  $\underline{X}$ ,  $\underline{Y}$  in the  $\Lambda$  and the  $\omega$  rest frames, respectively. If we could do our experiment at a unique  $E$  and a unique  $\theta$ , then, in each frame, any choice would be related to any other by a simple rotation around the normal. However, since we must average over regions of  $E$  and  $\cos\theta$ , it behooves us to choose our axes carefully. The choice is determined by the characteristics of the model being tested.

Most current theories have as a basis the idea of exchanged particles, as expressed, for example, in Feynman diagrams or unitarity graphs. Figures 7a, b, and c represent exchanges of the least-massive particles allowed in the three possible channels in  $Kp \rightarrow \Lambda\omega$  -- the s, t, and u channels. With this model the correct choice of axes is apparent. In the u channel the appropriate axes in the  $\omega$  rest frame are  $\underline{P}_1$ ,  $\underline{N} \times \underline{P}_1$ , and  $\underline{N}$ , and in the  $\Lambda$  rest frame they are  $\underline{q}_1$ ,  $\underline{N} \times \underline{q}_1$ ,

and  $\underline{N}$ . In the  $t$  channel, the two sets are  $\underline{q}_1$ ,  $\underline{N} \times \underline{q}_1$ , and  $\underline{N}$ , and  $\underline{P}_1$ ,  $\underline{N} \times \underline{P}_1$ , and  $\underline{N}$ . In the  $s$  channel we have  $\underline{P}_2$ ,  $\underline{N} \times \underline{P}_2$ , and  $\underline{N}$  and  $\underline{q}_2$ ,  $\underline{N} \times \underline{q}_2$ , and  $\underline{N}$ . In this article we concern ourselves, in the section on the absorption model, with  $t$ -channel exchanges of pseudoscalar and vector mesons. We therefore tabulate our data with the  $\{f_i\}$  determined with axes appropriate to the  $t$  channel. We iterate that the  $\{f_i\}$ , if they were obtained at a unique  $E$  and  $\theta$ , would be related to the  $u$  and  $s$  channel  $\{f_i\}$  by a simple rotation.

### E. Experimental Calculations

The quantities  $d\sigma/d\cos\theta$  and  $\sigma_T$  were obtained by a simple counting of events in a given region of  $E$  and  $\cos\theta$ . The only problem here is background subtraction, which is discussed in the next section.

The maximum-likelihood technique was used to determine  $\{f_i\}$ . For each event we have a probability density that is a function of the twelve  $f_i$ ,

$$P_k = f_1(\underline{n} \cdot \underline{N})_{\underline{k}}^2 + f_2(\underline{n} \cdot \underline{X})_{\underline{k}}^2 + f_3(\underline{n} \cdot \underline{Y})_{\underline{k}}^2 + \dots,$$

where the vectors have been evaluated for the particular event, as the subscript  $k$  indicates.

For a sample of  $N$  events, the likelihood is,

$$\mathcal{L} = \prod_{k=1}^N P_k$$

We maximize  $\mathcal{L}$  by maximizing

$$\ln \mathcal{L} = \sum_{k=1}^N \ln P_k.$$

We vary only eleven of the parameters  $\{f_i\}$ , since there is one constraint. Only one extremum can exist for our likelihood, and it is a maximum. Both of these facts are a consequence of the linearity of  $P_k$  as a function of the parameters  $\{f_i\}$ .

### F. Background

If we look at the three-pion mass spectra in the reaction  $Kp \rightarrow \Lambda \pi^+ \pi^- \pi^0$ , we see a prominent  $\omega$  peak (see Figs. 14 through 17). Under this peak we also see a significant background, which we judged from the regions adjacent to the peak. By sketching a curve through the regions next to the  $\omega$  peak, we estimate the number of non- $\omega$  events in the region of the three-pion mass between 750 and 810 MeV. We assume that the remainder in this region represent events of the reaction  $Kp \rightarrow \Lambda \omega$ .

Let us call the 750- to 810-MeV region the  $\omega$  region, and the two regions 690 to 750 MeV and 810 to 870 MeV, combined, the control region. Let  $N_B$  be the number of background events in the  $\omega$  region, and  $N_C$  the number of events in the control region. We are dealing with a spectrum at a given  $E$ . To find the number of  $\omega$  events  $N_\omega$  in a certain region of  $\cos \theta$ , we use  $N_\omega = N - (N_B/N_C)(M)$ , where  $N$  is the number of events in the  $\omega$  region and in the region of  $\cos \theta$  under discussion, and  $M$  is the number of events in the control region and in the region of  $\cos \theta$  under discussion.

Treating background in determinations of  $\{f_i\}$  is only slightly more complicated. The  $\{f_i\}_\omega$  for events in the  $\omega$  region is determined with the technique described in Sec. E, and another set  $\{f_i\}_C$  is determined for the events in the control region. Both sets are normalized to a total integral of one, so that the expression for the  $\{f_i\}$  for the  $\omega$  events is

$$f_i = \frac{1}{N_\omega} \left\{ N f_{i\omega} - (N_B/N_C)(M) f_{iC} \right\}.$$

### G. Errors

The errors on cross sections are treated in the usual manner for counting experiments. The errors on the  $f_i$  are more complex. The maximum-likelihood routine we used yields an error matrix (obtained from inverting the second-derivative matrix) for the eleven

$f_1$  that were varied in the search. Thus we have all the correlated errors, and since the twelfth parameter is a function of the other eleven ( $f_1 = 1 - f_2 - f_3$ ), we may find its error correlations also. When we list the error,  $\sigma_{f_i}$ , for an  $f_i$ , we are listing the square root of the diagonal element of the error matrix corresponding to that  $f_i$ . Thus the error matrix is

$$\langle \delta f_i \delta f_j \rangle = \frac{1}{(N_\omega)^2} \left[ N^2 \langle \delta f_{i\omega} \delta f_{j\omega} \rangle - \left( \frac{N_B M}{N_C} \right)^2 \langle \delta f_{ic} \delta f_{jc} \rangle \right]$$

and

$$\sigma_{f_i} = \left( \langle (\delta f_i)^2 \rangle \right)^{1/2}.$$

Because of space limitations we have provided only the error matrices for the forward direction at 2.6 BeV/c.

#### H. Presentation of Data

Figure 8 shows that our data lie in four distinct regions of c. m. energy. The exposures at 1.95 and 2.41 BeV do not comprise enough data to meaningfully determine the many parameters of the angular distributions. Therefore we have separated the data into four sections corresponding to the beam momentum settings 1.2 to 1.5 BeV/c, 1.6 to 1.7 BeV/c, 2.1 BeV/c, and 2.4 to 2.7 BeV/c. Figures 3 through 6 show the distributions in production angle for each of the four regions. Figures 18 through 21 as well as Tables III through VIII give the decay correlations  $\{f_i\}$  determined in many intervals of production angle for each of the four regions. Thus Figs. 3 through 6 and 18 through 21, along with the total cross sections shown in Fig. 2, present the entire range of knowledge available about this reaction in our experiment, and in fact represent the entire extent of the information obtainable about the production mechanisms in this reaction from film of a bubble chamber with unpolarized protons.

## IV. THEORY OF THE ABSORPTION MODEL

### A. Introduction

Reactions involving two particles in the initial state and two particles in the final state generally show a peaking at small momentum transfers, or equivalently, at forward production angles, at least at energies large enough to avoid threshold effects. The characteristic dominance of small production angles has been explained on the basis of long-range forces--the one-particle-exchange model.<sup>17</sup> (See Fig. 1a.) However, the quantitative calculation of the appropriate Feynman diagrams generally results in a production-angle distribution that is not as forward-peaked as the data and in a cross section that is larger by an order of magnitude than the data. One can say equivalently that the theoretical predictions with low partial waves removed would fit the data.

A natural explanation for a dearth of low partial waves is absorption. That is, more complicated reactions go through small impact parameters and thus compete with the two-body final state in low partial waves. This competition effectively reduces the low partial-wave components of the two-body final state. The absorption model is a quantitative treatment of the foregoing idea.<sup>3</sup>

### B. Formalism

We use a formalism developed by Huff,<sup>7</sup> which uses a linear momentum representation rather than the more usual impact-parameter (angular momentum) representation. Since Huff's results have not been published, we briefly outline his methods and equations.

First we must find the amplitudes for the Feynman diagrams corresponding to Fig. 1a with particle  $e$  being either a  $K$  or a  $K^*$  meson. These amplitudes are called the Born amplitudes. Let  $B_{ij}$  be the Born amplitude taking into account both  $K$  and  $K^*$  exchange where the initial proton has helicity plus, the final  $\omega$  has helicity  $i$ , and the final  $\Lambda$  has helicity  $j$ . The amplitudes for an initial proton with

helicity minus are related to the  $B_{ij}$  by parity conservation. Then the six independent Born amplitudes are given by:

$$B_{++} = G_1 \{(C_-/\sqrt{2}) p_2 m_3 \sin \theta\} \\ + G_2 \{-(C_+/\sqrt{2}) Q_2 - \sqrt{2} D_+ Q_1 - \sqrt{2} D_- E_{c.m.} p_2 p_4 \sin \theta [(m_2 + m_4)/a]\} \\ + G_3 \{\sqrt{2} D_- E_{c.m.} p_2 p_4 \sin \theta (m_2 + m_4)^{-1}\}$$

$$B_{0+} = G_1 \{-C_- Q_2\} + G_2 \{-C_+ p_2 m_3 \sin \theta\}$$

$$B_{-+} = G_1 \{-(C_-/\sqrt{2}) p_2 m_3 \sin \theta\} \\ + G_2 \{(C_+/\sqrt{2}) Q_2 - \sqrt{2} D_+ Q_1 - \sqrt{2} D_- E_{c.m.} p_2 p_4 \sin \theta [(m_2 + m_4)/a]\} \\ + G_3 \{\sqrt{2} D_- E_{c.m.} p_2 p_4 \sin \theta (m_2 + m_4)^{-1}\}$$

$$B_{+-} = G_1 \{(C_+/\sqrt{2}) p_2 m_3 \sin \theta\} \\ + G_2 \{(C_-/\sqrt{2}) Q_2 - \sqrt{2} D_- Q_1 + \sqrt{2} D_+ E_{c.m.} p_2 p_4 \sin \theta [(m_2 + m_4)/a]\} \\ + G_3 \{-\sqrt{2} D_+ E_{c.m.} p_2 p_4 \sin \theta (m_2 + m_4)^{-1}\}$$

$$B_{0-} = G_1 \{-C_+ Q_2\} + G_2 \{C_- p_2 m_3 \sin \theta\}$$

$$B_{--} = G_1 \{-(C_+/\sqrt{2}) p_2 m_3 \sin \theta\} \\ + G_2 \{-(C_-/\sqrt{2}) Q_2 - \sqrt{2} D_- Q_1 + \sqrt{2} D_+ E_{c.m.} p_2 p_4 \sin \theta [(m_2 + m_4)/a]\} \\ + G_3 \{-\sqrt{2} D_+ E_{c.m.} p_2 p_4 \sin \theta (m_2 + m_4)^{-1}\},$$

where:

$$G_1 = \frac{G(\omega K^+ K^-) G(\bar{p} K^+ \Lambda)}{4\pi(m_K^2 - t)} \quad (2\beta)$$

$$G_2 = \frac{G(\omega K^+ K^{*-}) [G_V(\bar{p} K^{*+} \Lambda) + G_T(\bar{p} K^{*+} \Lambda)]}{4\pi(m_{K^*}^2 - t)} \quad \beta$$

$$G_3 = \frac{G(\omega K^+ K^{*-}) G_T(\bar{p} K^{*+} \Lambda)}{4\pi(m_{K^*}^2 - t)} \quad \beta$$

$$Q_1 = \frac{1}{2} [E_{c.m.} (E_2 + E_4) - \frac{1}{2} a + 2 m_2 m_4] \\ - [(m_1 + m_3)/a] E_{c.m.} (m_2 E_4 + m_4 E_2)$$

$$Q_2 = E_1 p_4 - E_3 p_2 \cos \theta$$

$$a = 2 E_2 E_4 + 2 m_2 m_4 - 2 p_2 p_4 \cos \theta$$

$$C_{\pm} = (1 \mp \cos \theta)^{1/2} \{ [p_4/(m_4 + E_4)] \pm [p_2/(m_2 + E_2)] \}$$

$$D_{\pm} = (1 \mp \cos \theta)^{1/2} \{ 1 \pm [p_2 p_4 / (m_4 + E_4)(m_2 + E_2)] \}$$

$$\beta = \left[ \frac{(E_4 + m_4)(E_2 + m_2)}{2 m_3^2} \right]^{1/2}$$

$$t = m_1^2 + m_3^2 - 2 E_1 E_3 + 2 p_1 p_3 \cos \theta$$

$$\cos \theta = \underline{p}_1 \cdot \underline{p}_2$$

$$p_i = |\underline{p}_i|$$

$$E_i = (p_i^2 + m_i^2)^{1/2}$$



Here the  $\underline{p}_i$  with  $i = 1, 2, 3, \text{ or } 4$  correspond to the c. m. momenta for the  $K^-$ ,  $p$ ,  $\omega$ , and  $\Lambda$  respectively;  $E_{\text{c. m.}}$  is the total energy in the c. m. system, and the coupling constants  $G(abc)$  are as defined by Jackson and Pilkuhn.<sup>18</sup>

The differential cross section in terms of these Born amplitudes is

$$\frac{d\sigma}{d\Omega} = \frac{1}{4(E_{\text{c. m.}})^2} \frac{q'}{q} \sum_{i,j} |B_{ij}|^2,$$

where  $q$  and  $q'$  are the c. m. momenta in the initial and final states, respectively.

We agree that the  $B_{ij}$  are not the correct amplitudes for the reaction  $K^- p \rightarrow \Lambda \omega$  even if this reaction takes place only by  $K$  and  $K^*$  exchange. The Born amplitudes must be modified by absorption.

The basic formula relied upon to correctly give the amplitude  $A_{ij}$  is, in matrix form,

$$A = S_f^{1/2} B S_i^{1/2} \quad (1)$$

which is the high-energy equivalent of the distorted-wave Born approximation.<sup>19</sup> The  $S_i$  ( $S_f$ ) is the S-matrix element for elastic scattering between the two particles in the initial (final) state. In other words, this extension of the distorted-wave Born approximation is equivalent to including in our calculations the Feynman diagrams represented in Fig. 1b. Omnes<sup>20</sup> has asserted that this equation is not valid in high-energy peripheral collisions involving low partial waves. However he admits that the general effect of the modification to the Born amplitude that this equation implies, namely depletion of low partial waves, should indeed actually appear due to absorption. In the opinion of Ball and Frazer,<sup>4</sup> this equation is fairly plausible within the S-matrix theory when the exchanged meson has spin zero. They find it impossible to justify for vector-meson exchange. They also assert, along with Omnes, that the approximations are easier to justify for high than

for low partial waves. The marked success of the equation is reason enough to try it here.

We may exhibit the matrix character of Eq. (1) by expanding the equation in either the angular- or linear-momentum representations.

(1) Angular-momentum representation. Let  $|a\rangle \equiv |i, J, M, \lambda_1, \lambda_2\rangle$  and  $|b\rangle \equiv |f, J, M, \lambda_3, \lambda_4\rangle$  be initial and final states, with angular momentum  $J, M$ , and let the helicity of particle  $i$  be  $\lambda_1$ . The remaining quantum numbers are contained in  $i, f$ :

$$\langle f J M \lambda_3 \lambda_4 | A | i J M \lambda_1 \lambda_2 \rangle = \sum_{\{\lambda_i'\}} \langle f J M \lambda_3 \lambda_4 | S_f^{1/2} | f J M \lambda_3' \lambda_4' \rangle \\ \cdot \langle f J M \lambda_3' \lambda_4' | B | i J M \lambda_1' \lambda_2' \rangle \langle i J M \lambda_1' \lambda_2' | S_i^{1/2} | i J M \lambda_1 \lambda_2 \rangle.$$

(2) Linear-momentum representation. Expanding in a linear-momentum representation, we have

$$\langle f \Omega \lambda_3 \lambda_4 | A | i 0 \lambda_1 \lambda_2 \rangle = \int d\Omega_f d\Omega_i \sum_{\{\lambda_i'\}} \langle f \Omega \lambda_3 \lambda_4 | S_f^{1/2} | f \Omega_f \lambda_3' \lambda_4' \rangle \\ \cdot \langle f \Omega_f \lambda_3' \lambda_4' | B | i \Omega_i \lambda_1' \lambda_2' \rangle \langle i \Omega_i \lambda_1' \lambda_2' | S_i^{1/2} | i 0 \lambda_1 \lambda_2 \rangle$$

where  $|i \Omega \lambda_i \lambda_i\rangle$  is the state vector with the c. m. momentum vector in direction  $\Omega$ .

The two representations are, of course, equivalent. However, significant differences arise in their application because different approximations are made. In the treatment of the  $S^{1/2}$  matrix elements, for the angular-momentum applications<sup>3</sup> it is assumed that the absorptivity is a function of the total angular momentum, whereas the more relevant variable is probably the orbital angular momentum. We do not have this problem with the linear-momentum application, but we must approximate the  $S^{1/2}$  matrix elements in another way

(see below). The angular-momentum applications approximate the partial-wave decomposition of the Born terms, while the linear-momentum application uses the exact Born terms. That is, the angular-momentum applications have approximated partial-wave summations by integrals. (Jackson, however, is now using exact partial-wave sums.<sup>21</sup>) This is completely avoided in the linear-momentum application. The effects of these approximations are discussed in Section V.

To continue with the linear-momentum representation, we make the usual simplifying assumption that  $\lambda_1' = \lambda_1$ ; that is, the elastic scattering in the initial and final states is all nonhelicity flip. Since the helicity-flip amplitudes must vanish in the forward direction, and the elastic-scattering differential cross sections extrapolate smoothly to near the optical-theorem point, this appears to be a reasonable assumption for the forward directions.

We must evaluate the matrix elements of  $S^{1/2}$ . We know that  $S = 1 - T$ , where  $T$  is the transition matrix, and the partial differential elastic cross sections are given by

$$d\sigma_{\lambda_1 \lambda_2} / d\Omega = |(2\pi/q) \langle i\Omega \lambda_1 \lambda_2 | T | i0 \lambda_1 \lambda_2 \rangle|^2.$$

Hence we have

$$\langle i\Omega \lambda_1 \lambda_2 | T | i0 \lambda_1 \lambda_2 \rangle = e^{i\psi} (q/2\pi) (d\sigma_{\lambda_1 \lambda_2} / d\Omega)^{1/2},$$

where  $\psi$  is an unknown phase that is a function of production angle. If the elastic scattering is due completely to the absorption of inelastic scattering (i. e., elastic scattering is "shadow" scattering), then  $\psi$  is 0. However, even if the elastic cross sections extrapolate exactly to the optical-theorem point,  $\psi \neq 0$  is still possible at  $\theta \neq 0$  deg. Calculations up until now have assumed  $\psi = 0$ ; however, we shall see in Section V that a nonzero value of  $\psi$  plays a crucial role in applying the absorption model to our data.

Under our assumptions, the differential elastic-scattering cross sections may be expressed by

$$\left(\frac{d\sigma}{d\Omega}\right)^{1/2} = \frac{\sigma_T^q}{4\pi} e^{-At/2},$$

where  $\sigma_T$  is the total cross section for interaction between the two particles in the initial state. An analogous formula holds for the final state.

We approximate  $S^{1/2}$  by

$$S^{1/2} = 1 - T/2.$$

This approximation is equivalent to considering at most one elastic scatter in each of the blobs in Fig. 1b.

We have now given enough information to construct the absorbed amplitude A. After properly taking into account the necessary rotations from various helicity frames to other helicity frames, Huff's final result is

$$\begin{aligned} \langle f\Omega\lambda_3\lambda_4 | A | i0\lambda_1\lambda_2 \rangle &= \langle f\Omega\lambda_3\lambda_4 | B | i0\lambda_1\lambda_2 \rangle \\ &- 1/2 \int d\Omega' \eta [\langle f\Omega'\lambda_3\lambda_4 | B | i0\lambda_1\lambda_2 \rangle \langle i\Omega\lambda_1\lambda_2 | T | i0\lambda_1\lambda_2 \rangle \\ &+ \langle f\Omega'\lambda_3\lambda_4 | T | i0\lambda_3\lambda_4 \rangle \langle f\Omega\lambda_3\lambda_4 | B | i0\lambda_1\lambda_2 \rangle], \end{aligned} \quad (2)$$

where

$$\eta = (\eta')^{2(\lambda_3 - \lambda_4)} e^{i(\lambda_1 - \lambda_2)\phi'}$$

$$\eta' = [e^{-i\phi'/2} \cos \theta'/2 \cos \theta/2 + e^{i\phi'/2} \sin \theta'/2 \sin \theta/2] / \cos \theta''/2,$$

and

$$\cos \theta'' = \cos \theta \cos \theta' + \sin \theta \sin \theta' \cos \phi'.$$

We have discarded the product term containing two T-matrix elements because it represents the Feynman diagram where one elastic scatter takes place in the initial state and one takes place in the final state. We already neglected the diagrams, presumably of the same order of magnitude, where two elastic scatters take place in the initial (or final) state, and none take place in the other state; therefore we must neglect the product term also.

Given the helicity amplitudes of Eq. (2), it is straightforward to give the theoretical values for the  $\{f_i\}$ , defined in Section III, and the differential cross section. However, to give numerical values we need the coupling constants and the elastic cross section behavior for  $\Lambda\omega$  scattering. We do not know the exact values for many of these parameters; we have therefore varied them in our application of the theory.

