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K_{l3} FORM FACTORS*

V. S. Mathur,[†] L. K. Pandit,[‡] and R. E. Marshak Department of Physics and Astronomy, University of Rochester, Rochester, New York (Received 28 March 1966)

Great interest attaches to the prediction of the form factors $F_+(s)$ and $F_-(s)$ for the K_{l3} decays since experimentalists are finding quite a variety of energy dependences and ξ values for the neutral and charged decay modes.¹ Recent calculations of these form factors on the basis of the algebra of currents² have either related³ $[F_+(m_K^2) + F_-(m_K^2)]$ to $K_{\mu 2}$ decay, or have predicted a value of $F_+(0)$ in terms of the K^* , ρ , and κ meson widths. In this note, we apply the current algebra and dispersion techniques⁴ to a direct calculation of $F_+(s)$ and $F_-(s)$, and fix the absolute scale as well as the energy dependence.

We introduce the K_{l3}^+ form factors as follows⁵:

$$\langle \pi^{0}(q') | (V_{\mu}(0))_{1}^{3} | K^{+}(q) \rangle$$

= $(4q_{0}q_{0}'V^{2})^{-1/2}$
 $\times [F_{\mu}(s)(q+q')_{\mu} + F_{\mu}(s)(q-q')_{\mu}], \qquad (1)$

where $s \equiv -(q - q')^2$. We next define

$$F_{1}(s) = F_{+}(s),$$
 (2a)

$$F_0(s) = F_{(s)} + [(M_K^2 - M_\pi^2)/s]F_{(s)},$$
 (2b)

where now $F_1(s)$ receives contributions only from the $J = 1^-$ states and $F_0(s)$ only from $J = 0^+$ states. These form factors are supposed to satisfy unsubtracted dispersion relations:

$$F_{1}(s) = \frac{1}{\pi} \int_{(M_{\pi} + M_{K})^{2}}^{\infty} ds' \frac{\operatorname{Im} F_{1}(s')}{s' - s - i\epsilon}, \qquad (3a)$$

$$F_0(s) = \frac{M_K^2 - M_{\pi}^2}{s} F_1(0)$$

$$+\frac{1}{\pi}\int_{(M_{\pi}+M_{K})^{2}}^{\infty} ds' \frac{\mathrm{Im} F_{0}(s')}{s'-s-i\epsilon}.$$
 (3b)

Following the standard method⁶ for calculating the absorptive parts and using the $K^*(891)$ pole for $F_1(s)$ and $\kappa(725)$ for $F_0(s)$, one obtains

$$F_{+}(s) = G_{K*}G_{K*K\pi}/(M_{K*}^{2}-s),$$
 (4a)

$$F_{-}(s) = -\frac{M_{K}^{2} - M_{\pi}^{2}}{M_{K}^{2}} \frac{G_{K} + G_{K} + K\pi}{M_{K}^{2} - s} + \frac{G_{\kappa} G_{\kappa\kappa\pi}}{M_{\kappa}^{2} - s}, \quad (4b)$$

with the following definitions⁷:

$$\langle 0 | (V_{\mu}(0))_{1}^{3} | K^{*+}(p, \epsilon) \rangle = [G_{K^{*}}/(2p_{0}V)^{1/2}]\epsilon_{\mu}, \quad (5a)$$
$$\langle K^{*+}(p, \epsilon) | j_{\pi^{0}}(0) | K^{+}(q) \rangle |_{(p-q)^{2} = 0}$$

$$= -\frac{2G_{K^{*K}\pi}}{(4p_{0}q_{0}V^{2})^{1/2}}(\epsilon \cdot q);$$
(5b)

$$\langle 0 | (V_{\mu}(0))_{1}^{3} | \kappa^{+}(p) \rangle = [G\kappa/(2p_{0}V)^{1/2}]p_{\mu},$$
 (6a)

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$$\langle \kappa^{+}(p) | j_{\pi^{0}}(0) | K^{+}(q) \rangle | (p-q)^{2} = 0$$

$$= -G_{\kappa K \pi} / (4p_{0}q_{0}V^{2})^{1/2}.$$
(6b)

Up to this point, we have not gone beyond the well-known dispersion-theoretic treatment. But now the algebra of currents, through its nonlinear structure, enables us to relate the constants G_{K*} and G_{κ} to the coupling constants $G_{K*K\pi}$ and $G_{\kappa K\pi}$, respectively, and thereby to fix the scale and <u>a fortiori</u> the energy dependence of the form factors. Following the standard reduction technique and using partial conservation of axial-vector current (PCAC) for both the strangeness-changing as well as the strangeness-conserving axial-vector currents, we obtain,⁸ in the limit of zero four-momenta for the K^+ and π^0 ,

