# **Critical look at rescattering effects on**  $\gamma$  from  $B^+ \rightarrow K \pi$

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Three ways of dealing with rescattering effects in  $B^{\pm} \to K^0 \pi^{\pm}$  are compared, in order to determine the weak phase  $\gamma$  from these processes and  $B^{\pm} \to K^{\pm} \pi^0$ . We find that neglecting these contributions altogether may involve sizeable errors in  $\gamma$ , depending on the rescattering amplitude and on the value of a certain measurable strong phase. We show that an attempt to eliminate these effects by using the charge-averaged rate of  $B^{\pm} \rightarrow K^{\pm} K^0$  suffers from a large theoretical error due to SU(3) breaking, which may be resolved when using also the processes  $B^{\pm} \rightarrow \pi^{\pm} \eta_8$ .

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## **I. INTRODUCTION**

The weak phase  $\gamma = \text{Arg}(V_{ub}^*)$  is presently the least well known quantity among the four parameters (three angles and a phase) of the Cabibbo-Kobayashi-Maskawa (CKM) matrix. Its determination, which is regarded to be more difficult than that of the other two angles of the CKM unitarity triangle  $[1]$ , can provide a crucial test of the CKM mechanism for *CP* violation in the standard model. Several methods have been proposed to determine  $\gamma$  from hadronic two-body *B* decays. The methods which seem to be experimentally most feasible in the near future are based on applications of  $SU(3)$ flavor symmetry in *B* decays into two light charmless pseudoscalars [2]. These methods involve certain theoretical uncertainties, which are expected to be reduced when more data become available and when better theoretical understanding of hadronic *B* decays is achieved.

In a first paper in a series, Gronau, London, and Rosner (GLR) [3] proposed to extract  $\gamma$  by combining decay rate measurements of  $B^+ \to K\pi$ ,  $B^+ \to \pi\pi$  with their chargeconjugates.  $SU(3)$  breaking, occurring in a relation between  $B \rightarrow \pi \pi I = 2$  and  $B \rightarrow K \pi I = 3/2$  amplitudes, was introduced through a factor  $f_K/f_\pi$  when assuming that these amplitudes factorize. In its original version, suggested before the observation of the heavy top quark, the method of Ref. [3] neglected electroweak penguin (EWP) contributions and certain rescattering effects. Subsequently, model-calculations showed that due to the heavy top quark the neglected EWP terms were significant  $[4]$ ; and recently these terms were related by SU(3) to the  $B \rightarrow K \pi I = 3/2$  current-current amplitudes  $[5,6]$ . This led to a modification  $[7]$  of the GLR method, to be referred to as the Gronau-London-Rosner-Neubert (GLRN) method, which in the limit of flavor  $SU(3)$ symmetry includes EWP effects in a model-independent way. Corrections from  $SU(3)$  breaking, affecting the relation between EWP terms and current-current terms, were argued to be small  $[5,8]$ .

Assuming that the above  $SU(3)$  breaking effects are indeed under control, there is still an uncertainty due to rescattering effects. To determine  $\gamma$  from the above rates, one takes the  $B^+ \rightarrow K^0 \pi^+$  amplitude to be pure penguin, involving no term with weak phase  $\gamma$ . This assumption, which neglects quark annihilation and rescattering contributions from charmless intermediate states, was challenged by a large number of authors [9]. Several authors proposed ways of controlling rescattering effects in  $B^{\pm} \rightarrow K^0 \pi^{\pm}$  by relating them through SU(3) to the much enhanced effects in  $B^{\pm}$  $\rightarrow K^{\pm}\bar{K}^{0}$  [10–12] (see also Refs. [13–15]). The chargeaveraged rate of the latter processes can be used to set an upper limit on the rescattering amplitude in  $B^{\pm} \rightarrow K^{0} \pi^{\pm}$ . While present limits are at the level of  $20-30\%$  of the dominant penguin amplitude  $[6,8]$  (depending somewhat on the value of  $\gamma$ ), they are expected to be improved in the future. The smaller the rescattering amplitude is, the more precisely can  $\gamma$  be determined from the GLRN method. A recent demonstration  $[8]$ , based on a few possible rate measurements, seems to show that if the rescattering amplitude is an order of magnitude smaller than the dominant penguin amplitude in  $B^+ \rightarrow K^0 \pi^+$ , the uncertainty in  $\gamma$  is only about 5°.

