Low Energy Constants from $K_{\ell 4}$ Form-Factors

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Abstract

We have calculated the form-factors $F$ and $G$ in $K \to \pi\pi\ell\nu$ decays ($K_{\ell 4}$) to two-loop order in Chiral Perturbation Theory (ChPT). Combining this together with earlier two-loop calculations an updated set of values for the $L_i^r$, the ChPT constants at $O(p^4)$, is obtained. We discuss the uncertainties in the determination and the changes compared to previous estimates.

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1. The theoretical study of $K \to \pi\pi\ell\nu$ decays, $K_{\ell 4}$ decays, provides an interesting possibility to test our understanding of the long-distance dynamics of the strong sector in the Standard Model. Perturbative QCD can not be applied at energies lower than the spontaneous symmetry scale, $\Lambda_\chi \sim m_\rho$. One has to resort to an effective approach to the full theory, Chiral Perturbation Theory (ChPT) [1], to obtain a reliable model independent description of physical processes. The main purpose in the study of $K_{\ell 4}$ decays is twofold: $i$) This decay is one of the cleanest sources of $\pi\pi$ pairs at low-energies and thus provides us with the possibility to check $\pi\pi$-scattering near threshold. $ii$) The form-factors themselves are also directly of interest. They provide a direct test of our understanding of the three-flavour sector and are in addition one of the major inputs to determine the needed constants to predict other quantities. The early history can be found in the review [2]. $K_{\ell 4}$ have been treated in the context of ChPT at one-loop [3] and in a dispersively improved one-loop representation [4]. These studies delivered the standard values of the low-energy constants $L_1^r, L_2^r$ and $L_3^r$.

On the other hand, $\pi\pi$-scattering has been studied at two-loop order in two-flavour ChPT, direct [5] and via dispersive methods [6]. The main remaining uncertainty on the prediction of $\pi\pi$-scattering is the value of the low-energy constants. In [5] the standard values from [4] were used. A Roy equation analysis using high energy $\pi\pi$-data, gave significantly different values for these constants [7]. There are several possible sources for this discrepancy: the Omnès representation of [4] is not sufficient for $K_{\ell 4}$; the presence of large systematic effects in the $\pi\pi$-scattering data base and Roy analysis; or large corrections to the one-loop relation between two- and three-flavour ChPT constants. Therefore, and since new experiments are underway, a full two-loop calculation of $K_{\ell 4}$ is necessary. A first step was the estimate of the double logarithm corrections of $O(p^6)$ to $K_{\ell 3}$ and $K_{\ell 4}$ [8]. It was shown there they could be large. In addition pions in the $I=0, S$-wave have strong final state interactions. We describe here first results of the full $K_{\ell 4} O(p^6)$ calculation.

2. $K_{\ell 4}$ decays are described by the matrix element

$$M^{ij} = \frac{G_F}{\sqrt{2}} V_{us}^* \bar{u}(p_\nu) \gamma_\mu (1 - \gamma_5) v(p_\ell) \langle \pi^i(p_i)\pi^j(p_j)|V^\mu - A^\mu|K(k) \rangle. \quad (1)$$

Lorentz invariance allows to parametrize the hadronic part by four form-factors

$$\langle \pi^i\pi^j|A^\mu|K \rangle = -\frac{i}{m_K} \left[ P^\mu F^{ij} + Q^\mu G^{ij} + L^\mu R^{ij} \right],$$

$$\langle \pi^i\pi^j|V^\mu|K \rangle = -\frac{1}{m_K^2} \epsilon^{\mu\nu\alpha\beta} L_\nu P_\alpha Q_\beta H^{ij}, \quad (2)$$


with $P = p_i + p_j$, $Q = p_i - p_j$ and $L = p_f + p_\nu$. $p_i$ is the momentum of $\pi_i$. All four form-factors are dimensionless functions of three variables, $s_\pi = P \cdot P, u_\pi = (k - p_i)^2$ and $t_\pi = (k - p_j)^2$. They can be decomposed, into a part symmetric or antisymmetric under $t_\pi \leftrightarrow u_\pi$ interchange, corresponding respectively to the isospin 0 or 1 part for $F, R$ and isospin 1 or 0 for $G, H$. These processes are dominated by $F_s$ and $R_s$ and are real functions with argument $F, R$ corresponding respectively to the isospin 0 or 1 part for $F, R$. The term $R_s$ only contributes proportional to $m_\pi^2$ and can thus be neglected in the decays with an electron.

