COMPARISON OF K^{\pm} P DISPERSION RELATIONS WITH EXPERIMENT

M. Lusignoli, M. Restignoli and G. Violini Istituto di Fisica dell'Università - Roma

and

G. A. Snow *),**)

CERN - Geneva

ABSTRACT

Once subtracted forward angle dispersion relations are used to predict the real parts of the $K^{\pm}P$ scattering amplitudes as a function of energy, using previously derived values of the NYK coupling constants. These predictions are compared with experiment.

^{*)} John S. Guggenheim Fellow on sabbatical leave from the University of Maryland, College Park, Md., U.S.A.

^{**)} Supported in part by the U.S. Atomic Energy Commission.

Recently we have evaluated $^{1)}$ the NYK coupling constants using unsubtracted forward dispersion relations for the K-N system. As described in I, the essential data that was used in this calculation was the K $^{\pm}$ N total cross-sections at all available energies and the real part of the K $^{\pm}$ N forward scattering amplitudes at very low energies $^{2)}$. In this note we want to present the predictions of the oncesubtracted dispersion relations for the real parts of the K $^{\pm}$ p forward scattering amplitudes at higher energy and compare them with experiment. A recent work by N. Zovko $^{3)}$ has questioned the consistency of the low energy K $^{\pm}$ N scattering data with the dispersion relations. We find that such a discrepancy does not exist if one includes, in a consistent manner, (K $^{-}$ p, $^{-}$ R 0 n) mass difference effects in the Dalitz-Tuan formalism 4 for the low energy K $^{-}$ p scattering.

Since the publication of I, more precise K^{\pm} N total cross-sections between 1.0 and 2.45 GeV/c have become available 5). We have redone the calculations of I including these data and find the same results for the NYK coupling constants namely $g_{p \wedge K}^2 = 4.8 \pm 1.0$, $g_{p \leq 0K}^2 \lesssim 3.2$.

The once subtracted dispersion relations can be written in the form 6)

$$D_{\pm}(\omega) = \frac{1}{2} \left[D(m_{K}) + D_{+}(m_{K}) \right] + \frac{\omega}{2m_{K}} \left[D(m_{K}) - D_{+}(m_{K}) \right] + \frac{k^{2}}{4\pi^{2}} \int_{m_{K}}^{\infty} \frac{\omega'(\sigma_{-} + \sigma_{+}) + \omega(\sigma_{-} - \sigma_{+})}{k'(\omega'^{2} - \omega'^{2})} d\omega' + \frac{k^{2}}{4\pi^{2}} \int_{m_{K}}^{m_{K}} \frac{(\omega'_{+} \omega) A_{-}}{k'^{2}(\omega'^{2} - \omega^{2})} d\omega' + \frac{k^{2}}{y(\omega_{y} \pm \omega)} \frac{\chi(y)}{\omega_{y}^{2} - m_{K}^{2}}$$

$$(1)$$

where D and A are, respectively, the real and imaginary parts of the (K^{\pm}, p) scattering amplitude in the laboratory system,

$$w_y = \frac{m_y^2 - m_p^2 - m_k^2}{2m_p}$$
 and $X(Y) = g_{pyk}^2 \frac{(m_y - m_p)^2 - m_k^2}{4m_p m_y}$

 $Y=\Lambda$, and \sum_{pYK}^{0} is the renormalized, rationalized pseudoscalar coupling constant, ω and k are the total energy and the momentum of the K meson in the laboratory system respectively. As in I, A_ in the unphysical region is assumed to have its S wave part given by the Dalitz-Tuan formalism and a p wave part only in I=1 coming from a Y_1^* (1385) pole term, which can be written in the form $k^2/(\omega_Y \pm \omega) \sum_{i=1}^{\infty} Z_p(Y_1^*)$.