### C. Coupling Constants

In this subsection we present what is known, either theoretically or experimentally, about the magnitudes of the coupling constants involved in the reaction  $K^-p \rightarrow \Lambda\omega$  proceeding via K and  $K^*$  exchange.

(1)  $G(\omega K^+ K^-)$ . From the decay of  $\phi$  into  $K^+ K^-$ , we can determine the  $G(\phi K^+ K^-)$  coupling constant by using

$$\Gamma_{\phi \rightarrow K^+ K^-} = \frac{2}{3} \frac{G^2(\phi K^+ K^-)}{4\pi} \frac{p^3}{m_\phi^2}$$

Thus we have

$$\frac{G^2(\phi K^+ K^-)}{4\pi} = 1.2.$$

Then from SU(3), where  $\theta$  is the vector-meson mixing angle ( $\theta = 40$  deg), we have

$$\frac{G^2(\omega K^+ K^-)}{4\pi} = \tan^2 \theta \frac{G^2(\phi K^+ K^-)}{4\pi} = 0.8.$$

Alternatively, using the  $\rho \rightarrow \pi\pi$  decay, we find

$$\frac{G^2(\omega K^+ K^-)}{4\pi} = \frac{G^2(\rho\pi\pi)}{4\pi} \left(\frac{3}{4} \sin^2 \theta\right) \cong 0.7.$$

(2)  $G(\bar{p}K^+\Lambda)$ . Again we may use SU(3) to relate  $G(\bar{p}K^+\Lambda)$  to  $G(\bar{p}\pi^0 p)$ :

$$\frac{G^2(\bar{p}K^+\Lambda)}{4\pi} = \left(\frac{3 - 2\alpha}{\sqrt{3}}\right)^2 \frac{G^2(\bar{p}\pi^0 p)}{4\pi},$$

where  $G^2(\bar{p}\pi^0 p)/4\pi = 14.6$ , and  $\alpha$ , the fraction of the interaction going through the "d" (symmetric) coupling, is known to be  $\cong 0.75$ .<sup>22</sup> We find

$$\frac{G^2(\bar{p}K^+\Lambda)}{4\pi} \cong 10.$$

(3)  $G(\omega K^+ K^{*-})$ . Using the  $\phi \rightarrow \rho\pi$  decay and the  $\rho\pi$  model of the  $\omega \rightarrow 3\pi$  decay, we may discover an approximate value for  $G(\omega K^+ K^{*-})$  through SU(3). Let  $\epsilon = G(\phi\pi^+\rho^-)/G(\omega\pi^+\rho^-)$ . Then the allowed SU(3) couplings lead to

$$\frac{G(\omega K^+ K^{*-})}{G(\omega\pi^+\rho^-)} = 1 - \frac{3}{2} \sin^2 \theta - \frac{3}{2} \epsilon \sin \theta \cos \theta.$$

Glashow and Socolow have predicted from their nonet coupling scheme that  $\epsilon = -0.08$ .<sup>23</sup> They have calculated a phase-space factor of 17 favoring  $\phi \rightarrow \rho\pi$  over  $\omega \rightarrow \rho\pi$ ; hence the determination of  $\Gamma(\phi \rightarrow 3\pi) = 0.4 \pm 0.3$  MeV by Lindsey and Smith<sup>24</sup> leads to  $|\epsilon| = 0.05 \pm 0.03$ . We will use  $\epsilon = -0.08$  to find  $G(\omega K^+ K^{*-})$ . Thus we have

$$\frac{G(\omega K^+ K^{*-})}{G(\omega\pi^+\rho^-)} = 0.64.$$

An expression for  $G(\omega\pi^+\rho^-)$  in terms of the width of the  $\omega$  has been derived by Gell-Mann, Sharp, and Wagner.<sup>25</sup> Their expression leads to

$$\frac{G^2(\omega\pi^+\rho^-)}{4\pi} \cong 14;$$

therefore

$$\frac{G^2(\omega K^- K^{*+})}{4\pi} = 5.7.$$

(4)  $G_V(\bar{p}K^{*+}\Lambda)$  and  $G_T(\bar{p}K^{*+}\Lambda)$ . The couplings of the  $\rho$ ,  $\omega$ , and  $\phi$  vector mesons to the baryons can be deduced from nucleon-nucleon forces. However, in view of the wide variations in determinations, [e. g., Scotti and Wong<sup>26</sup> find  $G_V^2(\bar{p}\omega p)/4\pi \cong 3$ ; Bryan and Scott<sup>27</sup> find  $G_V^2(\bar{p}\omega p)/4\pi \cong 22.$ ] we probably should restrict ourselves to saying that  $G_V^2/4\pi$  and  $G_T^2/4\pi$  are  $\sim 10$ .

Cabibbo<sup>28</sup> has suggested a scheme that predicts the ratio  $G_T/G_V$ . The interaction of baryons with vector particles can be written

$$\langle \bar{B} | V | B \rangle = \text{Tr} \left[ \mathcal{O} V [\bar{B}, B] + \mathcal{E} V \{B, B\} \right]$$

where  $V$ ,  $\bar{B}$ , and  $B$  are matrices representing the vector-meson, anti-baryon, and baryon octets,  $\mathcal{O}$  has the form  $a \gamma_\mu + b \sigma_{\mu\nu} k_\nu$  and  $\mathcal{E}$  has the form  $a' \gamma_\mu + b' \sigma_{\mu\nu} k_\nu$ . If we assume that the electromagnetic current has the same transformation properties as the vector-meson octet, which is another way of saying that the photon and the vector mesons are all coupled to the same SU(3) conserved currents, we can write the electromagnetic interaction of the proton and neutron as follows:

$$\langle \bar{p} | j_{e.m.} | p \rangle \sim \mathcal{E} - 3\mathcal{O}$$

$$\langle \bar{n} | j_{e.m.} | n \rangle \sim -2\mathcal{E}.$$

But we know that

$$\langle \bar{p} | j_{e.m.} | p \rangle \sim \gamma_\mu + (\mu_p/2M) \sigma_{\mu\nu} k_\nu$$

and

$$\langle \bar{n} | j_{e.m.} | n \rangle \sim (\mu_n/2M) \sigma_{\mu\nu} k_\nu.$$

Hence we can solve for the forms of  $\xi$  and  $\mathcal{O}$ :

$$\xi = -(1/2)(\mu_n/2M) \sigma_{\mu\nu} k_\nu$$

$$\mathcal{O} = -(1/3) \gamma_\mu - (1/6)[(\mu_n + 2\mu_p)/2M] \sigma_{\mu\nu} k_\nu.$$

Now the  $\bar{p}\Lambda K^*$  interaction is

$$\langle \bar{p} | j_{K^*} | \Lambda \rangle \sim 3\mathcal{O} - \xi \sim \gamma_\mu + (\mu_p/2M) \sigma_{\mu\nu} k_\nu.$$

In our theory we have used the expression

$$\langle \bar{p} | j_{K^*} | \Lambda \rangle \sim G_V \gamma_\mu + [G_T/(m_2 + m_4)] \sigma_{\mu\nu} k_\nu.$$

Thus the prediction is

$$G_T/G_V \approx \mu_p = 1.79.$$

#### D. Relationship between $K^- p \rightarrow \Lambda\omega$ and $K^- p \rightarrow \Lambda\phi$

If the  $K$  and  $K^*$  exchange model is valid for  $K^- p \rightarrow \Lambda\omega$ , then we expect the same model to hold for  $K^- p \rightarrow \Lambda\phi$ , with coupling constants related through SU(3). For example, we have

$$\frac{G(\omega K^+ K^-)}{G(\phi K^+ K^-)} = \tan \theta.$$

Also Glashow has pointed out that the ratio

$$R \equiv \left( \frac{G(\phi K^* \bar{K})}{G(\phi K \bar{K})} \right) \left( \frac{G(\omega K^* \bar{K})}{G(\omega K \bar{K})} \right)^{-1}$$



can be expressed as<sup>29</sup>

$$R = -\tan \theta \left\{ \frac{(3/2) \sin \theta \cos \theta - \epsilon [1 - (3/2) \cos^2 \theta]}{1 - (3/2) \sin^2 \theta - (3/2) \epsilon \sin \theta \cos \theta} \right\}.$$

However,  $\Lambda\phi$  elastic scattering in general is not determinable from  $\Lambda\omega$  elastic scattering alone; the absorption parameters used for  $\Lambda\phi$  in general would be different than those used for  $\Lambda\omega$ .

Lindsey has studied the reaction  $K^-p \rightarrow \Lambda\phi$  in the same energy region as we have studied  $K^-p \rightarrow \Lambda\omega$ .<sup>30</sup> Some comments on the relationship of his results to ours are made in Section V.

## V. APPLICATION OF THE ABSORPTION MODEL

### A. Introduction

A computer program which puts Huff's treatment of the absorption model to practical use has been written at Lawrence Radiation Laboratory by J. Friedman (for the reaction  $K^-p \rightarrow K^*p$ ) and modified by L. Hardy and S. Flatté (for  $\pi^-p \rightarrow YK^*$  and  $K^-p \rightarrow \Lambda\omega$ ).

In this section we first compare the results of our treatment with previously published results that used the angular-momentum treatment of the absorption model: We show that the different approximations that are used in the two treatments lead to qualitatively similar, though quantitatively somewhat different, predictions.

After satisfying ourselves that our method and our computer program are valid and useful within the context of the absorption model, we proceed to test the applicability of the absorption model to the reaction  $K^-p \rightarrow \Lambda\omega$ . We attempt to find confirmation of the idea, expressed in Section I, that t-channel exchange mechanisms do not become dominant until the highest momentum region, 2.6 BeV/c, the lower momenta being dominated by threshold and perhaps resonance effects.

The product of the coupling constants for K exchange which is found in the best solution at 2.6 BeV/c compares quite well with the SU(3) prediction derived in Section IV. It then becomes of great interest to see if the  $\Lambda\omega$  coupling constants, appropriately modified, can explain the characteristics of the reaction  $K^-p \rightarrow \Lambda\phi$ . We use some recently available data on  $K^-p \rightarrow \Lambda\phi$  at 2.6 BeV/c to test the absorption model further, and we compare our results with the parameters obtained in  $K^-p \rightarrow \Lambda\omega$ .

### B. Comparison with the Angular-Momentum Method

The two reactions similar to ours whose production characteristics have been explained by using the absorption model are  $\pi^-p \rightarrow \rho^-p$  and  $Kp \rightarrow K^*p$ .<sup>5</sup> Polarization information on the final

fermion in these two reactions is not available; hence the only parameters that have been determined for these reactions are  $f_2, f_3, f_4$  and the differential cross section. The parameters have usually been given in terms of a density-matrix notation<sup>3</sup> for the final vector meson. The equations relating the two notations are

$$\rho_{00} = f_2$$

$$\rho_{1, -1} = 1/2 - f_3 - f_2/2$$

and

$$\text{Re } \rho_{10} = -f_4/2\sqrt{2}.$$

In the angular-momentum treatment, the absorption parameters are expressed in terms of the parameters  $C$  and  $\gamma$  where the absorption factor is

$$\exp(2i\delta) = 1 - C \exp(-\gamma J^2).$$

The correspondence with the total cross section,  $\sigma_T$ , and the slope of the elastic differential cross section,  $A$ , is

$$C = \frac{\sigma_T}{4\pi A}$$

$$\gamma = \frac{1}{2q^2 A}.$$

If  $\sigma_T(i), A_i$  and  $\sigma_T(f), A_f$  are the parameters of initial- and final-state scattering respectively, then we have

$$C_+ = \frac{\sigma_T(i)}{4\pi A_i}$$

$$\gamma_+ = \frac{1}{2q^2 A_i}$$

$$C_- = \frac{\sigma_T(f)}{4\pi A_f}$$

$$\gamma_- = \frac{1}{2q^2 A_f}$$

Figure 22a shows the predictions of the angular-momentum method taken from Jackson et al. for  $\pi^- p \rightarrow \rho^- p$  at 4 BeV/c with  $\pi^0$  exchange.<sup>5</sup> The parameters they used are  $C_+ = 0.76$ ,  $\gamma_+ = 0.04$ ,  $C_- = 1.0$ , and  $\gamma_- = 0.03$ , which translate as  $\sigma_T(i) = 28$  mb,  $A_i = 7.5 (\text{BeV})^{-2}$ ,  $\sigma_T(f) = 56.7$  mb, and  $A_f = 11.6 (\text{BeV})^{-2}$ . In Fig. 22a the squares are the results of our method; the agreement is excellent. This comparison checks only pseudoscalar exchange. To check vector exchange, we take the predictions given by Jackson et al. for the same reaction with some vector exchange added.<sup>5</sup> The curves in Fig. 22b are the predictions of the angular-momentum treatment with the parameters  $\xi$  and  $\eta$ , given by Jackson et al. as

$$\xi = \frac{G(\pi^+ V \rho^-) [G_V(\bar{p} V p) + G_T(\bar{p} V p)]}{2G(\pi^+ \pi^0 \rho^-) G(\bar{p} \pi^0 p)}$$

and

$$\eta = \frac{G(\pi^+ V \rho^-) G_T(\bar{p} V p)}{G(\pi^+ \pi^0 \rho^-) G(\bar{p} \pi^0 p)},$$

set at  $\eta = 0$  and  $\xi = \pm 0.25$  (lower curves) and  $\xi = \pm 0.50$  (upper curves). Here R is the ratio of the results with nonzero  $\xi$  to the results for  $\xi = 0$ . We have determined that  $\xi = \pm 0.25$  corresponds to  $G_V = \pm 34$ ,  $G_T = 0$  and that  $\xi = \pm 0.50$  corresponds to  $G_V = \pm 68$ ,  $G_T = 0$ . Our results are shown as squares (for positive  $G_V$ ) and circles (for negative  $G_V$ ). The agreement in  $\rho_{00}$  seems good, but the differential cross section appears to be in disagreement, our curves being higher than those of Jackson et al. To check further, we look at Fig. 23 which compares our results for  $K^+ p \rightarrow K^* p$  at 3 GeV/c with those of Jackson et al. Again our agreement in the decay parameters is extremely satisfactory, but our differential cross-section curves lie higher than those of Jackson et al.

By comparing results of our program with those of Jackson<sup>21</sup> with no absorption, we find that the definitions of our coupling constants are indeed consistent. Therefore our disagreement arises

from our method of absorption. In Fig. 23 at  $\cos \theta = 0.7$ , our value for the differential cross section is 0.3 mb and that of reference 5 is 0.12 mb. How much of a disagreement is this? We must remember that the crux of the calculation we are making is the calculation of how much the amplitude is absorbed. The unabsorbed cross section at this point is near 6 mb. Therefore we calculated the absorption as 95.0%, and reference 5 calculated 98.0%. In the amplitude this means we calculated 78%, and reference 5 used 86%. Not such a large disagreement when considered in this way! We have discovered two important facts; our calculations and those of the angular-momentum treatment are acceptably close considering the completely different methods used, and the small differences between our answers result in large changes in the differential cross-section predictions. We can now explain why we agree in the forward direction: it is because the absorption is relatively small there, and the calculated cross section is much less sensitive to differences in the absorption calculation. We can also explain why we agree on pseudoscalar exchange results--in fact we don't agree in the nonforward directions, but both our results are so small compared to the forward peak that a large percentage difference goes unnoticed.

We now say the following; the two different treatments of the absorption model agree closely on the effect of pseudoscalar exchange, but disagree by large factors (in two cases, by 2 or 3) on the effect of vector exchange in the differential cross section. The decay correlations are not very sensitive to the difference in the two methods. The differences in the cross section will be buried in the variation of vector coupling constants which are not known. In other words the vector couplings found in reference 5 would need to be reduced by a factor  $\sim 2$ , if used in our treatment. We do not believe that either answer is inherently right; the results are too sensitive to the calculational technique. However, it seems reasonable that the results of one program will be internally consistent, and therefore the ratios of vector couplings determined by one program will have an approximate meaning.

### C. Comparison with Experiment for $K^- p \rightarrow \Lambda \omega$

To predict an experimental result, we must provide the theory with the following parameters:

$$g_P^2 = \frac{G^2(K^- K^+ \omega) G^2(\bar{p} K^+ \Lambda)}{(4\pi)^2} = \text{K-meson-exchange coupling}$$

$$g_V = G(K^- K^* \omega) G_V(\bar{p} K^* \Lambda) = \text{K}^* \text{- exchange vector coupling}$$

$$g_T \equiv G(K^- K^* \omega) G_T(\bar{p} K^* \Lambda) = \text{K}^* \text{ exchange tensor coupling}$$

$$\sigma_T(i) \{ \sigma_T(f) \} = \text{K}^- p \{ \Lambda \omega \} \text{ total cross section}$$

$$A_i \{ A_f \} = \text{K}^- p \{ \Lambda \omega \} \text{ elastic differential cross-section slope in the forward direction}$$

$$\psi = \text{possible nonzero phase of the transition matrix element for elastic scattering.}$$

One of the predictions arising from these parameters is the differential cross section for  $K^- p \rightarrow \Lambda \omega$ . Since the data give cross sections for  $K^- p \rightarrow \Lambda \omega \rightarrow \Lambda \pi^+ \pi^- \pi^0$  we have multiplied all experimental cross sections by 1.1 to account for other  $\omega$  decay modes.

An important comment which we must make immediately is that if we assume  $\psi$  is zero, then the theory will predict that  $f_5$  through  $f_{12}$  are identically zero everywhere. The data at the highest energy are actually not too inconsistent with this prediction; however a non-zero  $\psi$  does significantly improve the fits obtained. In all our fits we have assumed that the  $\psi$  for  $K^- p$  elastic scattering is the same as the  $\psi$  for  $\Lambda \omega$  elastic scattering, and that  $\psi$  is not a function of production

angle. These are drastic approximations, but the effect of  $\psi$  is only felt significantly by  $f_5$  through  $f_{12}$ , which are not very well determined anyway. We note that the values of  $\psi$  for the best fits at 2.6 BeV/c are small, in keeping with our belief that  $\psi$  is close to zero in the very forward direction.

We have set  $\sigma_T(i) = 30$  mb and  $A_i = 7.5$  (BeV)<sup>-2</sup> everywhere. This is certainly a good approximation in the case of  $\sigma_T(i)$ ; Lynch<sup>31</sup> has shown us preliminary data from 1.5 to 2.6 BeV/c in which  $A_i$  varied from 7.0 to 8.5 (BeV)<sup>-2</sup>. Any deviation from 7.5 (BeV)<sup>-2</sup> can easily be taken into account by a small variation in the final-state absorption parameters  $\sigma_T(f)$  and  $A_f$ .

At each momentum we have tried two different fits. First, we have tried K exchange only, varying  $g_p^2$ ,  $\sigma_T(f)$ ,  $A_f$ , and  $\psi$ . Then we have included K\* exchange, adding  $g_V$  and  $g_T$  as parameters. Jackson et al.<sup>5</sup> have already observed that two regions of vector-meson exchange couplings often give comparably good fits to the data; one corresponds to constructive and the other to destructive interference between the vector exchange and the pseudoscalar exchange. We find similar results, and we have tabulated both fits where necessary.

The parameters and  $\chi^2$  for the best fits are given in Table IX, and the curves corresponding to the fits at 2.6 and 1.7 BeV/c are shown in Figs. 24 and 25. It is difficult to state errors on the parameters at 2.6 BeV/c, because the curves are in qualitative but clearly not quantitative agreement with the data. This also results in  $\chi^2$  which are certainly higher than would be acceptable for a perfect theory; one must judge by the curves whether one agrees that "qualitative agreement" has been reached. We prefer to show the curves for the best fits and state that changes of the order of 20% in the coupling constants would definitely give much worse fits. At the lower momenta, errors would not be meaningful, since we wish to argue that the theory is not applicable.

Some comments on the fits at each momentum are made below.

1. 2.6 BeV/c

First we note that  $C_- \equiv \sigma_T(f)/4\pi A_f$  is 1.2. To be consistent with our assumption that the elastic scattering is almost entirely the shadow of the inelastic processes,  $C$  should be  $\leq 1$ . However, since so far  $C_- = 1$  has given the best results in the absorption model, and since we have in no way constrained our parameters to satisfy  $C \leq 1$ , we feel that a value of 1.2 is quite reasonable and acceptable. In fact, since our formalism essentially averages the initial- and final-state S-matrix elements, and since we fixed our initial-state  $C_+ = \sigma_T(i)/4\pi A_i$  at 0.8, then  $C_- = 1.2$  corresponds to total absorption of the s wave, a prominent feature of other successful absorption-model fits.

The total cross section for  $\Lambda\omega$  scattering of  $\sim 80$  mb may be compared with estimates of  $\sim 80$  mb for  $\sigma_T(\rho N)$  made by Drell and Trefil.<sup>32</sup>

Next we note that the K-exchange coupling,  $g_p^2 = G^2(K^- K^+ \omega) G^2(\bar{p} K^+ \Lambda)/(4\pi)^2$ , for the best fit is 7.4, in remarkable agreement with the SU(3) prediction derived in Section IV,  $g_p^2 \approx 7$  to 8. The  $K^*$ -exchange couplings are certainly of a reasonable magnitude.

The curves show a qualitative agreement with the data; the differential-cross-section fit is excellent. The worst quantitative discrepancies occur in  $f_3$  and  $f_{12}$ , but even in  $f_3$  the shape is similar. On the whole, the absorption model appears to give a reasonable qualitative picture at 2.6 BeV/c.