$$\langle K^{+}(q) | j_{\pi^{0}}(0) | K^{*+}(p,\epsilon) \rangle |_{(p-q)^{2}=0} = -\frac{M_{\pi^{2}M_{K}^{2}}}{C_{\pi}C_{K}} \frac{(p-q)_{\mu}}{(2q_{0}V)^{1/2}} \langle 0 | [B_{1}^{3}(0), (P_{\mu}(0))_{1}^{1} - (P_{\mu}(0))_{2}^{2}] | K^{*+}(p,\epsilon) \rangle, \quad (7)$$

where⁹

$$B_{1}^{3}(0) = \int d^{4}x \ \theta(x_{0}) \partial_{\nu} (P_{\nu}(x))_{1}^{3}$$
(8)

is the strangeness-changing axial "charge." Using the equal-time commutation relation

$$[B_{1}^{3}(0), (P_{\mu}(0))_{1}^{1} - (P_{\mu}(0))_{2}^{2}] = (V_{\mu}(0))_{1}^{3}, \qquad (9)$$

and Eqs. (5a) and (5b), we obtain the desired relation

$$G_{K*} = 2 \frac{C_{\pi} C_{K}}{M_{\pi}^{2} M_{K}^{2}} G_{K*K\pi}.$$
 (10)

Similarly,

$$G_{\kappa} = 2 \frac{C_{\pi} C_{K}}{M_{\pi}^{2} M_{K}^{2}} \frac{G_{\kappa K \pi}}{M_{\kappa}^{2}}.$$
 (11)

Equations (4a) and (4b) together with Eqs. (10) and (11) determine completely the form factors in terms of the coupling constants $G_{K * K \pi}$ and $G_{\kappa K \pi}$, which in turn are given by the K^* and κ decay widths.

Using the experimental K^* width $\Gamma(K^*) = 50$ MeV, and the values^{5,7} $C_{\pi}/M_{\pi}^2 = -0.16$ BeV, $C_K/M_K^2 = 0.14$ BeV, we obtain

$$F_+(0) = -0.60. \tag{12}$$

A comparison with the SU(3)-symmetry value $F_+(0) = -1/\sqrt{2} = -0.71$, which is expected to be correct to second order in SU(3) breaking in view of the Ademollo-Gatto theorem, ¹⁰ shows that the K^* dominance model for $F_+(s)$ is indeed very reasonable. If we use the form factor $F_+(s) = -0.6/(M_K *^2 - s)$ to calculate $\Gamma(K_{e3})$, we find for the "renormalized" Cabibbo¹¹ angle $\theta_V^M = 0.238$, to be compared with the "bare"

value $\theta = 0.222 \pm 0.006$. In vew of the approximations made,¹² this result can also be regarded as compatible with the Ademollo-Gatto theorem.¹³

It is of interest to compare our result of Eq. (12) with a previous evaluation of $F_{+}(0)$ on the basis of the current algebra.³ In the paper of Mathur, Okubo, and Pandit, ${}^{3}F_{+}(0)$ was determined in terms of the single-particle intermediate states K^*, κ, ρ with the result that $F_{+}(0) = -0.77$, in reasonable agreement with the present evaluation. In the present work, of course, we do not encounter the ρ and the κ states for $F_{+}(s)$. The difference arises from the fact that one is really "dispersing" in different variables in the two methods, implying a "bootstrap" relation between the two. A detailed discussion of "bootstrapping" within the framework of the algebra of currents will be taken up elsewhere.

To determine $F_{-}(0)$, we must know the κ width, assuming that it exists¹⁴; if we choose the published value $\Gamma(\kappa) = 10$ MeV, we obtain

$$F_{-}(0) = 0.12,$$
 (13)

and hence

$$\xi \equiv F_{-}(0)/F_{+}(0) = -0.20. \tag{14}$$

It should be noted that the ratio ξ is independent of C_{π} and C_K and so does not suffer from any uncertainties in their numerical estimates. The parameter ξ is predicted to be negative and can only be reduced to 0 if $\Gamma(\kappa) \simeq 35$ MeV; a value of $\xi \simeq 1.2$ as suggested by the recent experiment of Carpenter et al.¹⁵ would require an unreasonably large value of $\Gamma(\kappa)$, and would imply a substantially larger SU(3) breaking than is indicated by other evidence.¹⁶ It will be of great interest to have an accurate measurement of ξ .

Our results depend on the use of an $I = \frac{1}{2}$ strangeness-changing vector current density. The $I = \frac{1}{2}$ hypothesis¹⁷ for the strangeness-changing weak hadron current is naturally taken over in the algebra of the octet of currents. In view of the many successes¹⁸ of this approach, it is hard to believe the recent experimental claims^{15,1} for a different energy dependence of the form factors and a different parameter ξ for K_{I3}^{0} and K_{I3}^{+} .