The numerical evaluation of $D_{\pm}(\omega)$ is insensitive to exactly how the I=1 pole term is divided between the ξ^0 pole and the Y_1^* pole as long as Eq. (3') of I is used. For definiteness we have assumed that $g_{p, \xi^0 K}^2 = \frac{1}{27} g_{p, k}^2 K$ corresponding to an f/d ratio of $\frac{2}{3}$.

Above 20 GeV/c, we used a Regge pole model extrapolation for the total cross-sections. We used the parameters kindly provided by C. Foggart and R.J.N. Phillips. The results for $D_{\pm}(\omega)$ for $\lesssim 3.5$ GeV/c are not sensitive to the exact form of the particular Regge fit used in the asymptotic region, since the total asymptotic contribution is less than the errors. We have not explored the sensitivity of $D_{\pm}(\omega)$ for higher momenta, to the fit used in the asymptotic region, so that the errors quoted for the contributions from the asymptotic region are only the statistical ones associated with this particular model.

The evaluation of the principal value integrals was carried out using a straight line interpolation between the most accurate experimental points.

<u>K</u>+p :

The calculated values of $\propto (\omega) = D_{+}(\omega)/A_{+}(\omega)$ for $k \gtrsim 0.7$ GeV/c, are presented in Fig. 1, together with all known experimental values $^{8)}$ of \propto $_{+}$. Above 1.2 GeV/c only the magnitudes of $lpha_{\!\scriptscriptstyle\perp}$ have been measured, but they have been plotted as negative quantities in Fig. 1. The values of $D_{+}(\omega)$ have been calculated using both the low energy $K^{-}p$ parameters of solution I of Kim 9) and of solution II of Sakitt et al. 10) (the K p parameters of Kittel et al. 11) are essentially the same as those of Sakitt et al.). In fact we find that these different solutions yield values for $\mathrm{D}_{+}(\omega$) that differ by a small fraction of the calculated error. Therefore only the points of Kim are plotted. This insensitivity of $\mathrm{D_+}(\omega)$ to the low energy K p solutions is a consequence of our use of the unsubtracted dispersion relation in I to derive the Λ and $(^{\circ} + Y_1^*)$ pole term contribution : that is we use different pole term contributions for Kim and for Sakitt parameters as given in Eq. (3!) of I. The agreement between dispersion theory predictions and experiment for $\mathbb{D}_+(\omega)$ and $\propto_+(\omega)$ is reasonably good except in the highest momentum region (k \gtrsim 10 GeV/c). Similar but not identical calculations for $D_+(\omega)$ have been made recently by N.M. Queen above 1 GeV/c. The general trend of his results for $\mathrm{D_+}(\omega)$ is the same as ours except that when he uses our value of the NYK coupling constants obtained from I his magnitude of $\mathrm{D_{+}(\,\omega\,)}$ is substantially larger than ours, and hence in less good agreement with the experimental data.

To illustrate the contributions to $D_+(\omega)$ and to $\Delta D_+(\omega)$ from different parts of Eq. 1, we divide the right-hand side of Eq. 1 into six different pieces c_i (1 = 1,...,6), defined as follows:

 $c_1 = // + \sum_{i=1}^{O} + Y_1^*$ pole term contributions

c₂ = the low energy K p contributions from $\omega_{\Lambda\pi}$ = 240 MeV to $\widetilde{\omega}_{-}$ = 574.2 MeV including the D_(m) term. (The uncertainty in c₂ includes all the errors from the low energy K p solutions.)

 $c_3 = \text{the } K^{-}p$ contributions from ω_{-} to 20 GeV

 c_4 = the low energy K^+p contributions from m_K to $\widetilde{\omega}_+$ = 914 MeV including the $D_+(m_K)$ term

 $c_5 = the K^+p$ contributions from $(2)_+$ to 20 GeV

 c_6 = the K⁺ and K⁻ cross-section integrals in the asymptotic region, $\omega > 20$ GeV.