2. 2.1 BeV/c

We find  $\chi^2 = 177$  if we use the parameters determined at 2.6 BeV/c. If the final-state absorption parameters are allowed to vary freely,  $A_f$  goes negative and  $\sigma_T(f)$  becomes small ( $< 5$  mb), a reflection of the lack of forward peaking in the differential cross section. We therefore set  $\sigma_T(f) = 0$  for our final fits. We then find that  $g_p^2$  is at least an order of magnitude below what we expect. (The two fits with  $K^*$  exchange really correspond to more or less the same region. One may think of it as positive  $g_V$  and  $g_T$  with small  $g_p$ ; in one case  $g_p$  is negative and small, in the other positive and small.)

Thus at 2.1 BeV/c, we find two very unpleasant facts, if we want to believe the absorption model. First, the  $\Lambda\omega$  total cross section is extremely small, in contradiction to estimates of the  $\rho N$  total cross sections ( $\sim 80$  mb at 4 BeV/c  $\rho$  laboratory momentum) made by Drell and Trefil,<sup>32</sup> and to other absorption model fits such as  $p\rho^-$  and  $pK^*$ .



Second, the K-exchange coupling is an order of magnitude smaller than one expects.

### 3. 1.7 BeV/c

The fits at 1.7 BeV/c are quite reasonable in all respects. For the final state, we have  $C = 0.6$  for the best fit and  $g_p^2 = 12$ . Since we do not expect the absorption model to apply here, the only comment to make, obviously, is that a theory is not required to fail where it is inapplicable, only to succeed where it is applicable. We find  $\chi^2 = 196$  if we use the parameters determined at 2.6 BeV/c.

### 4. 1.5 BeV/c

Here the model again has trouble. The best-fit value of  $C$  is 5.2, which is clearly unacceptable. Essentially  $A_f$  tends to be much too small. Also  $\psi$  is becoming rather large. Of course when  $C$  is this large, our approximation that  $(1 - T)^{1/2} \cong 1 - \frac{1}{2} T$  is no longer even approximately good. We can say that the 1.5-BeV/c data are not well explained when treated by our method for the absorption model. We find  $\chi^2 = 326$  if we use the parameters determined at 2.6 BeV/c.

## D. Comparison with Experiment for $K^- p \rightarrow \Lambda \phi$

The fact that  $g_p^2$ , the K-exchange coupling, came out quite reasonable for  $K^- p \rightarrow \Lambda \omega$  at 2.6 BeV/c is gratifying. It then becomes of great interest whether the characteristics of  $K^- p \rightarrow \Lambda \phi$  are consistent with these couplings also. We have translated the results of Lindsey<sup>30</sup> on  $K^- p \rightarrow \Lambda \phi$  at 2.6 BeV/c into our notation and plotted the results in Fig. 26. The solid curve is calculated from the parameters determined at 2.6 BeV/c for  $K^- p \rightarrow \Lambda \omega$ , appropriately modified. The modifications, given in Section IV, are

$$g_p(\Lambda \phi) = g_p(\Lambda \omega) \cot \theta_m = 1.19 g_p(\Lambda \omega)$$

and

$$g_V \text{ or } T(\Lambda \phi) = g_V \text{ or } T(\Lambda \omega) [-R \cot \theta_m] = -1.7 g_V \text{ or } T(\Lambda \omega).$$

The  $\chi^2$  for the solid curve is 552 for 28 data points, where we have assumed that the error matrix for the decay parameters is diagonal. This is not a bad assumption.<sup>30</sup>

We then allowed all parameters to vary and found as the best fit the dashed curve in Fig. 26. The  $\chi^2$  is 77.8. The parameters of the dashed curve are  $\sigma_T(f) = 81.6$  mb,  $A_f = 13.1 (\text{BeV})^{-2}$ ,  $\psi = 0.07$ ,  $g_p^2 = 6.9$ ,  $g_V = 32.0$ , and  $g_T = 20.3$ . Hence we have

$$\frac{G(K^-K^+\phi)}{G(K^-K^+\omega)} \cong \left( \frac{6.9}{7.4} \right)^{1/2} = 1.0$$

where we expect 1.2,

$$\frac{G_V(K^-K^*\phi)}{G_V(K^-K^*\omega)} \cong \frac{32.0}{-28.9} = -1.1$$

where we expect -1.7, and

$$\frac{G_T(K^-K^*\phi)}{G_T(K^-K^*\omega)} \cong \frac{20.3}{-8.8} = -2.3$$

where we expect -1.7. Of course we have chosen the fit to  $\Lambda\omega$  which best meets the predictions (it is also the best fit).

Glashow<sup>29</sup> has pointed out that the test that is least sensitive to kinematical effects is the ratio R. We find using  $G_V$  that

$$R = \left( \frac{32.0}{\sqrt{6.9}} \right) \left( \frac{\sqrt{7.4}}{-28.9} \right) = -1.1$$

where we expect -1.4, and using  $G_T$  we find

$$R = \left( \frac{20.3}{\sqrt{6.9}} \right) \left( \frac{\sqrt{7.4}}{-8.8} \right) = -2.4$$

where we expect -1.4.

Cabibbo's prediction (see Section IV) that  $G_T/G_V = 1.79$  is not verified since  $G_T$  in our best fits is about a factor of two smaller than

$G_V$ . Cabibbo's scheme also predicts  $G_T/G_V$  for the  $\bar{p}p\rho^0$  vertex as  $\mu_p - \mu_n = 4.7$ . Using the reaction  $K^+p \rightarrow K^*p$ , Jackson et al.<sup>5</sup> found  $G_T/G_V(\bar{p}p\rho^0) \lesssim 1$ , again not in agreement with the prediction.

Since the vector-meson exchange formalism is in much doubt, the vector-coupling comparisons may be academic; however, we have avoided the main problem of vector exchange -- its energy dependence -- by working always at the same energy (although not at the same distance from threshold).

The orders of magnitude seem to be clearly in order, and even the signs seem of some significance. (The signs are relative ones between  $g_p$  and either  $g_v$  or  $g_T$ .)

#### E. Conclusions

We have shown that the characteristics of the reactions  $K^-p \rightarrow \Lambda\omega$  and  $K^-p \rightarrow \Lambda\phi$  at 2.6 BeV/c at forward production angles are explained reasonably well by the absorption model with K and  $K^*$  exchange. The couplings obtained from best fits to the data are in remarkable qualitative agreement with the predictions of SU(3).

The qualitative features of the reaction  $K^-p \rightarrow \Lambda\omega$  at lower momenta (namely the lack of a strong forward peak) indicate that t-channel exchanges are not dominant; therefore we would not expect the absorption model to work. If it did work we could not fault the model, but a theory which works everywhere, regardless of whether it is applicable or not, is not a very testable theory. We find that at 1.5 and 2.1 BeV/c the absorption model does fail to explain the data, while at 1.7 BeV/c it works.

## VI. ACKNOWLEDGMENTS

I thank Professors M. Lynn Stevenson and Luis W. Alvarez for their continued support, encouragement, and guidance. I also thank Jerome Friedman, Dr. Ronald Ross, and Dr. Robert Huff for their many discussions on the absorption model.

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## FIGURE LEGENDS

- Fig. 1. (a) Diagram for  $K^-p \rightarrow \Lambda\omega$  representing one-particle exchange. We consider particle  $e$  as being a  $K$  or  $K^*$  meson.
- (b) Diagram used in the absorption-model calculations. The shaded blobs represent elastic scattering.
- Fig. 2. Cross sections of the reactions (top curve)  $K^-p \rightarrow \Lambda\pi^+\pi^-\pi^0$  and (bottom curve)  $K^-p \rightarrow \Lambda + (\omega \rightarrow \pi^+\pi^-\pi^0)$  from threshold to 3 BeV/c incident- $K^-$  momentum. The connecting lines are only to eliminate confusion.
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- Fig. 3. Production-angle distribution for  $K^-p \rightarrow \Lambda\omega$  at 1.5 BeV/c; background has been subtracted, leaving 3570 events.
- Fig. 4. Production-angle distribution for  $K^-p \rightarrow \Lambda\omega$  at 1.7 BeV/c; background has been subtracted, leaving 1570 events.



- Fig. 5. Production-angle distribution for  $K^-p \rightarrow \Lambda\omega$  at 2.1 BeV/c; background has been subtracted, leaving 1021 events.
- Fig. 6. Production-angle distribution for  $K^-p \rightarrow \Lambda\omega$  at 2.6 BeV/c; background has been subtracted leaving 1300 events.
- Fig. 7. Feynman diagrams representing exchanges in the three channels that affect  $K^-p \rightarrow \Lambda\omega$ . Exchanges of the least massive particles are shown. (a) s channel, (b) t channel, (c) u channel.
- Fig. 8. Distribution of beam momentum for 31 800 events of the type  $K^-p \rightarrow \Lambda\pi^+\pi^-\pi^0$ . Over one million pictures were taken to gather this data.
- Fig. 9. Distribution of  $\cos\psi = \frac{\underline{\Lambda} \cdot \underline{\pi}}{\Lambda \pi}$  in the  $\Lambda$  rest frame, where  $\underline{\Lambda}$  is the direction of the  $\Lambda$  in the laboratory. This graph, for  $K^-p \rightarrow \Lambda\pi^+\pi^-\pi^0$  events at 1.7 BeV/c in the  $\omega$  region, exhibits no bias against any opening angle for the  $\Lambda$ .
- Fig. 10. Distribution of length of the  $\Lambda$  path for  $K^-p \rightarrow \Lambda\pi^+\pi^-\pi^0$  events at 1.7 BeV/c in the  $\omega$  region, showing a loss of events at small lengths due to scanning bias. The dashed curve is the expected exponential if all  $\Lambda$ 's had a laboratory momentum of 1 BeV/c, which is about the average in this sample.
- Fig. 11. Production-angle distributions for  $K^-p \rightarrow \Lambda\pi^+\pi^-\pi^0$  events at 2.6 BeV/c with  $750 \text{ MeV} < M(\pi^+\pi^-\pi^0) < 810 \text{ MeV}$ . The boxes indicate points we obtained by imposing a 2-mm cutoff on the  $\Lambda$  laboratory length and appropriately weighting the remaining events.
- Fig. 12. Production-angle distributions for  $K^-p \rightarrow \Lambda\omega$  events at 2.6 BeV/c in the  $\omega$  region which (a) passed first measurement, and (b) failed first measurement but passed second measurement.
- Fig. 13. Schematic drawing of a  $K^-p \rightarrow \Lambda\omega$  reaction indicating momentum vectors and the production angle  $\theta$  in the c.m. system.
- Fig. 14. Three-pion mass distributions for 7720  $K^-p \rightarrow \Lambda\pi^+\pi^-\pi^0$  events at 1.5 BeV/c. The curves in this and the next three plots are hand-drawn estimates of background under the  $\omega$  peak.

Fig. 15. Three-pion mass distributions for 4900  $K^-p \rightarrow \Lambda \pi^+ \pi^- \pi^0$  events at 1.7 BeV/c.

Fig. 16. Three-pion mass distributions for 5560  $K^-p \rightarrow \Lambda \pi^+ \pi^- \pi^0$  events at 2.1 BeV/c.

Fig. 17. Three-pion mass distributions for 11 830  $K^-p \rightarrow \Lambda \pi^+ \pi^- \pi^0$  events at 2.6 BeV/c. The statistically significant peaks at 960 and 1020 MeV are the  $\pi^+ \pi^- \gamma$  decays of the  $\eta(959)$  and the  $3\pi$  decay of the  $\phi$ , respectively.

Fig. 18. Distributions of the number of events and the decay correlations  $\{f_i\}$  as a function of production-angle cosine for  $K^-p \rightarrow \Lambda \omega$  events at 1.5 BeV/c. Dotted lines are used to eliminate possible confusion.

Fig. 19. Distributions of the number of events and the decay correlations  $\{f_i\}$  as a function of production-angle cosine for  $K^-p \rightarrow \Lambda \omega$  events at 1.7 BeV/c. Dotted lines are used to eliminate possible confusion.

Fig. 20. Distributions of the number of events and the decay correlations  $\{f_i\}$  as a function of production-angle cosine for  $K^-p \rightarrow \Lambda \omega$  events at 2.1 BeV/c. Dotted lines are used to eliminate possible confusion.

Fig. 21. Distributions of the number of events and the decay correlations  $\{f_i\}$  as a function of production-angle cosine for  $K^-p \rightarrow \Lambda \omega$  events at 2.6 BeV/c. Dotted lines are used to eliminate possible confusion.

Fig. 22. Parameters predicted by the absorption model for the reaction  $\pi^- p + \rho^- p$  at 4 GeV/c. The curves are taken from Jackson et al;<sup>5</sup> and the points are from our method. (a) Pion exchange only; (b) some vector exchange added in the form of  $\xi = \pm 0.25$  (lower solid and dotted curves and lower squares and circles) and  $\xi = \pm 0.50$  (upper solid and dotted curves and upper squares and circles).

- Fig. 23. Parameters predicted by the absorption model for the reaction  $K^+p \rightarrow K^{*+}p$  at 3 GeV/c. The curves are taken from Jackson et al.,<sup>5</sup> and the points are from our method. The squares correspond to the solid lines and the circles to the dotted lines. See text for absorption parameters and coupling constants.
- Fig. 24. Data at forward production angles for 2.6 BeV/c  $K^-p \rightarrow \Lambda\omega$  events. The solid curves correspond to the best fits for K exchange only, the dashed curves to  $KK^*(1)$  best fit, the dash-dot curve to  $KK^*(2)$  best fit, and the dotted curve to K exchange only with no absorption and a K coupling equal to that used in the solid curves.
- Fig. 25. Data at forward production angles for 1.7 BeV/c  $K^-p \rightarrow \Lambda\omega$  events. The solid curves correspond to the best fits for K exchange only, the dashed curves to  $KK^*(1)$  best fit, the dash-dot curve to  $KK^*(2)$  best fit, and the dotted curve to K exchange only with no absorption and a K coupling equal to that used in the solid curves.
- Fig. 26. Data taken from Lindsey<sup>29</sup> for the reaction  $K^-p \rightarrow \Lambda\phi$  at 2.6 BeV/c. The solid curves are the predictions of the absorption model with absorption parameters identical to those determined by the  $KK^*(1)$  solution to  $\Lambda\omega$  at 2.6 BeV/c and with coupling constants obtained from the  $\Lambda\omega$  fit by invoking exact SU(3). The dashed curves correspond to the best fit with a free variation of parameters. See text for a comparison of couplings determined in the best fits.

Table I. Total cross sections for  $K^- p \rightarrow \Lambda \omega \rightarrow \Lambda \pi^+ \pi^- \pi^0$ .

Run	Momentum (BeV/c)	Path length (events/ $\mu$ b)	$\sigma_{\Lambda 3\pi}^a$ (mb)	N( $3\pi$ )	N( $\omega$ )	$\sigma_{\Lambda \omega}^a$ (mb)
K72	1.22	$1.23 \pm 0.06$	$0.68 \pm 0.05$	392	$0 \pm 30$	$0 \pm 0.05$
	1.32	$1.44 \pm 0.07$	$1.53 \pm 0.10$	965	502	$0.80 \pm 0.06$
	1.42	$0.83 \pm 0.04$	$2.10 \pm 0.06$	1093	505	$0.97 \pm 0.08$
	1.51	$5.09 \pm 0.20$	$2.26 \pm 0.08$	5847	2475	$0.96 \pm 0.05$
	1.60	$0.72 \pm 0.04$	$2.14 \pm 0.15$	1006	366	$0.78 \pm 0.05$
	1.70	$1.10 \pm 0.06$	$2.82 \pm 0.17$	1000	357	$1.01 \pm 0.06$
K63	1.7	$3.86 \pm 0.20^b$	$1.66 \pm 0.25$	2923	1060	$0.58 \pm 0.10$
	2.1	$6.04 \pm 0.30^b$	$1.98 \pm 0.20$	5563	1299	$0.46 \pm 0.05$
	2.6	$16.5 \pm 0.9^b$	$1.55 \pm 0.16$	11831	1660	$0.22 \pm 0.03$

a. Corrected for neutral  $\Lambda$  decay.

b. For 2.1 and 2.6 BeV/c path lengths were obtained from Lindsey.<sup>29</sup> At 1.7 BeV/c we used the same method of counting taus as he did.

Table II. Scanning and measuring correction factors for total-cross-section determinations of  $V +$  two-prong events.

	Momentum (BeV/c)		
	1.7	2.1	2.6
Events on first scan	17 988	28 326	71 722
Scanning efficiency (%)	$94 \pm 3$	$97 \pm 3$	$94 \pm 3$
Scanning correction factor <sup>a</sup>	$1.10 \pm 0.04$	$1.08 \pm 0.04$	$1.14 \pm 0.04$
Events not measured	1682	1064	5848
Events measured at least once	16 306	27 262	65 874
Events passing first measurement	10 652	19 266	38 420
First-measurement passing fraction (%)	65	71	58
Events measured at least twice	3745	0	14 877
Events passing second measurement	1720	0	7040
Second-measurement passing fraction (%)	46	47 <sup>b</sup>	47
Events measured at least thrice	0	0	2896
Events passing third measurement	0	0	1448
Third-measurement passing fraction	50 <sup>b</sup>	50 <sup>b</sup>	50
Total passing events	12 372	19 266	46 908
Projected passing events	15 500	25 300	59 800
Measuring correction factor	$1.26 \pm 0.04$	$1.31 \pm 0.04$	$1.28 \pm 0.04$

a. Includes corrections for short  $\Lambda$  and escaping  $\Lambda$ .

b. Where no information is available, the passing fraction of 2.6 BeV/c is used.

Table III. Correlation parameters for incident  $K^-$  laboratory momentum  $\approx 1.5$  BeV/c. The c.m. energy is  $\approx 2.02$  BeV. The total cross section for  $K^-p \rightarrow \Lambda_c^- + \Lambda^- + \pi^+\pi^-\pi^0$  near 1.5 BeV/c is  $0.96 \pm 0.05$  mb. The total number of  $\omega$  events at all  $\cos \theta$  is 3568.

$\cos \theta_{\max}$	1.00	0.75	0.50	0.25	0.00	-0.25	-0.50	-0.75
$\cos \theta_{\min}$	0.75	0.50	0.25	0.00	-0.25	-0.50	-0.75	-1.00
$N(E, \cos \theta)^a$	704	770	818	694	550	414	310	237
$C(E, \cos \theta)^b$	262	244	253	245	207	261	252	248
$f_1$	$0.279 \pm 0.034$	$0.260 \pm 0.031$	$0.222 \pm 0.029$	$0.188 \pm 0.031$	$0.302 \pm 0.039$	$0.426 \pm 0.053$	$0.359 \pm 0.070$	$0.489 \pm 0.100$
$f_2$	$0.392 \pm 0.034$	$0.500 \pm 0.033$	$0.599 \pm 0.032$	$0.571 \pm 0.035$	$0.551 \pm 0.039$	$0.420 \pm 0.054$	$0.479 \pm 0.072$	$0.336 \pm 0.091$
$f_3$	$0.329 \pm 0.033$	$0.240 \pm 0.031$	$0.179 \pm 0.028$	$0.241 \pm 0.032$	$0.147 \pm 0.034$	$0.154 \pm 0.048$	$0.162 \pm 0.066$	$0.175 \pm 0.092$
$f_4$	$0.288 \pm 0.059$	$0.313 \pm 0.053$	$0.369 \pm 0.048$	$0.385 \pm 0.057$	$0.350 \pm 0.064$	$0.376 \pm 0.087$	$0.148 \pm 0.116$	$-0.162 \pm 0.179$
$f_5$	$-0.053 \pm 0.064$	$-0.099 \pm 0.053$	$-0.101 \pm 0.050$	$-0.007 \pm 0.055$	$0.059 \pm 0.077$	$0.011 \pm 0.106$	$0.118 \pm 0.133$	$0.414 \pm 0.174$
$f_6$	$-0.299 \pm 0.062$	$-0.066 \pm 0.065$	$-0.137 \pm 0.066$	$-0.240 \pm 0.070$	$-0.250 \pm 0.079$	$-0.032 \pm 0.107$	$0.138 \pm 0.136$	$-0.219 \pm 0.166$
$f_7$	$0.027 \pm 0.059$	$0.021 \pm 0.057$	$0.084 \pm 0.052$	$0.101 \pm 0.055$	$-0.013 \pm 0.062$	$0.051 \pm 0.082$	$0.106 \pm 0.133$	$-0.249 \pm 0.163$
$f_8$	$-0.540 \pm 0.101$	$-0.118 \pm 0.091$	$-0.259 \pm 0.085$	$-0.130 \pm 0.098$	$-0.192 \pm 0.110$	$-0.332 \pm 0.154$	$-0.405 \pm 0.189$	$-0.147 \pm 0.294$
$f_9$	$0.204 \pm 0.101$	$0.244 \pm 0.102$	$0.318 \pm 0.091$	$0.005 \pm 0.104$	$-0.267 \pm 0.120$	$0.139 \pm 0.161$	$-0.143 \pm 0.228$	$-0.743 \pm 0.307$
$f_{10}$	$-0.077 \pm 0.110$	$-0.091 \pm 0.092$	$-0.087 \pm 0.080$	$-0.074 \pm 0.090$	$-0.244 \pm 0.105$	$0.160 \pm 0.156$	$0.471 \pm 0.208$	$0.118 \pm 0.284$
$f_{11}$	$-0.197 \pm 0.091$	$-0.179 \pm 0.094$	$-0.014 \pm 0.093$	$-0.016 \pm 0.105$	$0.289 \pm 0.119$	$0.072 \pm 0.164$	$0.439 \pm 0.225$	$0.298 \pm 0.288$
$f_{12}$	$0.191 \pm 0.096$	$0.139 \pm 0.083$	$0.210 \pm 0.083$	$0.333 \pm 0.088$	$0.127 \pm 0.103$	$0.407 \pm 0.148$	$0.396 \pm 0.186$	$0.178 \pm 0.300$

a. N is the number of events in the  $\omega$  region.

b. C is the number of events in the control region.