The remainder we concentrate on the $ij = ++$ channel. The others can be derived from it using isospin relations. The older available experiments are compatible with the $K^+ \rightarrow \pi^+\pi^-e^+\nu_e$ experiment [10]. The most recent other measurement is $K_L^0 \rightarrow \pi^+\pi^-e^+\nu_e$ [11].

The experiment [10] relied on a partial wave analysis of the form-factors. The analysis only kept $s$ and $p$ waves, the effect of $d$-waves as well as of $s_t$ was within the measurement errors and neglected in the further analysis. They used the parametrization

$$F = f_s e^{i\delta_s} + f_p e^{i\delta_p} \cos \theta_\pi + D\text{-wave} + \ldots, \quad G = g e^{i\delta_p} + D\text{-wave} + \ldots, \quad (3)$$

where $\delta_i$ are the $i$-wave strong two-pion final state phase shifts, and $f_s, f_p$ and $g$ are defined to be real functions with argument $s_\pi$. They then divided the data in 5 energy bins, observed an $s_\pi$ dependence in $f_s$ and one compatible with it for $g$. The final fit was performed using

$$f_s(q^2) = f_s(0) \left(1 + \lambda_f q^2\right), \quad g(q^2) = g(0) \left(1 + \lambda_g q^2\right), \quad q^2 = (s_\pi - 4m_\pi^2)/(4m_\pi^2). \quad (4)$$

with $\lambda_f = \lambda_g$. Furthermore, the $f_p$ value was compatible with zero. The results obtained, with $\sin \theta_C = 0.22$, are

$$f_s(0) = 5.59 \pm 0.14, \quad g(0) = 4.77 \pm 0.27, \quad \lambda_f = 0.08 \pm 0.02. \quad (5)$$

Using isospin the result of [11] is $g(0) = 5.50 \pm 0.50$ which using PDG procedures [12] leads to a combined

$$g(0) = 4.93 \pm 0.31. \quad (6)$$

Note that [11] neglected $f_p$ which can account for up to 5% of the value of $g(0)$.

3. As mentioned above, earlier calculations gave an indication of potentially sizeable corrections at $O(p^2)$. In order to obtain the full correction we have evaluated both form factors, $F$ and $G$, at next-to-next-to-leading order (NNLO) in ChPT following a diagrammatic approach. One faces the evaluation of a large number of diagrams but only one new topology, shown in Fig. 1, appears w.r.t. the vector-vector and axial-vector–axial-vector two point-functions [13]. It involves a new set of integrals, the vertex-integrals, that can be obtained in terms of a two-parameter integral representation [14].

The full evaluation leads to rather long expressions. More than half of the complexity stems from the topology of Fig. 1 due to the large number of possible mass and momentum combinations. We have performed several checks on our full expressions: i) As a basic test of the algebraic programs we recover the one-loop expressions [3, 4]. ii) All non-local divergences cancel when adding the full set of diagrams together with wave function renormalization. iii) The polynomial divergences also cancel against the counter-terms determined in general for the even-intrinsic parity representation [14].

Figure 1: Vertex topology. Dots refer to strong vertices or current insertions.
Here we follow the philosophy of most of the other renormalized physical ones and add the part coming from the matrix-elements. We fixed the parameters as much as possible from experiment or the comparison with diagrams involving three-point one-loop integrals can also be obtained via the renormalization of the one-loop graphs, and hence the three-point integrals should cancel in the final result. vi) The sunset-type integrals encountered in are in agreement with those already calculated in using different methods. vii) The two-particle discontinuities of the vertex-type integrals have been checked using the Cutkosky rules below the three-particle thresholds. viii) Our results satisfy the isospin relation $M^{+-} = M^{-0}/\sqrt{2} + M^{00}$. In view of these checks, we trust our calculations of the matrix-elements.

To obtain the final answer, we shift the bare quantities –masses and decay constants– to the renormalized physical ones and add the part coming from the $\mathcal{O}(p^6)$ Lagrangian as determined in [10].