The values of c_1 and Δc_1 at several representative momenta are listed in Table 1. We assume that the final error for $D_+(\omega)$ is given by a statistical combination of the errors of the six individual parts. This assumption is not rigorously correct since the errors in the hyperon pole terms are complicated functions of the experimental cross-sections used in I and again in Eq. (1). However this method of calculating the errors should give a reasonable order of magnitude estimate of the uncertainties.

к_р:

The calculated values of $\propto (\omega) = D(\omega)/A(\omega)$ are displayed in Fig. 2, along with the available experimental values 13) D (ω) is more sensitive to the S wave K p parameters so that there is a non-negligible difference between the results obtained using the Kim 9) and Sakitt 10) solutions, especially at low momenta. Again only the Kim solution is plotted. The Sakitt solution II prediction is rather close to the lower limit error curve of the Kim solution. The errors in the Sakitt case are about 30% larger than for Kim. The structure in D_(ω) versus ω reflects the resonant structure in the K p total crosssections at Y_0^* (1520), Y_0^* (1815), and Y_0^* (2100) (5). The agreement between the dispersion theory prediction and experiment is not very good at k = 1.97 GeV/c, Cook et al. 13), which is also the experimental point with the smallest apparent error. Confirmation of such a discrepancy would certainly require some substantial modifications of our analysis. The small values for \propto _ predicted in the vicinity of Y_0^* (1520 MeV), $k_z = 400$ MeV/c, are consistent with solution III of Watson et al. 13) but not with their solution I, which was their favoured one. This latter predicts values for α ~ +0.4 in the vicinity

of the resonance, which is not compatible with the dispersion relation predictions. Recently Kittel and Otter 14) have given arguments based on an effective range treatment of I=0, $\overline{K}n$ interactions that also favour Watson et al.'s solution III over solution I. The essential point is that the S wave parts of solution III of Watson et al. are much closer to the low energy solutions of Kim and Sakitt et al. than are the S wave parts of solution I.

Again, to illustrate the contributions to $D_{-}(\omega)$ and to $\Delta D_{-}(\omega)$ from different parts of Eq. (1), we divide the r.h.s. of Eq. (1) into the same six pieces c_1, \dots, c_6 as defined previously for K^+ with the necessary sign changes incorporated in the definition. The values of c_1 and Δc_1 at several representation momenta are listed in Table II.

In conclusion, we find that the dispersion relation predictions, for (), using the results of I, are in reasonable agreement with experiment. Less data for $\propto (\omega)$ is available, but there may be a discrepancy at k = 1.97 GeV/c. Clearly much more data on $D_+(\omega)$ would be of value for such comparisons, which can then also be used to make a more accurate determination of the KYN coupling constants. In the absence of such data, we think that the values of (ω) given in Figures 1 and 2 can be used to help decide between alternative (K^{\pm}, p) phase shift solutions in the momentum region below ~ 2 GeV/c, as illustrated by our discussion of the 400 MeV/c K, p analysis.

One of us, G.A.S., would like to thank the members of the Physics Department of the University of Rome for their kind hospitality during the academic year 1965-66, while he was the recipient of a Fulbright research scholarship. We are indebted to the members of the Theoretical Group at Rome for many helpful discussions and to Drs. Foggart and Phillips for providing us with a recent Regge Pole fit to the cross-section data before publication. Finally we would like to thank Professor J. Prentki and the Theoretical Study Division at CERN for hospitality while this paper was written.

Table I : Contributions to $D_{+}(\omega)$ from different parts of the dispersion relation given by Eq. (1) for $K^{+}p$.