Table IV. Correlation parameters for incident  $K^-$  laboratory momentum  $\approx 1.7$  BeV/c. The c.m. energy is  $\approx 2.10$  BeV. The total cross section for  $K^-p \rightarrow \Lambda\omega \rightarrow \Lambda\pi^+\pi^-\pi^0$  near 1.7 BeV/c is  $0.80 \pm 0.15$  mb. The total number of  $\omega$  events at all  $\cos\theta$  is 1566.

$\cos\theta_{\max}$	1.00	0.75	0.50	0.25	0.00	-0.25	-0.50	-0.75
$\cos\theta_{\min}$	0.75	0.50	0.25	0.00	-0.25	-0.50	-0.75	-1.00
$N(E, \cos\theta)^a$	389	425	394	366	330	335	306	266
$C(E, \cos\theta)^b$	218	174	163	194	161	163	238	232
$f_1$	$0.350 \pm 0.053$	$0.178 \pm 0.041$	$0.280 \pm 0.045$	$0.206 \pm 0.047$	$0.142 \pm 0.048$	$0.333 \pm 0.051$	$0.256 \pm 0.065$	$0.245 \pm 0.070$
$f_2$	$0.418 \pm 0.050$	$0.664 \pm 0.046$	$0.587 \pm 0.048$	$0.539 \pm 0.053$	$0.477 \pm 0.052$	$0.411 \pm 0.051$	$0.589 \pm 0.067$	$0.411 \pm 0.073$
$f_3$	$0.232 \pm 0.049$	$0.158 \pm 0.039$	$0.133 \pm 0.042$	$0.255 \pm 0.050$	$0.380 \pm 0.054$	$0.255 \pm 0.050$	$0.154 \pm 0.060$	$0.304 \pm 0.074$
$f_4$	$0.253 \pm 0.086$	$0.227 \pm 0.076$	$0.312 \pm 0.076$	$-0.109 \pm 0.091$	$0.189 \pm 0.098$	$0.419 \pm 0.087$	$0.211 \pm 0.117$	$-0.211 \pm 0.121$
$f_5$	$-0.092 \pm 0.100$	$0.100 \pm 0.076$	$-0.133 \pm 0.074$	$-0.124 \pm 0.081$	$-0.137 \pm 0.086$	$-0.006 \pm 0.098$	$0.068 \pm 0.107$	$0.370 \pm 0.137$
$f_6$	$-0.042 \pm 0.097$	$0.002 \pm 0.099$	$0.087 \pm 0.103$	$0.095 \pm 0.115$	$-0.190 \pm 0.113$	$0.136 \pm 0.103$	$0.070 \pm 0.146$	$-0.040 \pm 0.155$
$f_7$	$-0.048 \pm 0.086$	$-0.162 \pm 0.074$	$0.057 \pm 0.070$	$-0.008 \pm 0.090$	$0.149 \pm 0.106$	$0.058 \pm 0.082$	$0.087 \pm 0.117$	$0.124 \pm 0.139$
$f_8$	$-0.603 \pm 0.133$	$-0.009 \pm 0.128$	$-0.089 \pm 0.132$	$-0.002 \pm 0.151$	$-0.040 \pm 0.181$	$0.120 \pm 0.155$	$-0.259 \pm 0.195$	$-0.374 \pm 0.198$
$f_9$	$0.307 \pm 0.165$	$0.261 \pm 0.139$	$0.094 \pm 0.155$	$0.269 \pm 0.161$	$-0.081 \pm 0.173$	$0.033 \pm 0.165$	$-0.281 \pm 0.201$	$-0.040 \pm 0.251$
$f_{10}$	$-0.205 \pm 0.139$	$-0.199 \pm 0.131$	$-0.056 \pm 0.130$	$-0.205 \pm 0.141$	$-0.277 \pm 0.132$	$-0.276 \pm 0.153$	$-0.274 \pm 0.188$	$0.351 \pm 0.211$
$f_{11}$	$-0.033 \pm 0.148$	$0.179 \pm 0.131$	$0.189 \pm 0.147$	$-0.156 \pm 0.146$	$0.014 \pm 0.156$	$0.175 \pm 0.162$	$0.676 \pm 0.183$	$0.762 \pm 0.243$
$f_{12}$	$-0.069 \pm 0.153$	$0.145 \pm 0.121$	$0.041 \pm 0.113$	$0.091 \pm 0.144$	$0.190 \pm 0.133$	$0.351 \pm 0.141$	$0.370 \pm 0.174$	$0.462 \pm 0.194$

a. N is the number of events in the  $\omega$  region.

b. C is the number of events in the control region.

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Table V. Correlation parameters for incident  $K^-$  laboratory momentum  $\approx 2.1$  BeV/c. The c.m. energy is  $\approx 2.27$  BeV. The total cross section for  $K^-p \rightarrow \Lambda\omega \rightarrow \Lambda\pi^+\pi^-\pi^0$  near 2.1 BeV/c is  $0.48 \pm 0.05$  mb. The total number of  $\omega$  events at all  $\cos\theta$  is 1021.

$\cos\theta_{\max}$	1.00	0.75	0.50	0.25	0.00	-0.25	-0.50	-0.75
$\cos\theta_{\min}$	0.75	0.50	0.25	0.00	-0.25	-0.50	-0.75	-1.00
$N(E, \cos\theta)^a$	122	123	156	143	165	188	161	152
$C(E, \cos\theta)^b$	84	87	73	60	75	94	99	101
$f_1$	$0.572 \pm 0.090$	$0.433 \pm 0.097$	$0.370 \pm 0.072$	$0.331 \pm 0.070$	$0.181 \pm 0.059$	$0.190 \pm 0.056$	$0.263 \pm 0.073$	$0.261 \pm 0.072$
$f_2$	$0.105 \pm 0.071$	$0.237 \pm 0.090$	$0.249 \pm 0.069$	$0.299 \pm 0.066$	$0.521 \pm 0.066$	$0.624 \pm 0.061$	$0.366 \pm 0.075$	$0.302 \pm 0.075$
$f_3$	$0.323 \pm 0.085$	$0.330 \pm 0.089$	$0.381 \pm 0.071$	$0.370 \pm 0.073$	$0.298 \pm 0.059$	$0.186 \pm 0.055$	$0.371 \pm 0.071$	$0.437 \pm 0.079$
$f_4$	$0.013 \pm 0.145$	$0.334 \pm 0.135$	$0.271 \pm 0.119$	$0.547 \pm 0.113$	$-0.514 \pm 0.127$	$0.169 \pm 0.118$	$0.088 \pm 0.141$	$-0.092 \pm 0.136$
$f_5$	$-0.039 \pm 0.186$	$-0.002 \pm 0.195$	$-0.245 \pm 0.126$	$0.033 \pm 0.128$	$-0.009 \pm 0.123$	$-0.124 \pm 0.120$	$-0.143 \pm 0.136$	$0.193 \pm 0.150$
$f_6$	$-0.031 \pm 0.125$	$0.313 \pm 0.172$	$-0.000 \pm 0.122$	$0.253 \pm 0.111$	$0.146 \pm 0.138$	$0.209 \pm 0.128$	$0.003 \pm 0.148$	$-0.080 \pm 0.138$
$f_7$	$0.005 \pm 0.147$	$-0.222 \pm 0.152$	$0.022 \pm 0.147$	$-0.006 \pm 0.145$	$-0.062 \pm 0.125$	$0.173 \pm 0.093$	$-0.129 \pm 0.136$	$0.248 \pm 0.152$
$f_8$	$-0.313 \pm 0.235$	$-0.252 \pm 0.262$	$0.153 \pm 0.213$	$0.512 \pm 0.224$	$0.507 \pm 0.237$	$-0.039 \pm 0.186$	$-0.076 \pm 0.240$	$-0.144 \pm 0.228$
$f_9$	$0.257 \pm 0.255$	$0.114 \pm 0.245$	$0.083 \pm 0.207$	$0.472 \pm 0.205$	$0.021 \pm 0.194$	$0.265 \pm 0.182$	$-0.356 \pm 0.201$	$-0.158 \pm 0.235$
$f_{10}$	$-0.455 \pm 0.280$	$-0.123 \pm 0.272$	$0.292 \pm 0.231$	$-0.021 \pm 0.216$	$-0.281 \pm 0.212$	$-0.210 \pm 0.141$	$0.062 \pm 0.229$	$-0.019 \pm 0.262$
$f_{11}$	$0.515 \pm 0.245$	$0.263 \pm 0.229$	$0.187 \pm 0.199$	$0.371 \pm 0.214$	$-0.308 \pm 0.194$	$-0.760 \pm 0.205$	$-0.300 \pm 0.185$	$-0.095 \pm 0.224$
$f_{12}$	$-0.181 \pm 0.284$	$0.365 \pm 0.287$	$-0.173 \pm 0.202$	$-0.017 \pm 0.210$	$-0.082 \pm 0.208$	$-0.119 \pm 0.191$	$0.107 \pm 0.244$	$0.241 \pm 0.207$

a. N is the number of events in the  $\omega$  region.

b. C is the number of events in the control region.



Table VI. Correlation parameters for incident  $K^-$  laboratory momentum  $\approx 2.6$  BeV/c. The c.m. energy is  $\approx 2.49$  BeV. The total cross section for  $K^-p \rightarrow \Lambda\omega \rightarrow \Lambda\pi^+\pi^-\pi^0$  near 2.6 BeV/c is  $0.22 \pm 0.03$  mb. The total number of  $\omega$  events at all  $\cos\theta$  is 1300.

$\cos\theta_{\max}$	1.00	0.75	0.50	0.25	0.00	-0.25	-0.50	-0.75
$\cos\theta_{\min}$	0.75	0.50	0.25	0.00	-0.25	-0.50	-0.75	-1.00
$N(E, \cos\theta)^a$	435	180	132	123	118	180	210	201
$C(E, \cos\theta)^b$	145	115	74	79	84	111	122	168
$f_1$	$0.504 \pm 0.041$	$0.586 \pm 0.069$	$0.517 \pm 0.082$	$0.271 \pm 0.092$	$0.317 \pm 0.094$	$0.125 \pm 0.058$	$0.112 \pm 0.057$	$0.374 \pm 0.075$
$f_2$	$0.397 \pm 0.041$	$0.349 \pm 0.072$	$0.348 \pm 0.083$	$0.149 \pm 0.072$	$0.465 \pm 0.101$	$0.359 \pm 0.071$	$0.320 \pm 0.067$	$0.251 \pm 0.074$
$f_3$	$0.099 \pm 0.030$	$0.065 \pm 0.062$	$0.134 \pm 0.070$	$0.580 \pm 0.090$	$0.218 \pm 0.089$	$0.516 \pm 0.069$	$0.568 \pm 0.070$	$0.375 \pm 0.078$
$f_4$	$0.132 \pm 0.062$	$0.226 \pm 0.098$	$0.200 \pm 0.133$	$0.242 \pm 0.167$	$0.317 \pm 0.150$	$0.376 \pm 0.128$	$0.361 \pm 0.109$	$0.032 \pm 0.132$
$f_5$	$-0.185 \pm 0.085$	$-0.474 \pm 0.143$	$-0.436 \pm 0.178$	$0.210 \pm 0.163$	$0.150 \pm 0.202$	$-0.006 \pm 0.104$	$0.014 \pm 0.098$	$0.384 \pm 0.111$
$f_6$	$0.059 \pm 0.079$	$0.293 \pm 0.152$	$0.032 \pm 0.176$	$0.231 \pm 0.113$	$0.490 \pm 0.204$	$0.028 \pm 0.142$	$0.188 \pm 0.142$	$-0.039 \pm 0.134$
$f_7$	$-0.045 \pm 0.062$	$-0.140 \pm 0.102$	$0.300 \pm 0.170$	$0.379 \pm 0.187$	$0.076 \pm 0.148$	$-0.003 \pm 0.141$	$0.039 \pm 0.143$	$0.077 \pm 0.142$
$f_8$	$-0.248 \pm 0.103$	$-0.094 \pm 0.173$	$0.067 \pm 0.246$	$0.354 \pm 0.284$	$0.404 \pm 0.239$	$-0.089 \pm 0.224$	$-0.085 \pm 0.187$	$-0.025 \pm 0.237$
$f_9$	$-0.001 \pm 0.138$	$-0.509 \pm 0.235$	$-0.526 \pm 0.267$	$0.321 \pm 0.241$	$0.278 \pm 0.245$	$0.031 \pm 0.183$	$0.029 \pm 0.176$	$-0.211 \pm 0.231$
$f_{10}$	$-0.043 \pm 0.119$	$-0.265 \pm 0.212$	$0.213 \pm 0.289$	$-0.052 \pm 0.258$	$0.092 \pm 0.293$	$-0.351 \pm 0.218$	$-0.273 \pm 0.191$	$-0.086 \pm 0.221$
$f_{11}$	$-0.033 \pm 0.126$	$0.224 \pm 0.246$	$-0.226 \pm 0.222$	$0.436 \pm 0.252$	$-0.171 \pm 0.281$	$-0.286 \pm 0.185$	$-0.234 \pm 0.148$	$0.262 \pm 0.209$
$f_{12}$	$0.218 \pm 0.110$	$0.405 \pm 0.184$	$0.311 \pm 0.206$	$0.201 \pm 0.259$	$0.342 \pm 0.312$	$-0.189 \pm 0.229$	$0.370 \pm 0.143$	$0.523 \pm 0.223$

a. N is the number of events in the  $\omega$  region.

b. C is the number of events in the control region.

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Table VII. 2.6-BeV/c correlation parameters  
for forward production angles.

$\text{Cos } \theta_{\text{max}}$	1.00	0.875	0.75
$\text{Cos } \theta_{\text{min}}$	0.875	0.75	0.625
$N(E, \text{cos } \theta)^a$	252	183	118
$C(E, \text{cos } \theta)^b$	83	62	66
$f_1$	$0.464 \pm 0.054$	$0.546 \pm 0.060$	$0.667 \pm 0.080$
$f_2$	$0.414 \pm 0.053$	$0.378 \pm 0.062$	$0.325 \pm 0.090$
$f_3$	$0.121 \pm 0.041$	$0.076 \pm 0.039$	$0.008 \pm 0.067$
$f_4$	$0.197 \pm 0.086$	$0.077 \pm 0.090$	$0.098 \pm 0.119$
$f_5$	$-0.166 \pm 0.107$	$-0.297 \pm 0.137$	$-0.418 \pm 0.177$
$f_6$	$0.091 \pm 0.096$	$-0.022 \pm 0.127$	$0.148 \pm 0.181$
$f_7$	$0.085 \pm 0.099$	$-0.085 \pm 0.086$	$-0.029 \pm 0.119$
$f_8$	$-0.387 \pm 0.156$	$-0.186 \pm 0.166$	$0.014 \pm 0.218$
$f_9$	$0.133 \pm 0.180$	$-0.174 \pm 0.212$	$-0.163 \pm 0.298$
$f_{10}$	$-0.032 \pm 0.153$	$0.101 \pm 0.190$	$-0.508 \pm 0.306$
$f_{11}$	$0.068 \pm 0.166$	$-0.215 \pm 0.192$	$0.332 \pm 0.312$
$f_{12}$	$0.065 \pm 0.138$	$0.588 \pm 0.200$	$0.486 \pm 0.246$

a. N is the number of events in the  $\omega$  region.

b. C is the number of events in the control region.



Table IX. Best-fit parameters for  $K^-p \rightarrow \Lambda\omega$ .

Momentum (BeV/c)	Theory <sup>a</sup>	No. of data points	$\chi^2$	$\sigma_T(f)$ (mb)	$A_f$ (BeV) <sup>-2</sup>	$\psi$	$g_p^2$	$g_V$	$g_T$
1.5	K	26	144	45.0	4.7	0.47	10.4	0	0
1.5	KK*(1)	26	140	9.1	-6.0	-0.54	3.0	-24.6	-6.2
1.5	KK*(2)	26	54	41.9	1.7	0.36	16.6	36.7	22.1
1.7	K	26	71	61.9	9.8	0.51	12.7	0	0
1.7	KK*(1)	26	70	57.0	19.5	0.02	12.2	-28.2	-5.8
1.7	KK*(2)	26	49	60.8	10.5	0.43	14.9	20.5	10.1
2.1	K	26	302	0	0	-0.14	2.9	0	0
2.1	KK*(1)	26	66	0	0	-0.06	0.5	-11.8	-9.9
2.1	KK*(2)	26	76	0	0	-0.20	0.7	10.9	10.1
2.6	K	36	167	83.1	13.8	0.16	16.2	0	0
2.6	KK*(1)	36	68	83.0	14.2	-0.10	7.4	-28.9	-8.8
2.6	KK*(2)	36	83	83.2	13.7	0.12	11.8	25.6	3.4
2.6	KK*(K <sup>-</sup> p → Λφ)	28	78	81.6	13.1	0.07	6.9	32.0	20.3

a. K means K exchange only. KK\*(1) means K and K\* exchange with the relative sign between the K\* and K couplings negative; KK\*(2) means the sign is positive.