The encouraging result obtained for $F_+(0)$ in terms of the observed K^* width suggests applying the same technique to the evaluation of the widths of the vector mesons φ and ω . An earlier calculation¹⁹ has already been made for the ρ -meson width using the conserved isospin current. In the present case, we make use of the conserved hypercharge current²⁰ to which the φ and ω are coupled. Thus, we define the matrix element of the hypercharge current

$$\langle K^{+}(q') | (V_{\mu}(0))_{3}^{3} | K^{+}(q) \rangle$$

= $(4q_{0}q_{0}'V^{2})^{-1/2}(q+q')_{\mu}H_{+}(s).$ (15)

From the conservation of $(V_{\mu}(x))_{3}^{3}$, we have

$$H_{+}(0) = 1. \tag{16}$$

Following the method already discussed we obtain

$$H_{+}(s) = \frac{4}{3} \frac{C_{K}^{2}}{M_{K}^{4}} \left[\frac{G_{\varphi K \overline{K}}}{M_{\varphi}^{2} - s} + \frac{G_{\omega K \overline{K}}}{M_{\omega}^{2} - s} \right].$$
(17)

For the purpose of a numerical estimate based on this sum rule, we may use the $\omega -\varphi$ mixing model²¹ of broken SU(3) to obtain $G_{\omega K \overline{K}} = G_{\varphi K \overline{K}} / \sqrt{2}$. Then, from Eqs. (16) and (17), we obtain a value for $G_{\varphi K \overline{K}}^2$ which leads to the φ width:

$$\Gamma(\varphi \rightarrow K^+ + K^-) = 2.0 \text{ MeV}, \qquad (18)$$

to be compared with the two experimental values 0.9 ± 0.2 and 1.6 ± 0.15 .²²

‡On leave of absence from The Tata Institute of Fun-

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[†]On leave of absence from The Centre of Advanced Studies in Theoretical Physics and Astrophysics, University of Delhi, Delhi, India.

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NEAR FORWARD PEAKS IN THE K^-p AND π^-p CHARGE-EXCHANGE SCATTERING*

A. O. Barut and H. Kleinert

Department of Physics, University of Colorado, Boulder, Colorado (Received 8 April 1966)

The peak observed for very small t (square of the momentum transfer in the center-of-mass frame) in the $\pi^- p^1$ and $K^- p^2$ charge-exchange scatterings has been explained recently on the basis of the contribution of the spin-flip amplitude.²⁻⁴ In the first analysis,³ for K^-p , the ratio of the spin-flip to spin-nonflip contribution necessary to explain the data comes out to be large (about 1.8 at t = 0.4) and cannot be accounted for by the mechanism of ρ exchange, not even with absorptive corrections. In the second analysis,⁴ for $\pi^- p$, the spin-flip amplitude is again assumed to be larger than the spin-nonflip amplitude and to have essentially a diffraction-type behavior with a maximum at small t. Also in $\pi^- p$, a ρ -exchange model with absorptive corrections fails completely to explain the behavior of the charge-exchange scattering.⁵

Such a large spin-flip contribution at these very small *t* values and the shape of the spinflip amplitude would be quite startling. For one thing, the spin-flip amplitude is expected to become appreciable only around $\sin\theta \approx 1$ or $t = -2q^2$ (-9.85 at 10 BeV/c) and is likely to cause the secondary diffraction peak, also observed in charge-exchange scattering.⁶ Furthermore, the spin-flip amplitude *g* involves the difference of the partial-wave amplitudes $\alpha_{l,+}$ and $\alpha_{l,-}$, whereas the spin-nonflip amplitude *f* is the sum of these partial-wave amplitudes:

$$f = \sum_{l} \{ la_{l, +} + (l+1)a_{l, -} \} P_{l},$$

$$g = \sum_{l} \{ a_{l, -} - a_{l, +} \} P_{l}^{1}(z).$$

Therefore, we expect only a few low partial waves to contribute to the spin-flip amplitude, whereas all partial waves together make up the diffraction peak of the spin-nonflip part of the scattering. In fact, the elastic scattering on the whole range of momentum transfer and the secondary diffraction peak near $\sin\theta = 1$ can be well explained by a single, constant and small, *p*-wave spin-flip amplitude in the case of $\pi^- p$ scattering.⁷ Thus, the spin-flip amplitude alone is not expected to show a diffraction behavior.

We want to point out that the charge-exchange scattering, being the difference of two isospin amplitudes, is very sensitive to the changes in the relative phase, as a function of t, of the two isospin amplitudes and accounts in a simple way for the behavior of the charge-exchange scattering. The elastic scatterings are not sensitive to this relative phase. Thus, it would be important to measure the relative phase in the charge-exchange scattering. A small spinflip contribution may be introduced which accounts for the secondary peaks, as in the case of elastic scattering. Furthermore, the parameters of the charge-exchange scattering can be related to those of the elastic scattering.

<u> K^-p scattering</u>.-We assume that each of the I=0 and I=1 amplitudes has a diffraction-peak behavior, for small t, of the form

$$A^{0} = ae^{\frac{1}{2}\alpha t}, \quad A^{1} = \eta ae^{\frac{1}{2}\alpha t}, \quad (1)$$

where η contains a relative phase $\varphi(t)$ between the two amplitudes $[\eta = |\eta| e^{i\varphi(t)}]$. For simplicity we have assumed the same exponent α in