[See text for detailed explanation of c_{1} (i = 1,...,6)]

	•		:				
k	^c 1	c ₂	° 3	. ^c 4	°5	c ₆ *)	D+-1)(GeV)
(GeV/c)	(Pole terms)	(K low energy)	(K ⁻ higher energy)	(K [†] low . energy)	(K ⁺ higher energy)	(Regge contrib)	(Total)
	0.457	-0.802	+0.905	-3•493	0.665	0.032	-2.24
0,520	<u>+</u> 0.087	<u>+</u> 0.027	±0.010	<u>+</u> 0.252	<u>+</u> 0.005	<u>+</u> 0 •001	<u>+</u> 0.27
	1.100	-2.067	2.620	-5.707	1.957	0.112	-1.99
0.970	<u>+</u> 0.21	<u>+</u> 0.074	<u>+</u> 0.03	<u>+</u> 0.600	<u>+</u> 0.043	<u>+</u> 0.003	+0.64
	2.56	- 5.13	7.71	-8.30	-1.00	0.46	-3.71
1.970	<u>+</u> 0 • 50	<u>+</u> 0.19	<u>+</u> 0.07	<u>+</u> 0.76	<u>+</u> 0.06	<u>+</u> 0.02	<u>+</u> 0.92
	4.78	-9.89	17.01	-13.03	-5.21	1.46	-4.88
3.500	<u>+</u> 0.93	±0.35	±0.15	±1.16	<u>+</u> 0.22	<u>+</u> 0.08	<u>+</u> 1.56
	47 07	-29.51	61.64	-33.00	-34.19	12.25	-8.92
9.800	13.87 +2.80	±1.10	±0.48	+2.88	±0.65	<u>+</u> 0.68	<u>+</u> 4.20

^{*)} The errors in ${\bf c}_6$ reflect only the errors in the Regge pole model of Foggart and Phillips and do not include uncertainties related to choosing different asymptotic models.

Table II : Contributions to D_(ω) from different parts of the dispersion relation given by Eq. (1) for K p. [See text for detailed explanation of c_i (i = 1,...,6)]

k	° 1	°2	° 3	c ₄	c ₅	e ₆ *)	D_ 1
(GeV/c)	(Pole terms)	(K low energy)	(K ⁻ higher energy)	(K ⁺ low energy)	(K ⁺ higher energy)	(Regge contrib)	(GeV ⁻¹) (Total)
0.415	- 0.61	-0.30	- 0.45	0.468	0.157	0.020	-0.71
	+ 0.13	<u>+</u> 0.14	+ 0.28	<u>+</u> 0.039	<u>+</u> 0.001	<u>+</u> 0.0004	<u>+</u> 0.34
0.950	- 1.48	3•41	- 2.25	1.80	0.711	0•107	-2.30
	+ 0.30	±0•17	± 0.23	<u>+</u> 0.15	<u>+</u> 0.003	<u>+</u> 0•003	+0.42
1.975	- 3.02	6.97	-11.96	4•84	2.401	0•466	-0.30
	<u>+</u> 0.60	<u>+</u> 0.26	± 0.16	<u>+</u> 0•41	±0.009	<u>+</u> 0•020	<u>+</u> 0.78
3•460	- 5.18	11.67	-22.58	9•48	5.64	1.44	+0.27
	± 0.73	±0.40	+ 0.37	<u>+</u> 0•81	<u>+</u> 0.02	<u>+</u> 0.06	+1.22
5.000	- 7.43	16.49	-36.45	14.34	9•55	3•06	-0.44
	+ 0.15	±0.55	+ 0.53	+1.23	<u>+</u> 0•04	<u>+</u> 0•15	+2.10
9.000	-13.18	28 .96	-76.99	27.05	21.06	10.56	-2.55
	± 2.65	<u>+</u> 0.93	<u>+</u> 0.69	<u>+</u> 2.33	<u>+</u> 0.08	<u>+</u> 0.49	<u>+</u> 3.74

^{*)} The errors in c₆ reflect only the errors in the Regge pole model of Foggart and Phillips and do not include uncertainties related to choosing different asymptotic models.