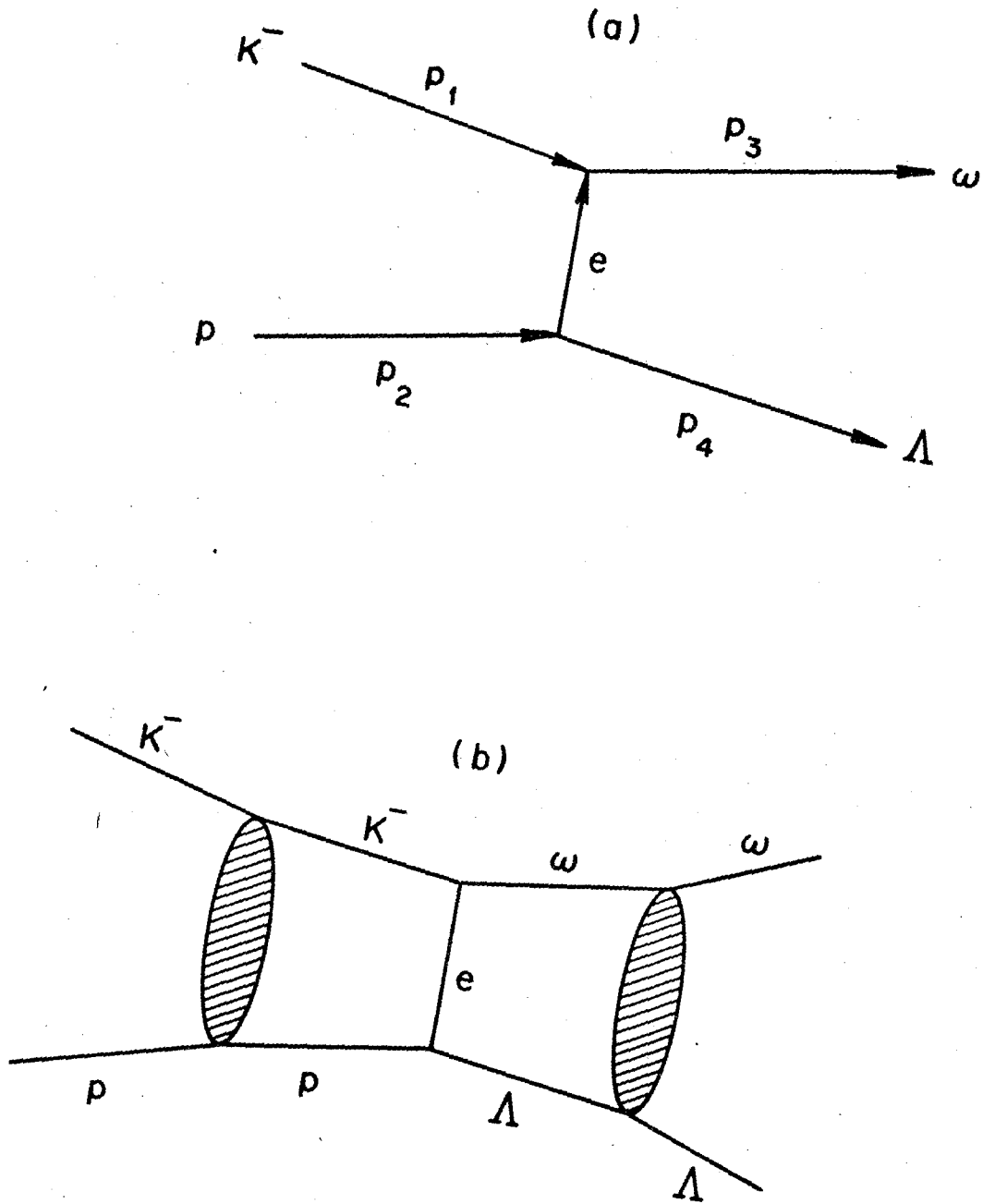


Fig. 1

MUB-9989

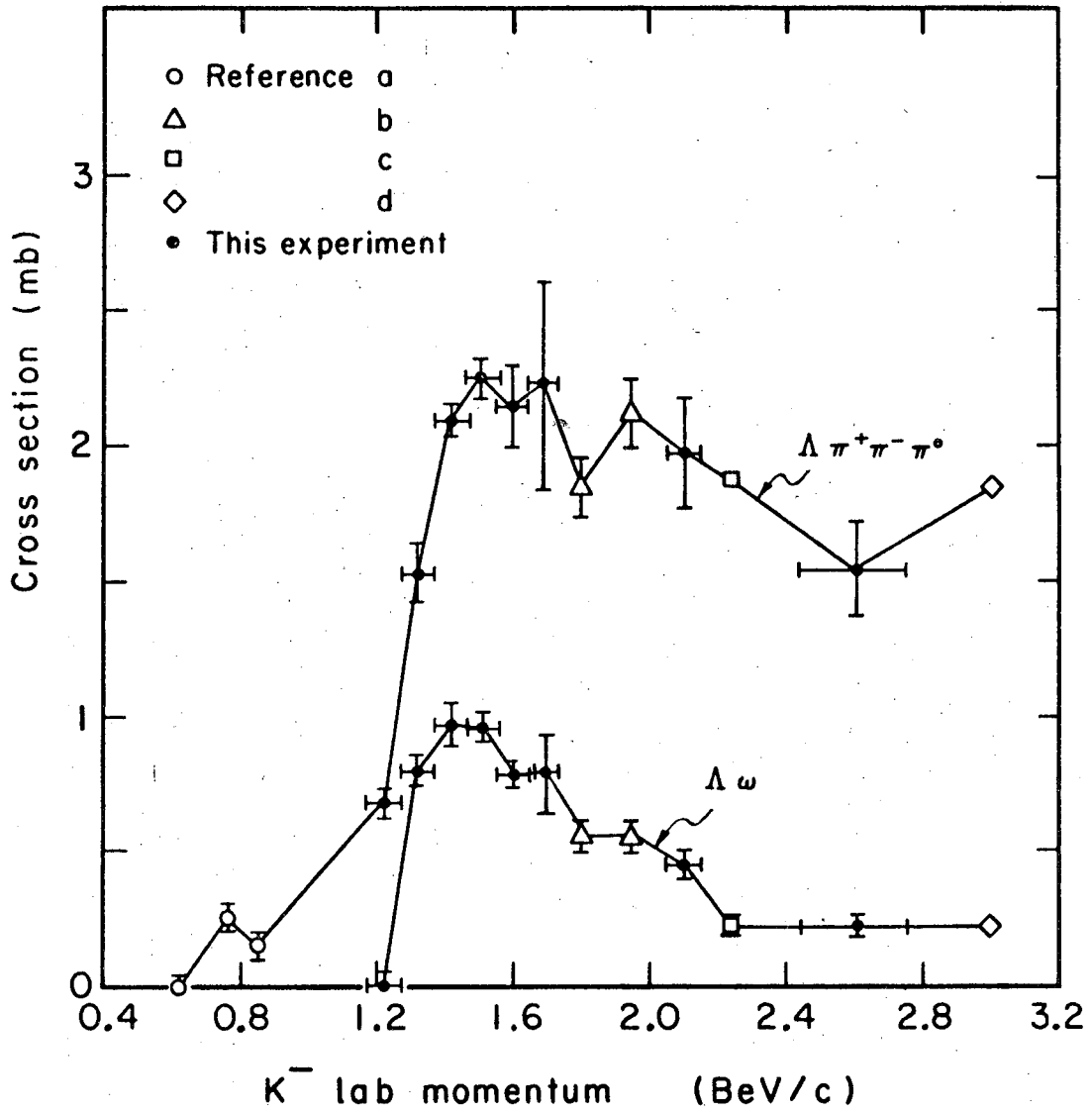


Fig. 2

MU-37233

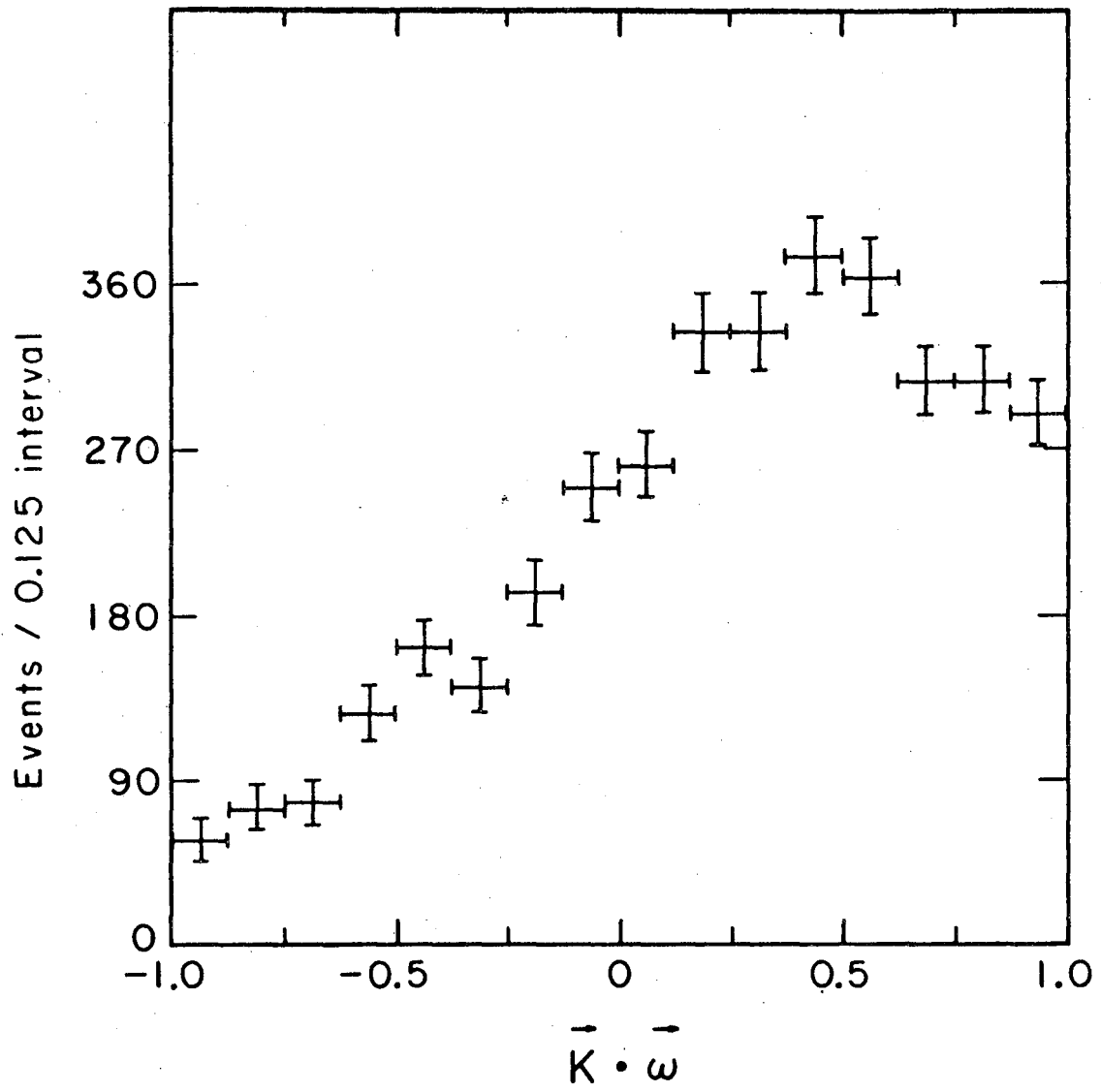


Fig. 3

MU.37227

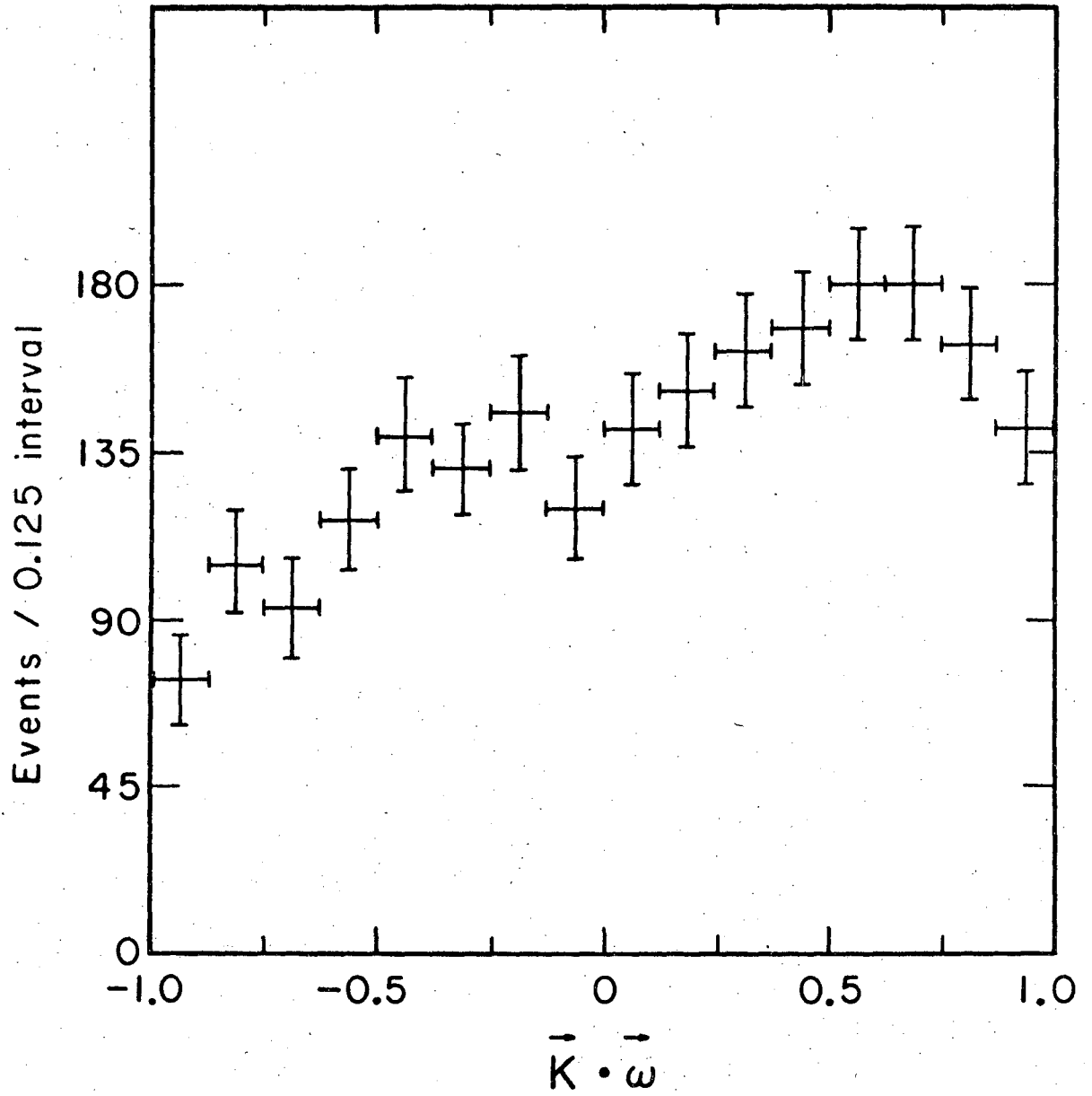


Fig. 4

MU-37228



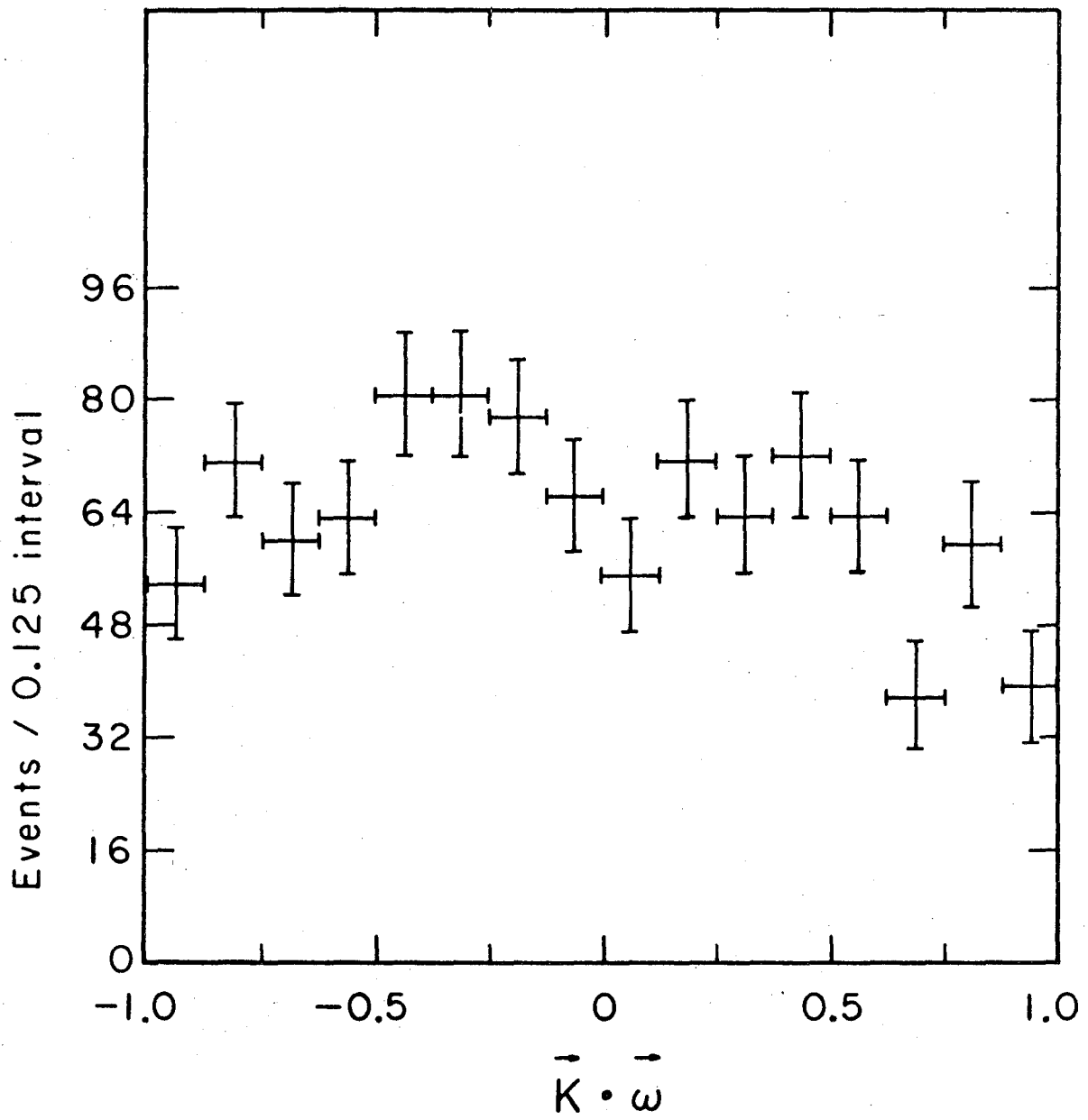


Fig. 5

MU-37230

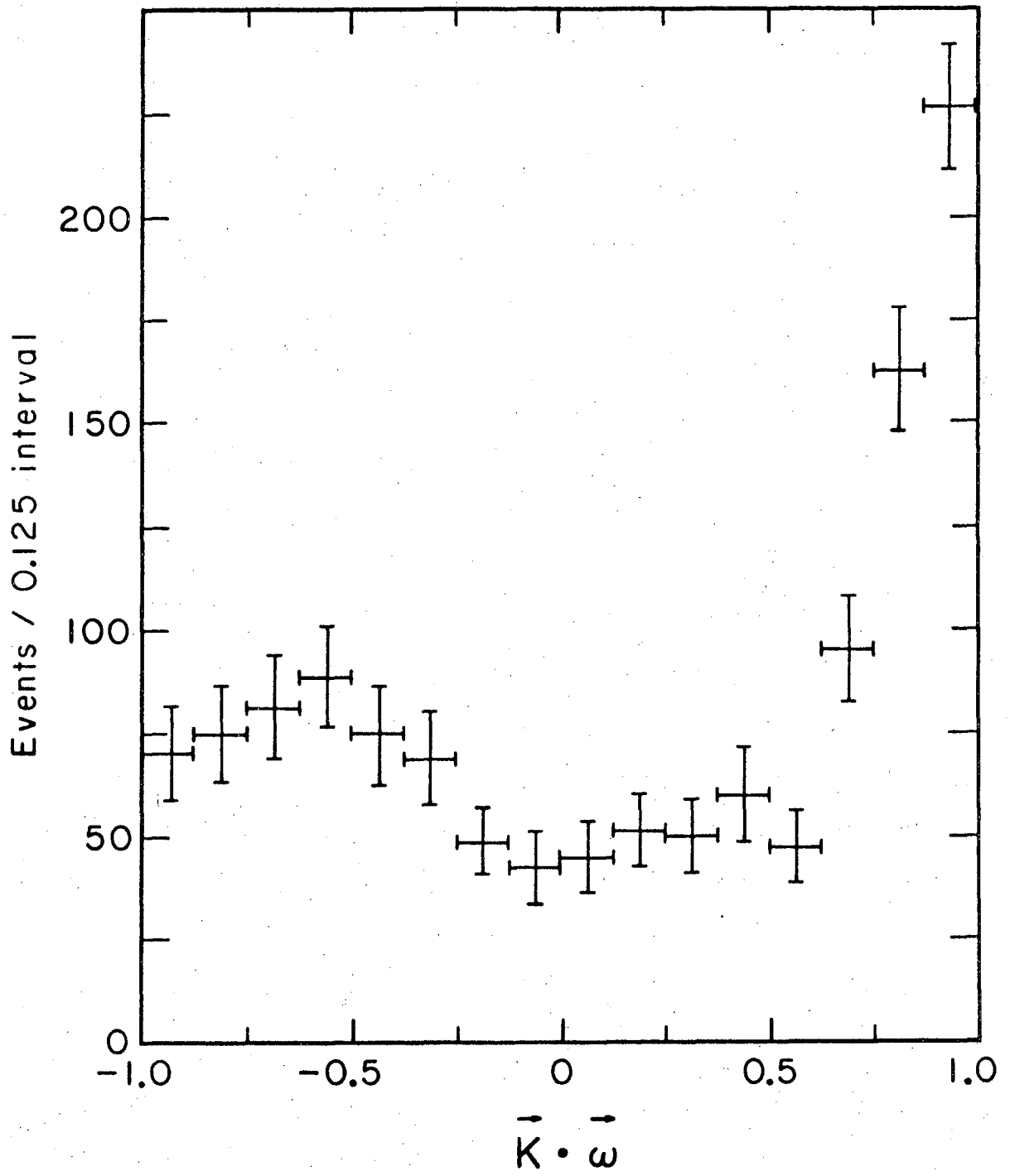


Fig. 6

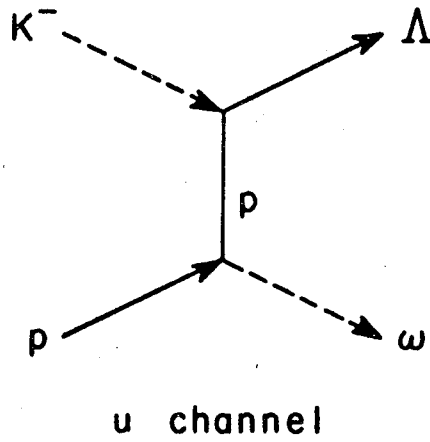
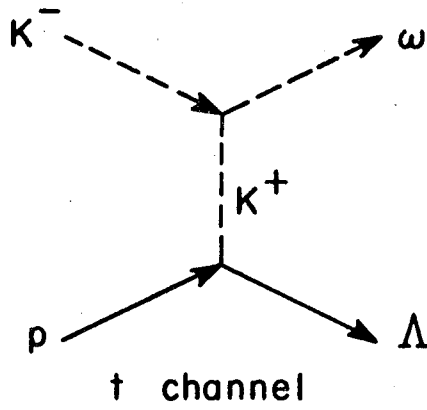
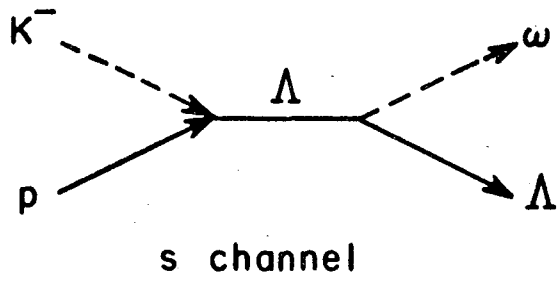