4. One of the main problems is how to deal with the coefficients of the $\mathcal{O}(p^6)$ Lagrangian. Here we follow the philosophy of most of the other $\mathcal{O}(p^6)$ calculations and estimate them using resonance exchange. This worked well at $\mathcal{O}(p^4)$ [17]. In the second paper of [18] this was found to agree reasonably well at $\mathcal{O}(p^6)$ for the form-factors considered there. We use the notation as defined in [13] but include also terms which did not contribute there. Specifically we use for the vector nonet matrix $V_{\mu}$

$$\mathcal{L}_V = -\frac{1}{4} (V_{\mu \nu} V_{\rho \sigma}) + \frac{1}{2} m_V^2 \langle V_{\mu \nu} V_{\rho \sigma} \rangle - \frac{f_V}{2\sqrt{2}} (V_{\mu \nu} f_{\mu \nu}^V) - \frac{i g_V}{2\sqrt{2}} \langle V_{\mu \nu} | u^\nu, u^\nu \rangle + f_\chi \langle V_{\mu \nu} | u^\nu, \chi_- \rangle$$

and for the axial-vector nonet $A_{\mu}$

$$\mathcal{L}_A = -\frac{1}{4} (A_{\mu \nu} A_{\rho \sigma}) + \frac{1}{2} m_A^2 (A_{\mu \nu} A_{\rho \sigma}) - \frac{f_A}{2\sqrt{2}} (A_{\mu \nu} f_{\mu \nu}^A) + \gamma_A^{(1)} \langle A_{\mu \nu} u^\nu u^\nu \rangle + \gamma_A^{(2)} \langle A_{\mu \nu} u^\nu, u^\nu \rangle.$$ (8)

We have chosen a representation where the vectors and the axial-vectors only start contributing to the mesonic Lagrangian at $\mathcal{O}(p^6)$. For the scalar nonet we take

$$\mathcal{L}_S = \frac{1}{2} \langle \nabla_{\mu} S \nabla^{\mu} S \rangle - \frac{1}{2} m_S^2 (S^2) + c_d \langle S w^\mu u^\mu \rangle + c_m \langle S \chi_+ \rangle .$$

We fixed the parameters as much as possible from experiment or the comparison with $\mathcal{O}(p^4)$ [13, 17]. For the remainder we use the values predicted by the Nambu-Jona-Lasinio model [13]. The specific inputs used are $m_V = 0.77$ GeV, $m_A = 1.23$ GeV, $m_S = 0.98$ GeV, $f_V = 0.20$, $f_\chi = -0.025$, $g_V = 0.09$, $\alpha_V = -0.014$, $f_A = 0.1$, $\gamma_A^{(1)} = 0.006$, $\gamma_A^{(2)} = -0.01$, $c_m = 42$ MeV and $c_d = 32$ MeV. These are used in all the fits of Table 1 except fit 6 where we kept only $f_V$, $f_\chi$ and $g_V$ as nonzero couplings.

5. ChPT in the meson sector has as parameters at $\mathcal{O}(p^2)$ $F_0$, $B_0$ and the quark masses $\hat{m} = m_u = m_d$ and $m_s$. The latter only appear multiplied with $B_0$ for a total of 3 parameters. At $\mathcal{O}(p^4)$, we have $L_1', \ldots, L_9'$ [13] and at $\mathcal{O}(p^6)$ there are 90 additional parameters [16]. In the two-flavour sector the equivalent of $L_6$ and $L_9$ has been determined to NNLO [18]. $L_1'$ has no influence on the other quantities considered here and the influence of $L_9'$ is small. We will therefore use the standard values of [13, 17]. For the $\mathcal{O}(p^6)$ we use the resonance saturation approximation described above.

As input we use $F_\pi = 92.4$ MeV, $F_K/F_\pi = 1.22 \pm 0.01$ and the physical masses of pion, kaon and eta [13] in addition to $f_\pi(0)$, $g(0)$ and $\lambda_f$. The final input is the ratio $m_s/\hat{m}$. We further assume that the parameters $L_{1}'$ and $L_9'$ are small because they vanish in the large $N_c$-limit. When calculating the form-factors for $K^+ \to \pi^+ \pi^- e^+ \nu$, $F_\pi$ and $F_K$ we use the physical charged masses $m_{\pi^+} = 139.56995$ MeV; $m_{K^+} = 493.677$ MeV and $m_{\eta} = 547.30$ MeV. When comparing with the quark-mass ratios we use the $\pi^0$-mass, $m_{\pi^0} = 134.9764$ MeV, the physical $\eta$-mass and $m_K^2 = 1/2(m_K^2 + m_K^2 - 1.8(m_{\pi^+}^2 - m_{\pi^0}^2)) = (494.53$ MeV)$^2$. This removes the electromagnetic mass corrections including estimated corrections to Dashen’s theorem [19].
The axial-vectors contribute only for the vector-exchange with the experimentally determined quantities. The rest is from the scalar vector meson part in the resonance estimate. As can be seen the main contribution comes from the renormalization scale, $\mu$ encoded in the matrix elements. We set $m^*= (336 \text{ MeV})$, which is the same as in [4], since we use the same data the errors have not changed.