R E F E R E N C E S

- 1) M. Lusignoli, M. Restignoli, G.A. Snow and G. Violini Phys.Letters 21, 229 (1966), hereafter called I.
- 2) See I for the references to the experimental data used, besides those listed below.
- 3) N. Zovko Z.für Physik 192, 346 (1966).
- 4) R.H. Dalitz and S.F. Tuan Ann. Phys. N.Y., 10, 307 (1960).
- 5) R.L. Cool, G. Giacomelli, T.F. Kycia, B.A. Leontic, K.K. Li, A. Lundby and J. Teiger Phys.Rev.Letters 16, 1228 (1966) (K); and Phys.Rev.Letters 17, 102 (1966) (K⁺).
- 6) See Refs. 2), 3) of I.
- 7) F. Buccella, M. Lusignoli and G. Violini Phys. Letters 21, 572 (1966).
- 8) a) T.F. Stubbs, H. Bradner, W. Chinowsky, G. Goldhaber, S. Goldhaber, W. Glater, D.M. Stork and H.K. Ticho Phys.Rev.Letters
 7, 188 (1966).
 - b) V. Cook, D. Keefe, L.T. Kerth, P.G. Murphy, W.A. Wenzel and T.F. Zipf Phys.Rev. 129, 2743 (1963).
 - c) W. Chinowsky, G. Goldhaber, S. Goldhaber, T. O'Halloran and B. Schwarzschild Phys. Rev. 139, B1411 (1965).
 - d) J. De Baisieux, F. Grard, J. Heughebaert, L. Pape, R. Windmolders, R. George, Y. Goldschmidt-Clermont, V.P. Henri, D.W.G. Leith, G.R. Lynch, F. Muller, J.M. Perreau, G. Otter and P. Sällstrom Nuovo Cimento 43, A142 (1966).

- e) W. De Baere, J. De Baisieux, P. Dufour, F. Grard, J. Heughebaert,
 L. Pape, P. Peeters, F. Verbeure, R. Windmolders, R. George,
 Y. Goldschmidt-Clermont, V.P. Henri, B. Jongejans, D.W.G. Leith,
 A. Moisseev, F. Muller, J.M. Perreau and V. Yarba Nuovo
 Cimento (to be published).
- f) K.J. Foley, S.J. Lindenbaum, W.A. Love, S. Ozaki, J.J. Russell and L.C.L. Yuan Phys.Rev.Letters 11, 503 (1963).
- 9) J.K. Kim Phys. Rev. Letters 14, 29 (1965).
- 10) M. Sakitt, T.B. Day, R.G. Glasser, N. Seeman, J. Friedman, W.E. Humphrey and R.R. Roos Phys.Rev. 139, 572 (1965).
- 11) W. Kittel, G. Otter and I. Wacek Phys. Letters 21, 349 (1966).
- 12) N.M. Queen University of Birmingham Preprint (1966).
- 13) a) M.B. Watson, M. Ferro-Luzzi and R.D. Tripp Phys.Rev. <u>131</u>, 2248 (1963).
 - b) A. Fridman, O. Benary, A. Michalon, B. Schiby, R. Strub and G. Zech Phys.Rev. 145, 1136 (1966).
 - c) V. Cook, B. Cork, T.F. Hoang, D. Keefe, L.T. Kerth, W.A. Wenzel and T. Zipf Phys.Rev. 123, 320 (1961).
 - d) M.N. Focacci, S. Focardi, G. Giacomelli, F. Serra, M.P. Zerbetto and L. Monari Phys. Letters 19, 441 (1965).
 - e) J. Gordon Phys. Letters 21, 117 (1966).
- 14) W. Kittel and G. Otter Phys. Letters 22, 115 (1966).

FIGURE CAPTIONS

- Figure 1 K^+P dispersion relation predictions of $\alpha_+ = D_+/A_+$ as a function of the K^+ laboratory momentum. The references for the experimental points are given in Ref. 8).
- Figure 2 K-P dispersion relation predictions of \propto _ = D/A as a function of the K laboratory momentum. The references for the experimental points are given in Ref. 13).