Fig. 7

MUB-8404

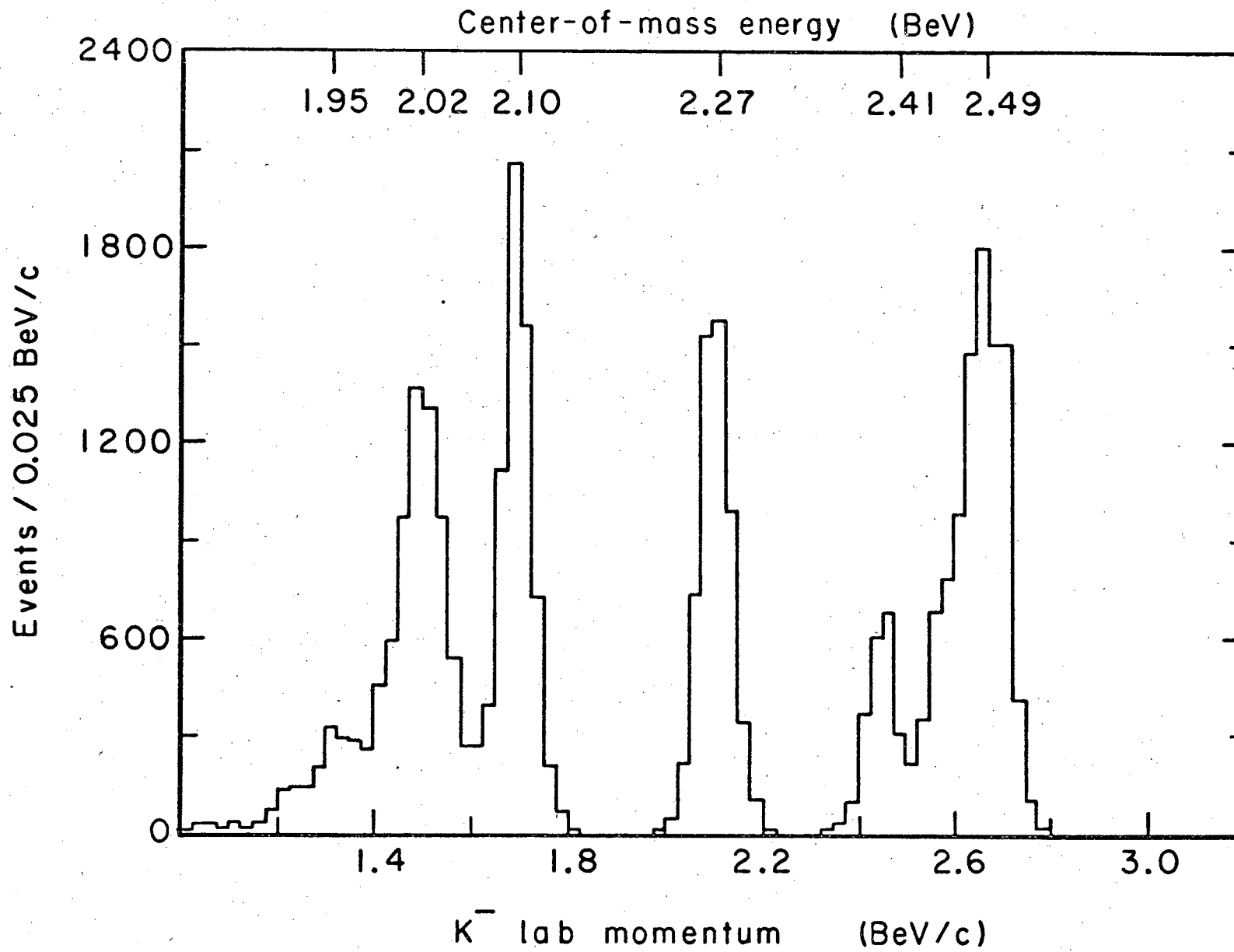


Fig. 8

MU-37237

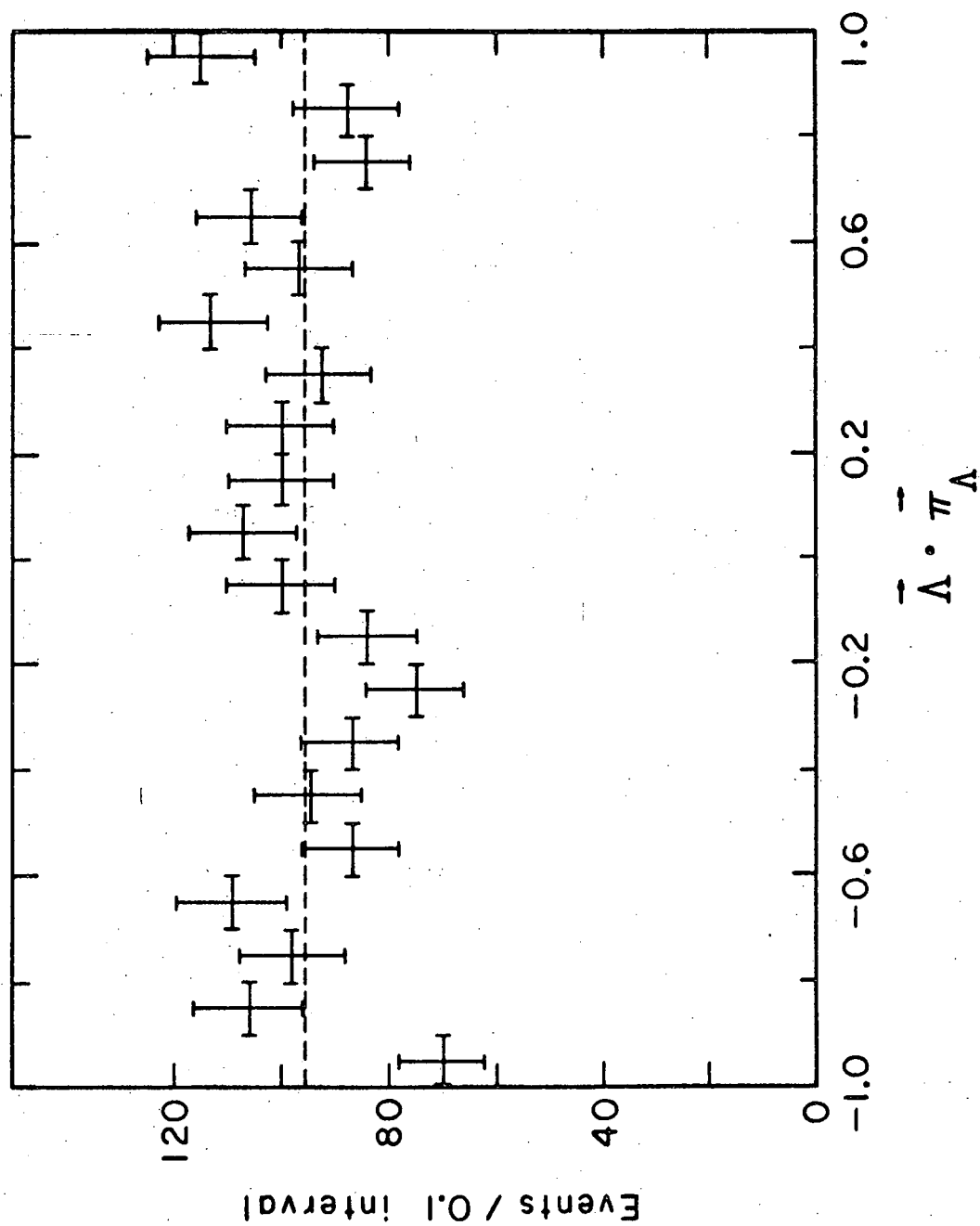


Fig. 9

MU-37226

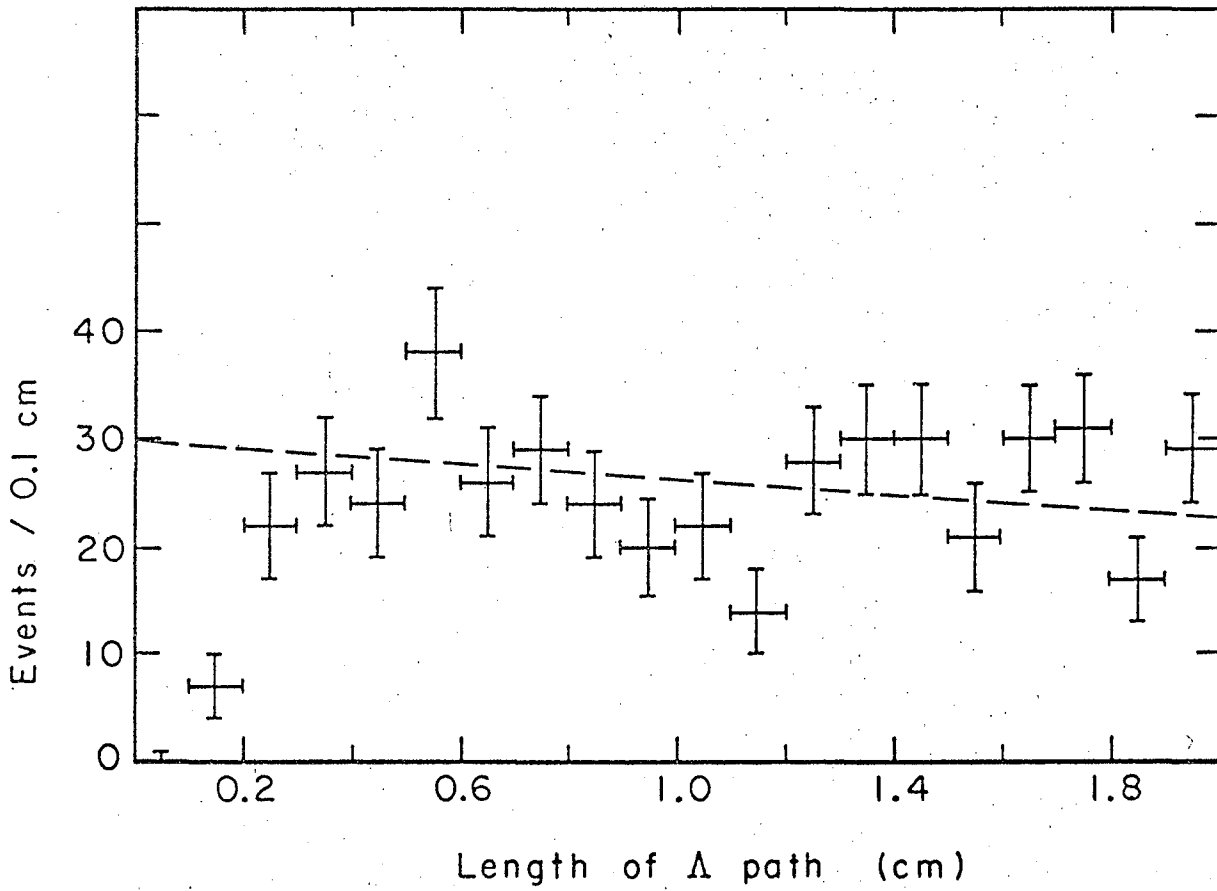


Fig. 10

MU.37231

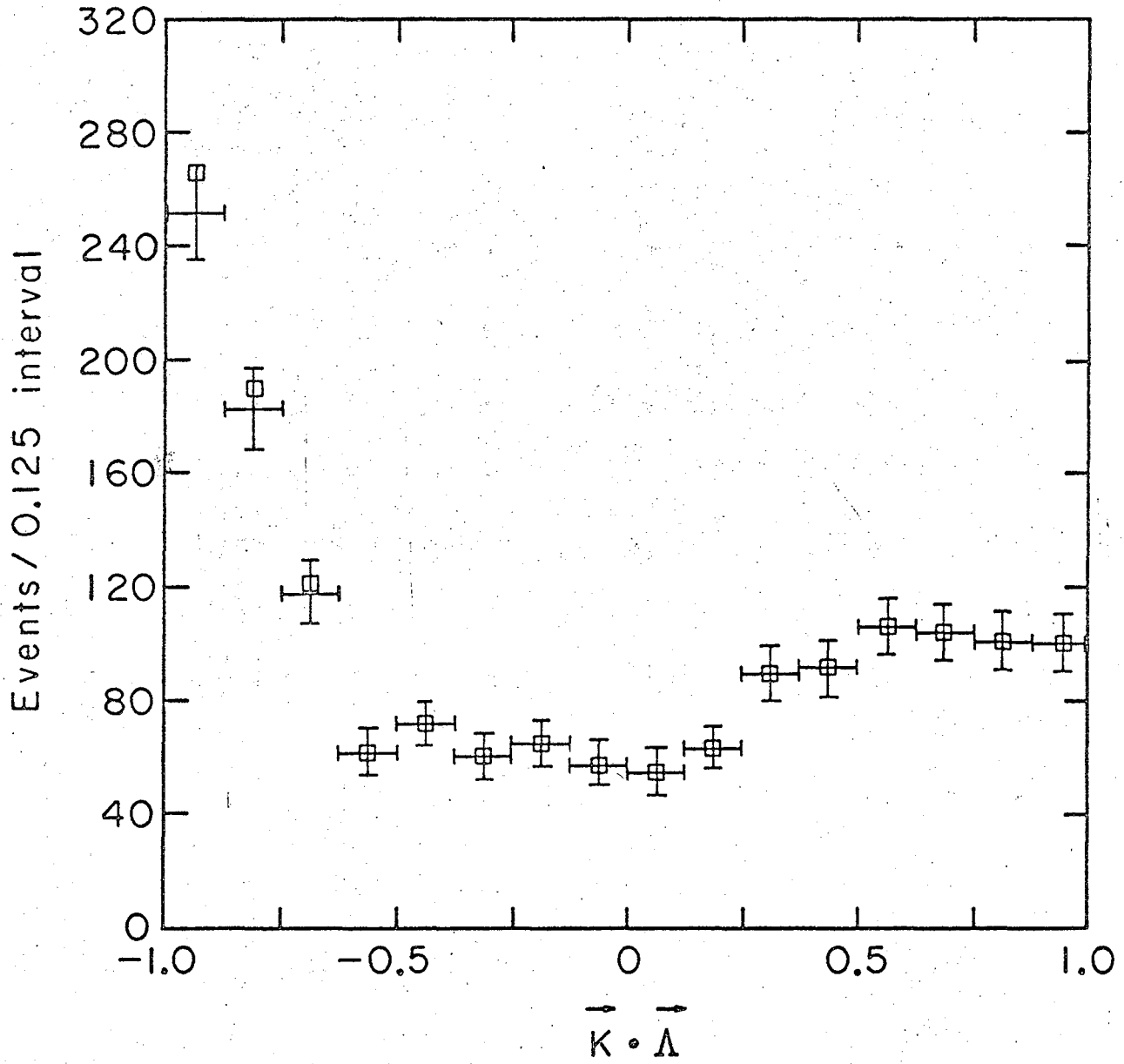


Fig. 11

MU-37235

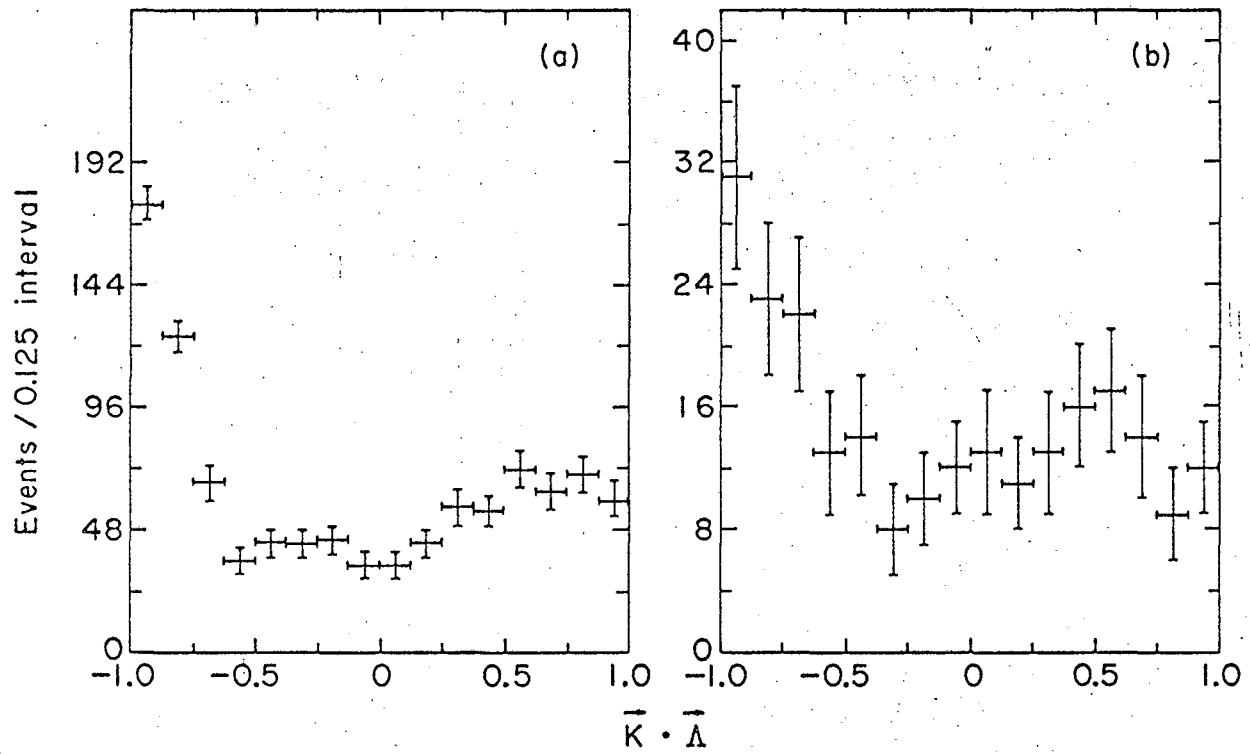


Fig. 12

MU-37234



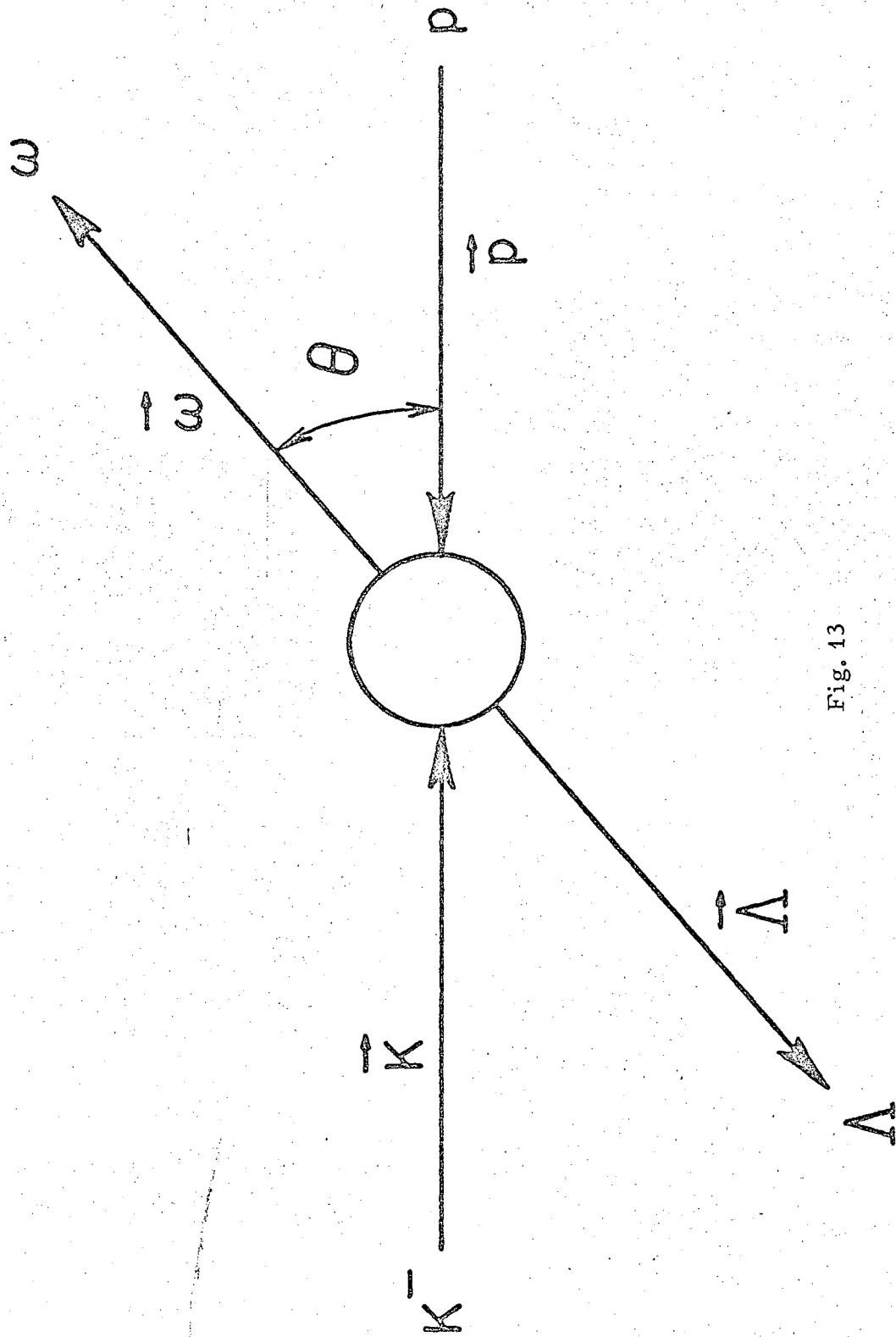