In conclusion, we have performed a NNLO calculation of the $F$ and $G$ form-factors in $K_{\ell4}$ decays. We then used this calculation together with earlier calculations of masses and decay constants to update the ChPT parameters at $O(p^6)$ with a certain range of choices for the fitting of the available $K_{\ell4}$ data and the estimates of the $O(p^6)$ parameters. None of the variations of input change the parameters outside the experimentally determined error except the rather extreme case of $\mu = 0.5$ GeV.

When looking at the $K_{\ell4}$ calculation it can be seen that for the form-factor $F$ the estimate of higher orders of [4] was in the right direction. The changes are somewhat larger than naively expected.
The errors should be taken with caution, \( L_1^r \) and \( L_2^r \) are strongly anti-correlated and \( L_3^r \) is strongly anti-correlated with \( L_1^r \) and \( L_2^r \). As an example of this we have shown in Fig. 2 a distribution of sets of \( L_i^r \) that fall within the 68\% confidence level limit. If we choose as fit variables \( X_1 = L_2^r - 2L_1^r - L_3^r \), \( X_2 = L_2^r \) and \( X_3 = (L_2^r - 2L_1^r)/L_3^r \), \( X_3 \) is little correlated with the others. We obtain

\[
X_3 = 0.12^{+0.08}_{-0.11}. \tag{10}
\]

The large \( N_c \)-prediction \(|X_3| \ll 1\) is obviously well satisfied. These questions will be discussed in more detail in a future publication.

With these values for the low-energy constants we can see how the various quantities behave. The numbers correspond to \( \mathcal{O}(p^2) \), \( \mathcal{O}(p^4) \) and \( \mathcal{O}(p^6) \) contributions.

\[
\begin{align*}
m_{\pi^0}/m_{\pi^0}\exp &= 0.740 + 0.007 + 0.253, \\
m_K^2/m_K^2\exp &= 0.689 + 0.024 + 0.287, \\
m_n^2/m_n^2\exp &= 0.735 - 0.066 + 0.331, \\
F_K/F_\pi &= 1.000 + 0.135 + 0.085, \\
F_\pi/F_0 &= 1.000 + 0.136 - 0.076, \\
f_\pi(0) &= 3.78 + 1.15 + 0.66, \\
g(0) &= 3.78 + 0.82 + 0.17, \\
\lambda_f &= 0.000 + 0.127 - 0.047. \tag{11}
\end{align*}
\]

These numbers were calculated before rounding the \( L_i^r \) to the number of significant digits given in Table 1. Contrary to what was observed in \([8, 21]\) the slope is now mainly from \( \mathcal{O}(p^4) \)-effects.

Making use of the central values of the main fit in Table 1 together with the \( \mathcal{O}(p^4) \) relations between the two- and three-flavour low-energy constants we estimate \( \bar{t}_1 \) and \( \bar{t}_2 \)

\[
\bar{t}_1 = 0.4(-0.2), \quad \bar{t}_2 = 4.9(5.2). \tag{12}
\]

The values in brackets are those using the values of fit 9. We remind the reader that the relation used has possibly large \( \mathcal{O}(p^6) \) corrections. These should be compared respectively with \(-1.7 \) and \( 6.1 \) using the \( L_i^r \) from \([2] \) and with \( \bar{t}_2 = 4.2 \) \([12] \)\).

In conclusion, we have performed a full \( \mathcal{O}(p^6) \) calculation of \( K_{\ell 4} \) decays. We then performed a full refit of the \( \mathcal{O}(p^4) \) parameters of ChPT and discussed the changes and the validity of a large-\( N_c \) prediction.

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**Figure 2:** Correlation plot of \( L_1^r \) versus \( L_3^r \) and \( L_2^r \) versus \( L_5^r \). Shown are the points from a distribution of \( L_i^r \) with the \( \chi^2 \) calculated from the observables described in the text as for our main fit. Only points within a 68\% confidence level are included.
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