Fig. 13

MU-37232

MU-37236

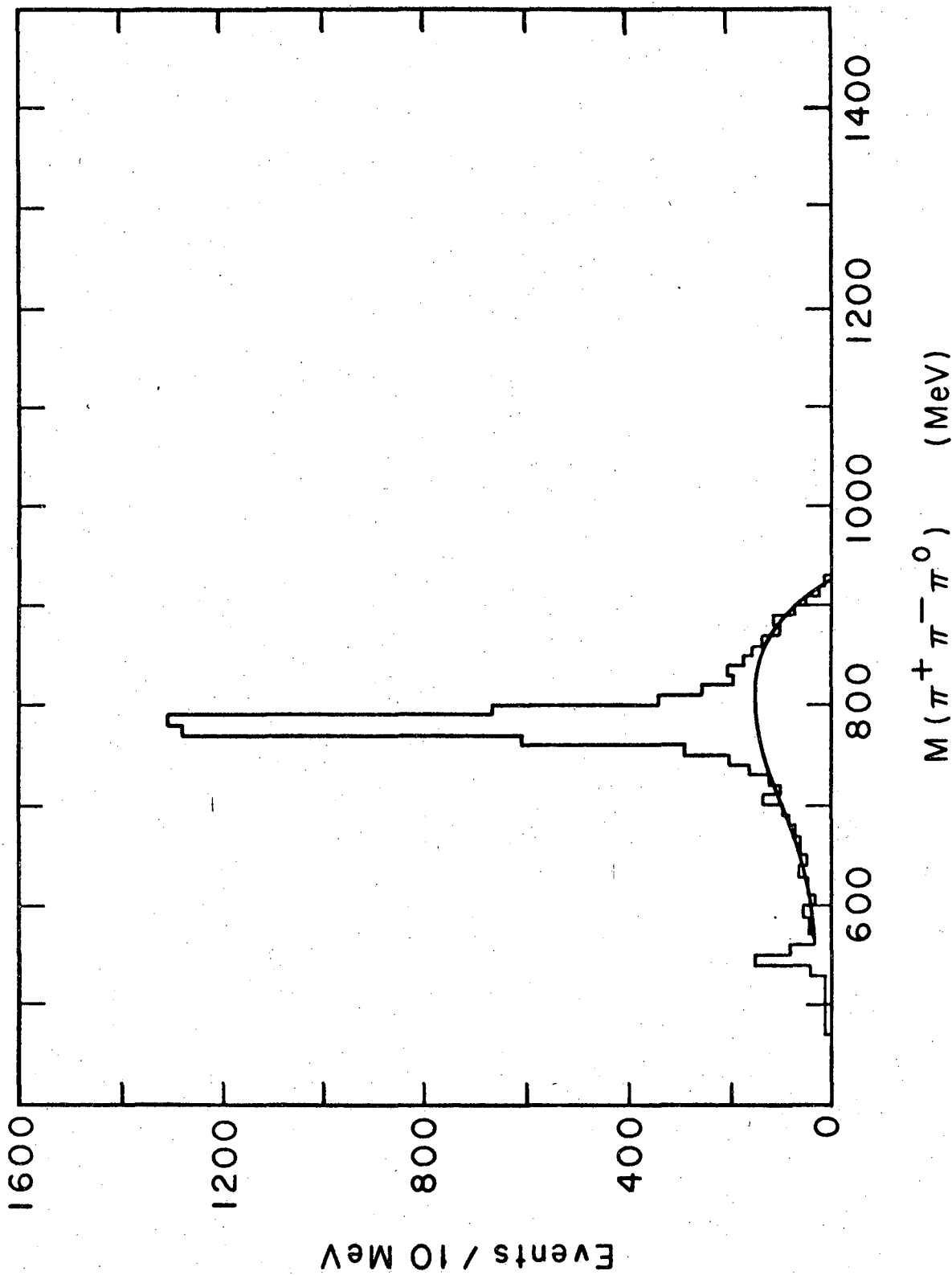


Fig. 14

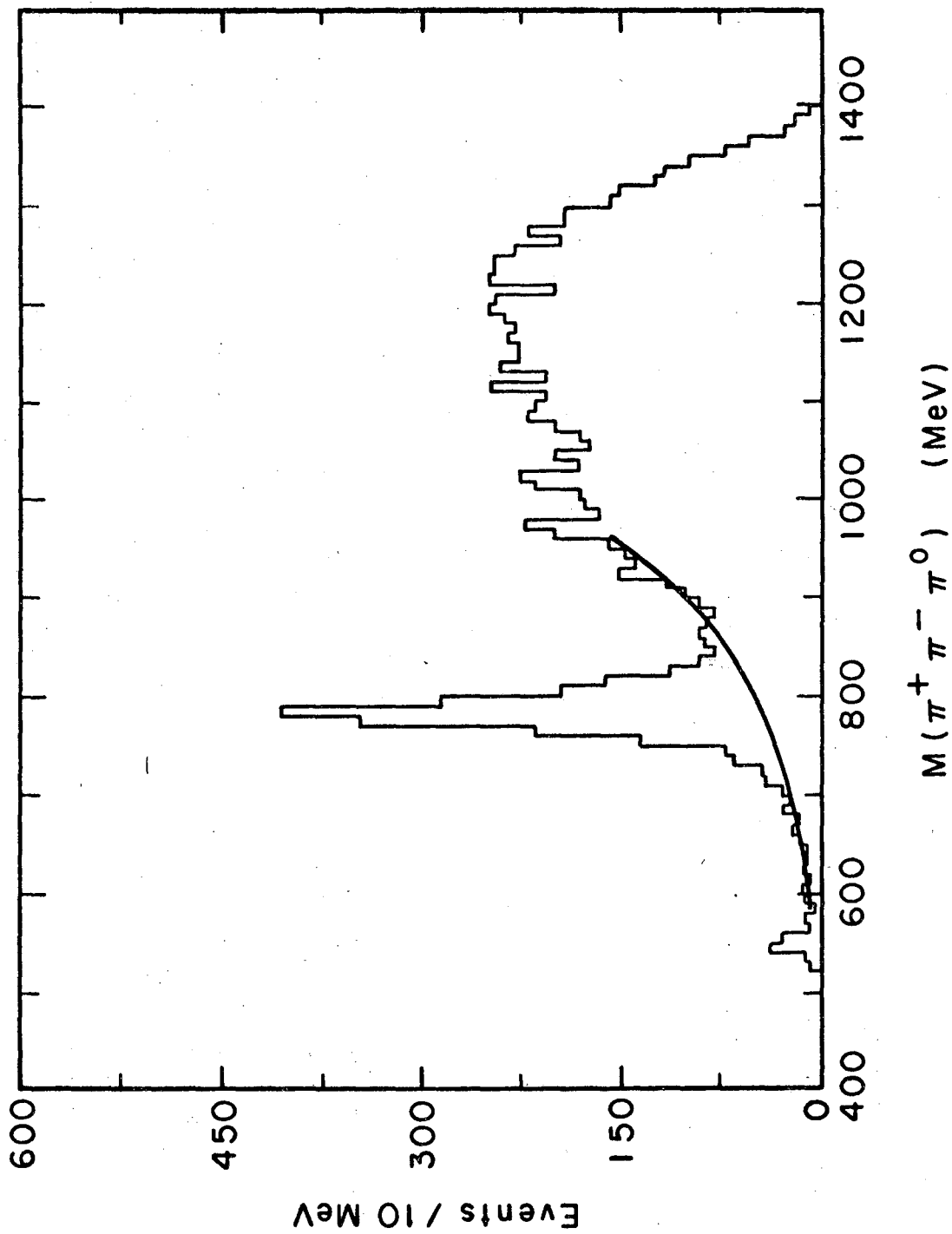


Fig. 15

MU-37240

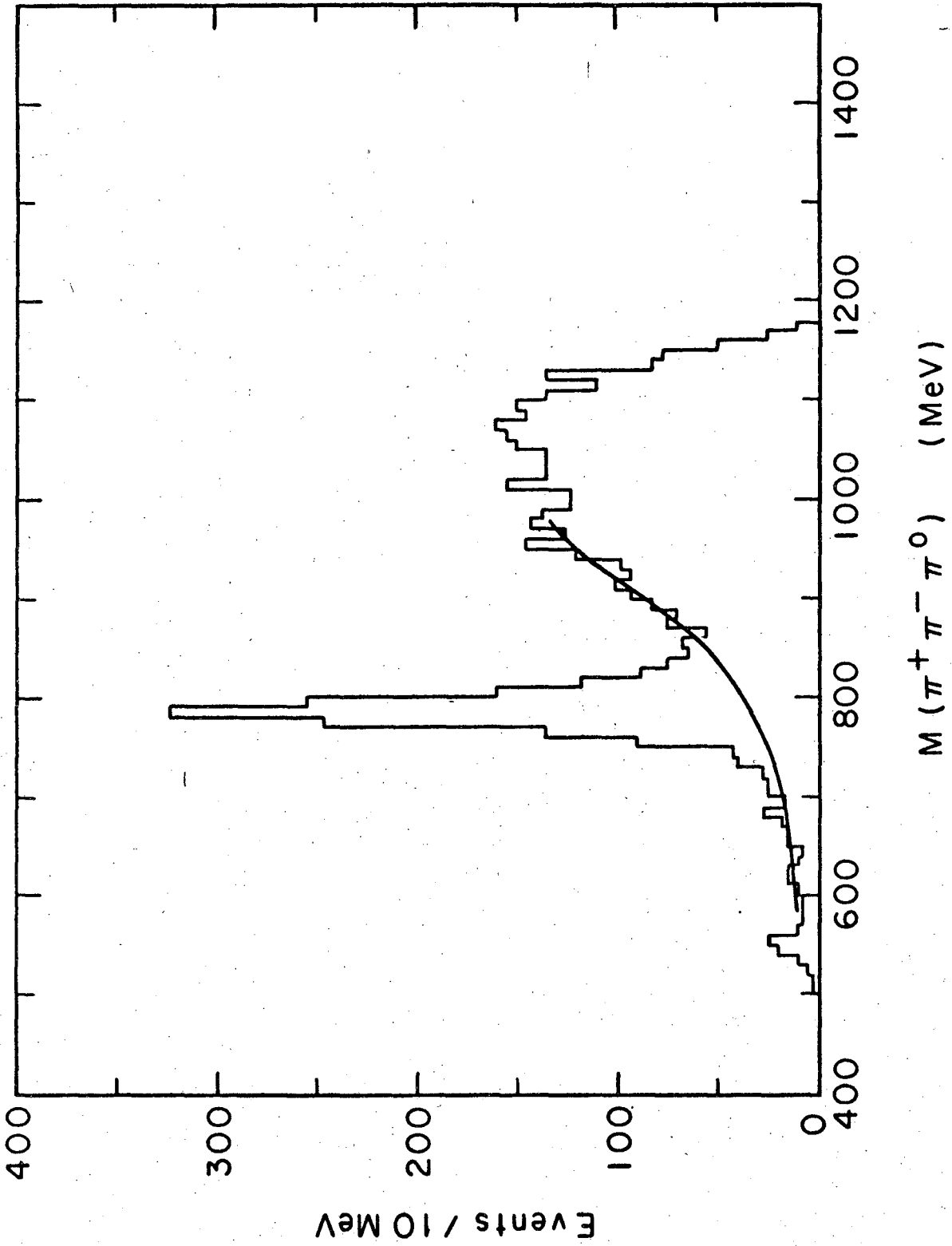


Fig. 16

MU 37239

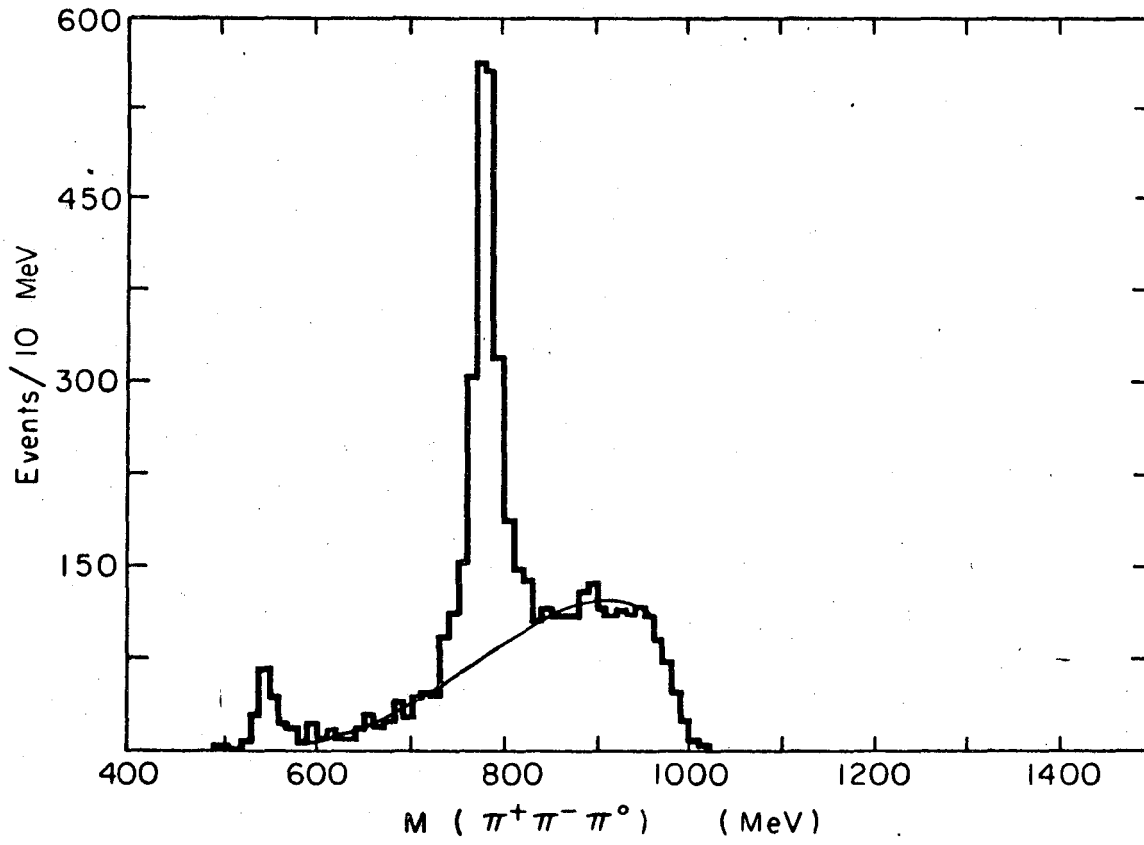


Fig. 17

MUB-12104



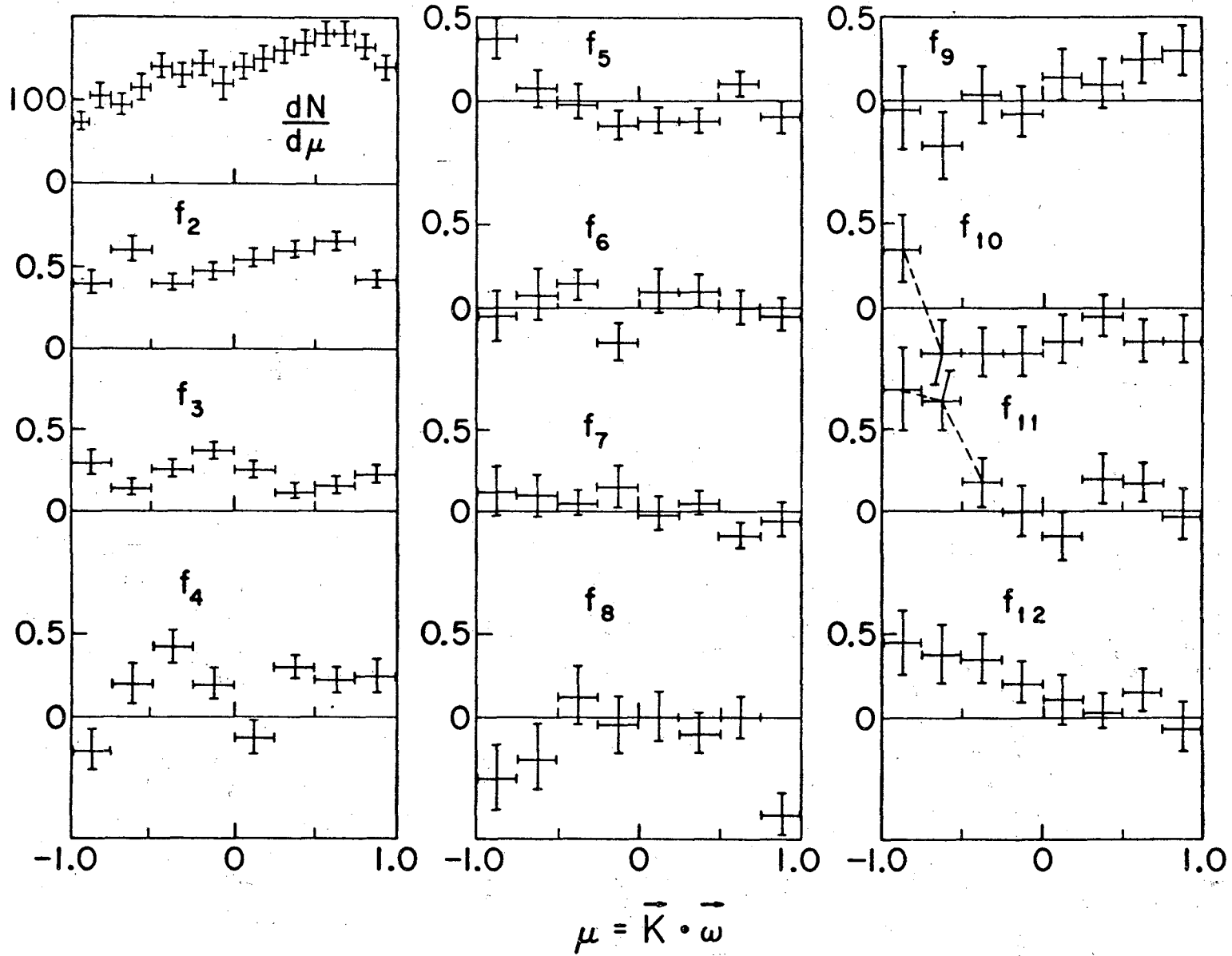
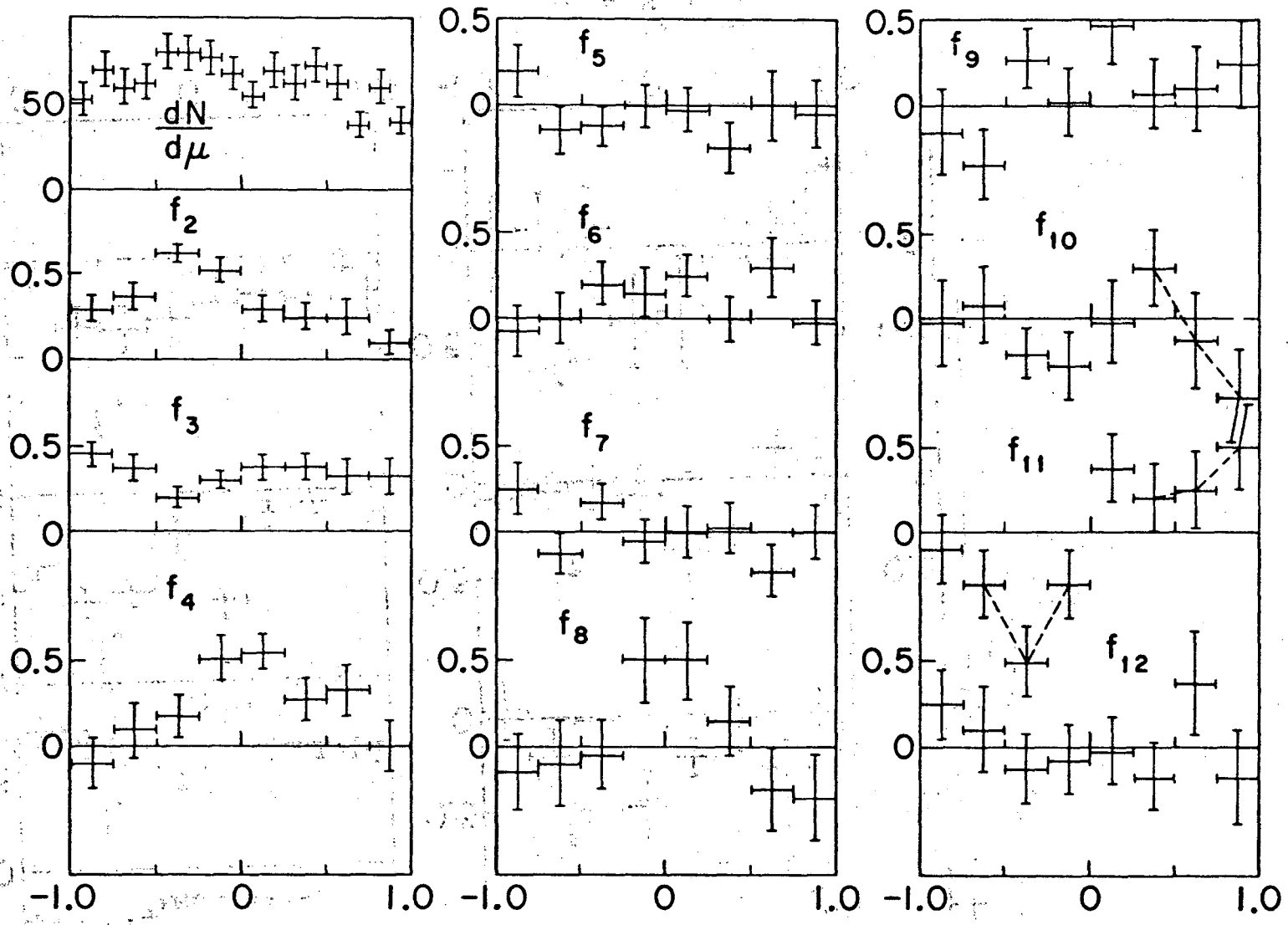


Fig. 19

MU-37242



$$\mu = \vec{K} \cdot \vec{\omega}$$

Fig. 20

MU-37243



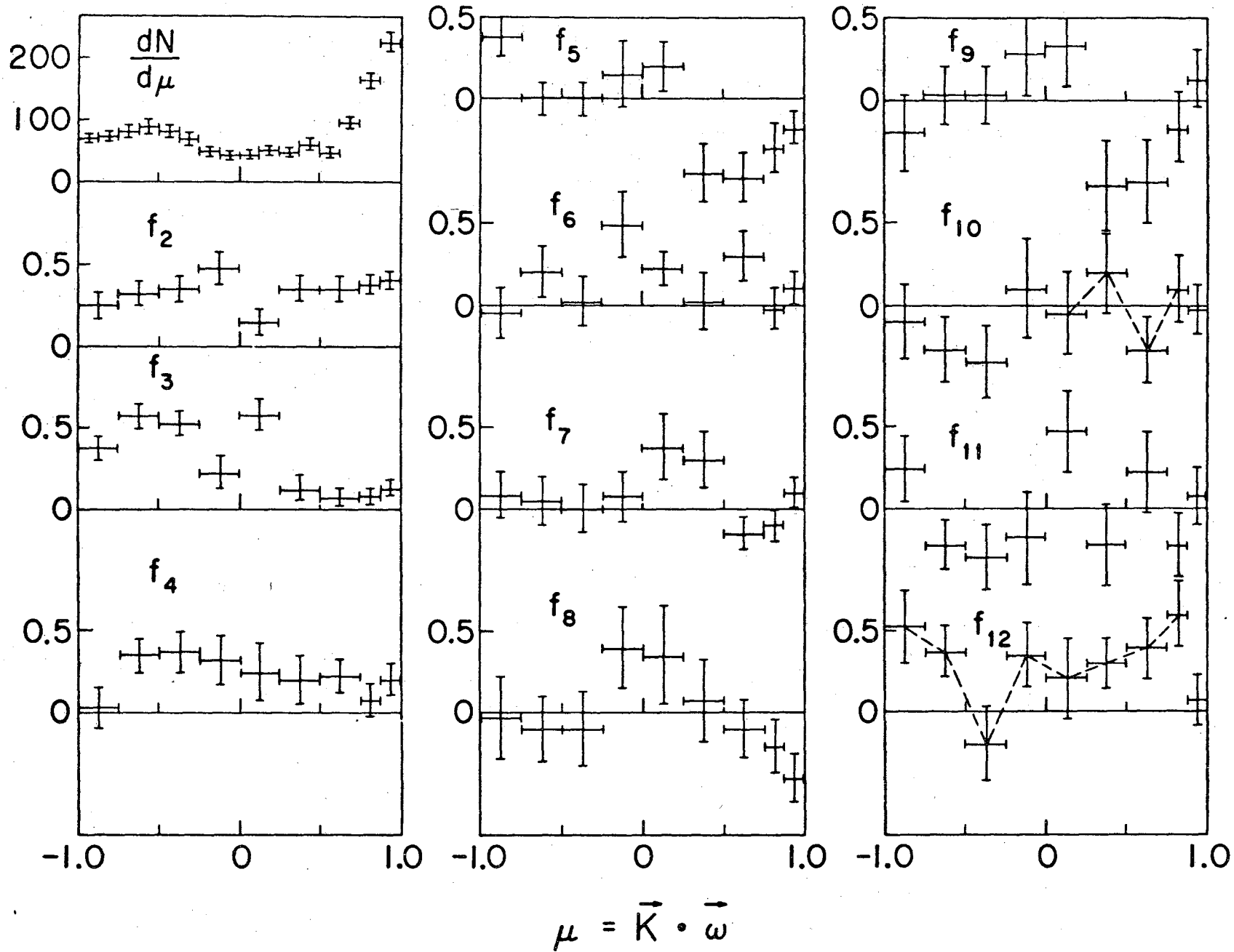


Fig. 21

MU-37244

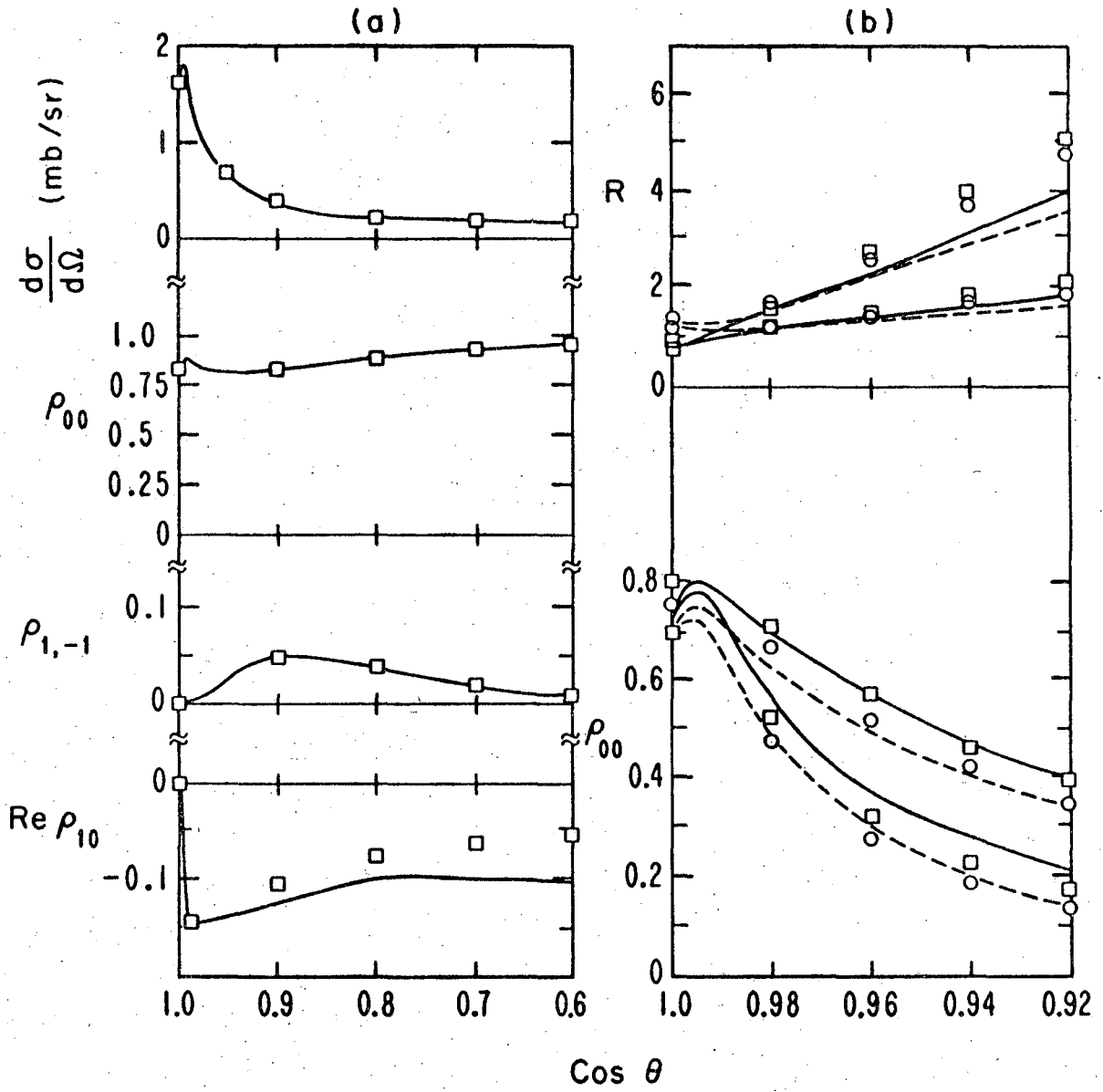
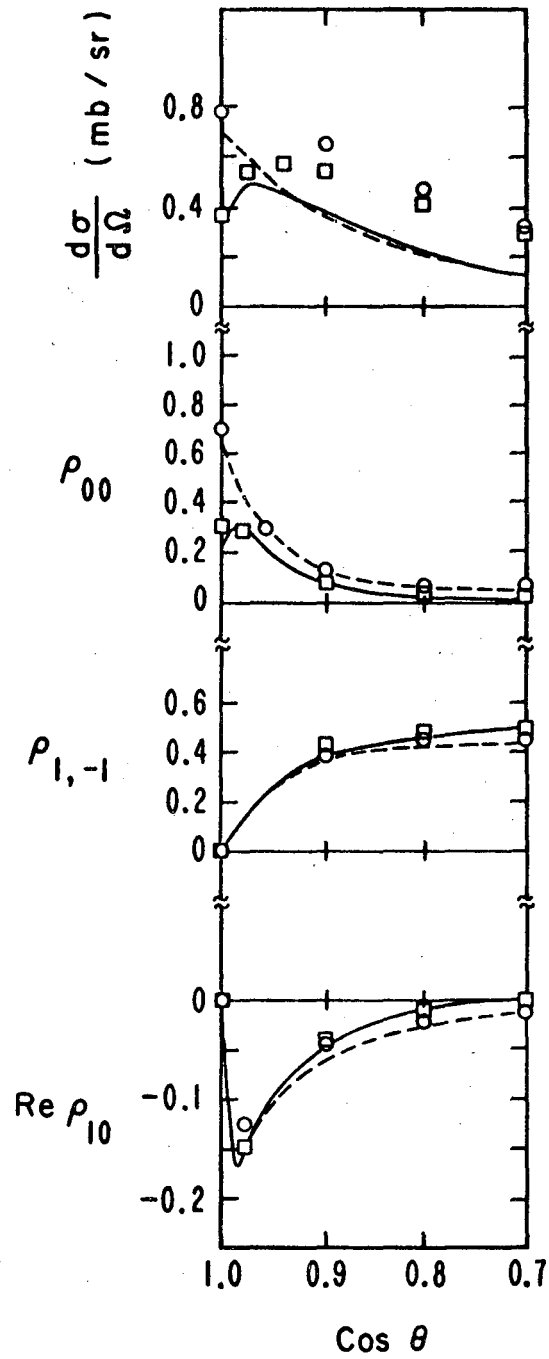


Fig. 22

MUB-9990



MUB-9991

Fig. 23

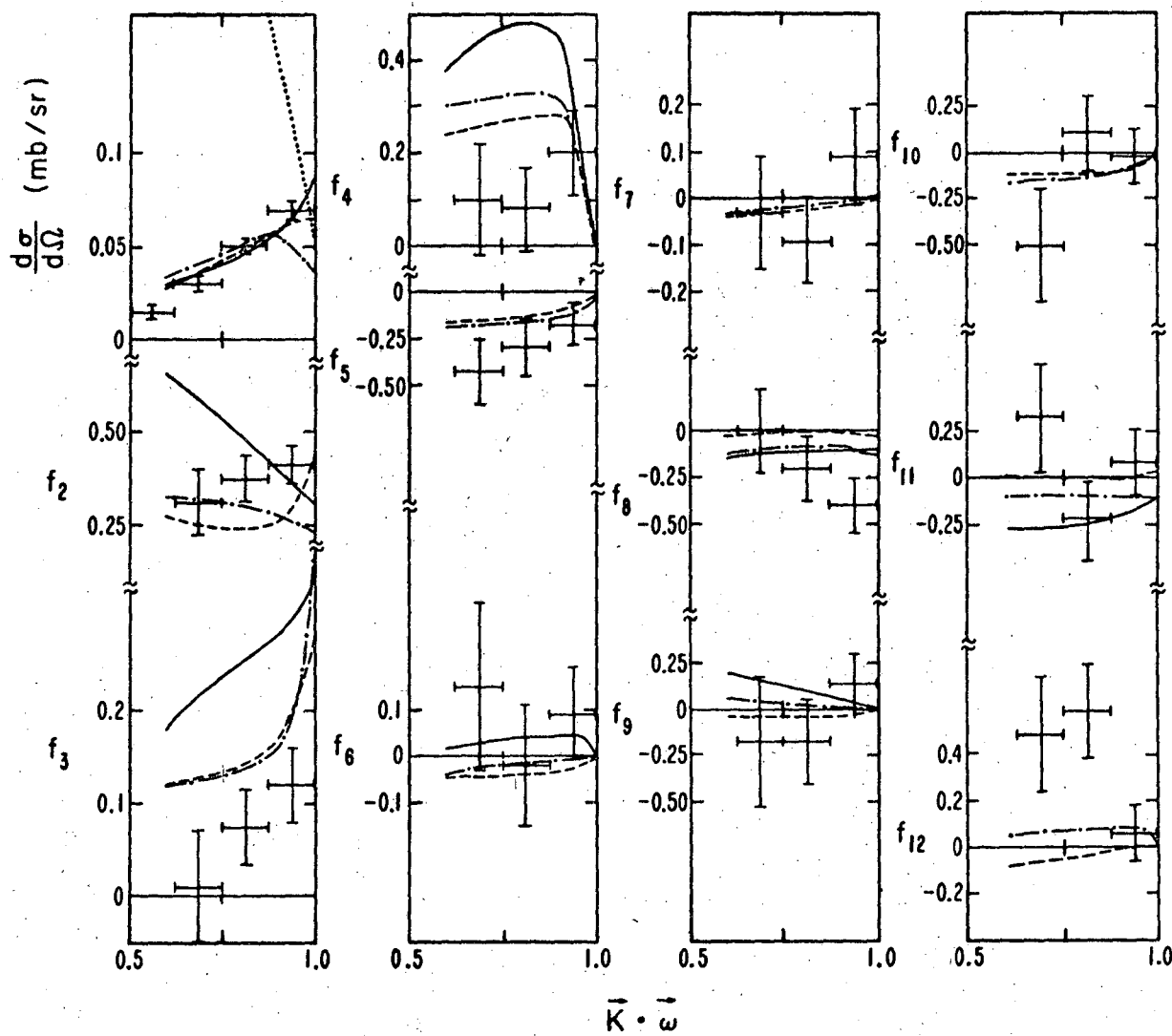


Fig. 24

MUB-9992

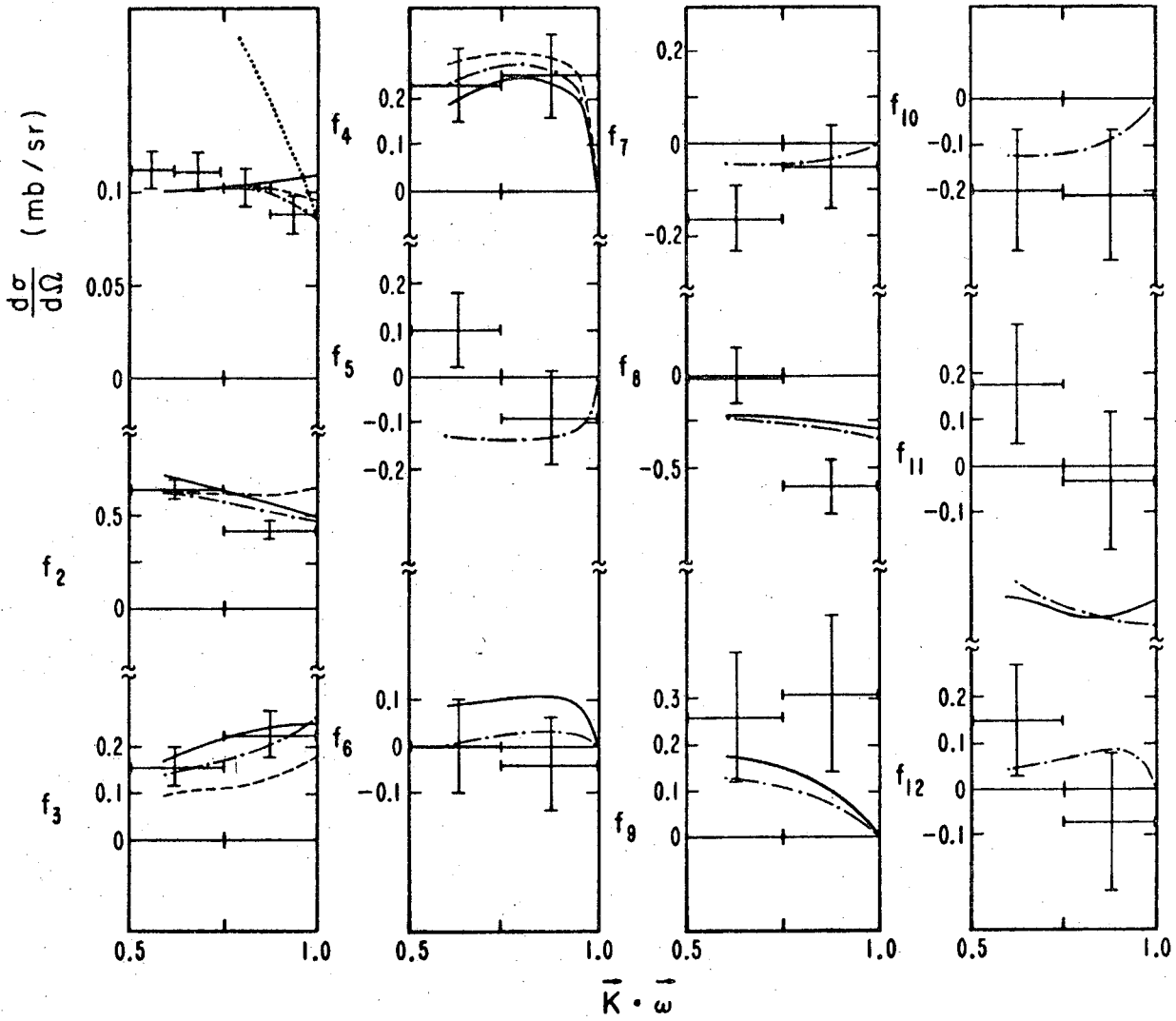


Fig. 25

MUB-9993

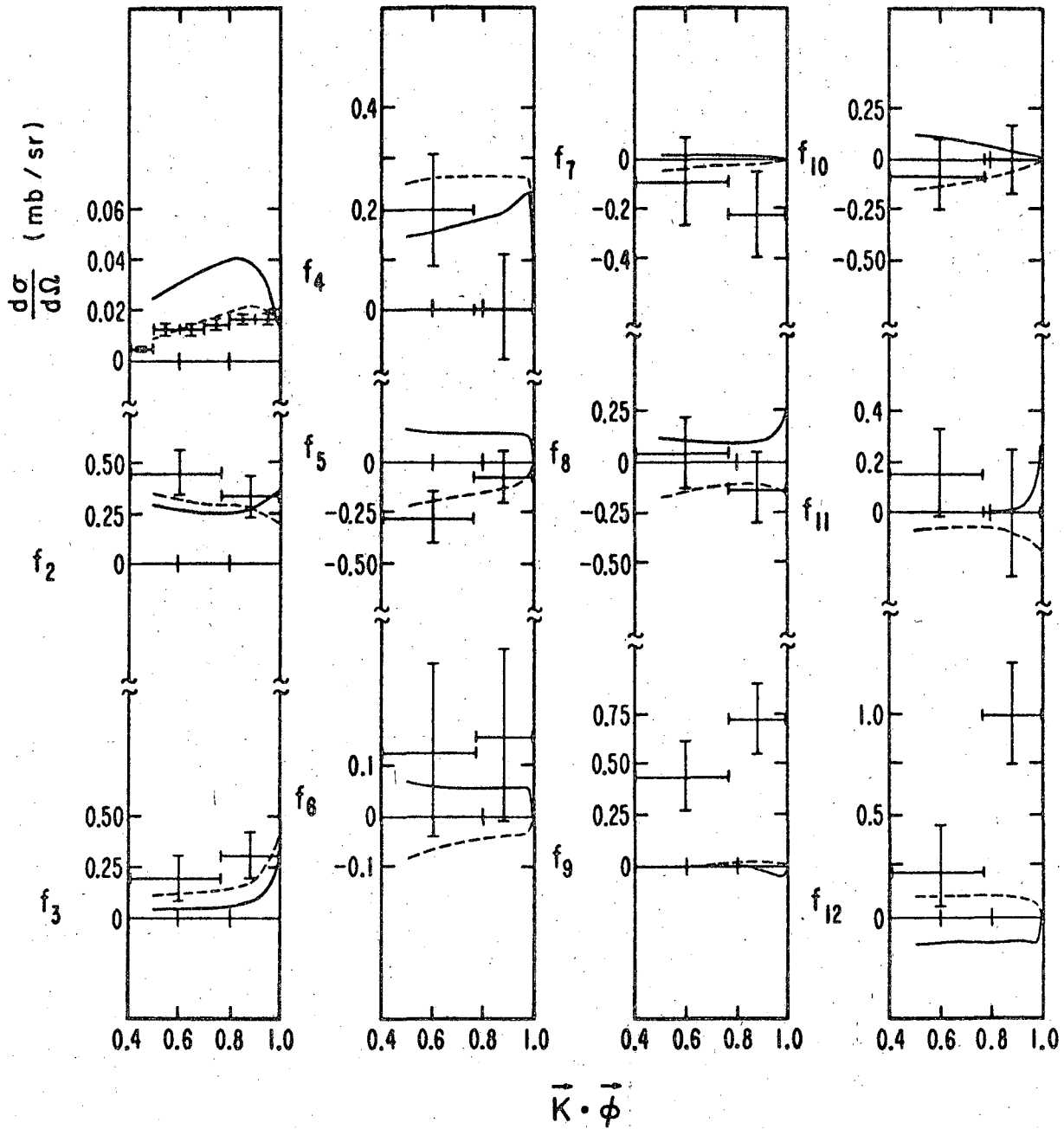


Fig. 26

MUB-9994

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