In the present Letter we reexamine in detail the uncertainty in  $\gamma$  due to rescattering effects. Using a geometrical interpretation for the extraction of  $\gamma$ , we perform in Sec. II numerical simulations which cover the entire parameter space of the two relevant strong phases, the rescattering phase  $\phi_A$  and the relative phase  $\phi$  between *I*=3/2 currentcurrent and penguin amplitudes. We find that, contrary to the demonstration made in Ref.  $\vert 8 \vert$ , a 10% rescattering amplitude leads to an uncertainty in  $\gamma$  as large as about 14 $\degree$  around  $\phi$  ~90°. For certain singular cases no solution can be found for  $\gamma$ . We show that  $\phi$  can be determined rather precisely from the  $B^{\pm} \rightarrow K\pi$  rate measurements [8], which could reduce substantially the error in  $\gamma$  if values far apart from  $\phi$  $=90^{\circ}$  were found.

It has been suggested  $[12]$  to go one step beyond setting limits on rescattering contributions in  $A(B^{\pm} \rightarrow K^0 \pi^{\pm})$  and to completely eliminate them by using the charge-averaged rate measurement of  $B^{\pm} \rightarrow K^{\pm} K^{0}$ . Applying our geometrical formulation, we will show in Sec. III that the resulting determination of  $\gamma$  is unstable under SU(3) breaking which can introduce very large uncertainties in  $\gamma$ .

Finally, in order to overcome these uncertainties, we have



FIG. 1. Relative orientation of the  $B^+\to K\pi$  amplitude triangles (4), normalized as described in the text. The triangle *OAA'* (*OBB'*) corresponds to the  $B^{+(-)} \rightarrow K\pi$  decays. The rescattering amplitude in  $B^{\pm} \rightarrow K^0 \pi^{\pm}$  is described by the line *OY*.

recently proposed to use in addition to  $B^{\pm} \rightarrow K^{\pm} \overline{K}^0$  also the processes  $B^{\pm} \rightarrow \pi^{\pm} \eta_8$  [16]. Although this may be considered an academic exercise, mainly due to complicating  $\eta$  $-\eta'$  mixing effects, we will examine in Sec. IV the precision of this method. We will show that when neglecting  $\eta$  $-\eta'$  mixing, the theoretical error in  $\gamma$  is reduced to a few degrees. We conclude in Sec. V. An algebraic condition, used in Sec. III to eliminate rescattering effects by  $B^{\pm}$  $\rightarrow$ *K*<sup> $\pm$ </sup>*K*<sup>0</sup> decays, is derived in an appendix.

## **II. RESCATTERING UNCERTAINTY IN**  $\gamma$  **FROM**  $B^{\pm} \rightarrow K \pi$

The amplitudes for charged *B* decays can be parametrized in terms of graphical contributions representing  $SU(3)$ amplitudes (we use the notations of Ref.  $[6]$ ):

$$
A(B^+\to K^0\pi^+) = |\lambda_u^{(s)}|e^{i\gamma}(A+P_{uc}) + \lambda_t^{(s)}(P_{ct}+P_3^{\text{EW}}),
$$
\n(1)

$$
\sqrt{2}A(B^+\to K^+\pi^0) = |\lambda_u^{(s)}|e^{i\gamma}(-T-C-A-P_{uc}) + \lambda_t^{(s)}(-P_{ct}+\sqrt{2}P_4^{EW}), \tag{2}
$$

$$
\sqrt{2}A(B^+\to \pi^+\pi^0) = |\lambda_u^{(s)}|e^{i\gamma}(-T-C),
$$
\n(3)

where  $\lambda_{q'}^{(q)}$  $(v_q^q) = V_{q'b}^* V_{q'q}$  are the corresponding CKM factors. These amplitudes satisfy a triangle relation  $[3,7]$ 

$$
\sqrt{2}A(B^+\to K^+\pi^0) + A(B^+\to K^0\pi^+)
$$
  
= 
$$
\sqrt{2}\tilde{r}_u|A(B^+\to \pi^+\pi^0)|e^{i(\gamma+\xi)}(1-\delta_{\text{EW}}e^{-i\gamma}).
$$
 (4)

Here we denote  $\tilde{r}_u = (f_K/f_\pi)\lambda/(1-\lambda^2/2) \approx 0.28$ ,  $\delta_{\text{EW}}$  $= - (3/2) |\lambda_t^{(s)} \rangle \lambda_u^{(s)} | \kappa \approx 0.66 \left[ \kappa \equiv (c_9 + c_{10})/(c_1 + c_2) \right] = -8.8$  $\times 10^{-3}$ ], while  $\xi$  is an unknown strong phase. The second term in the brackets represents the sum of EWP contributions to the amplitudes on the left-hand side  $[5,6]$ . The factor  $f_K/f_\pi$  accounts for factorizable SU(3) breaking effects.

The relation  $(4)$ , together with its charge-conjugate counterpart, written for  $\vec{A}(\vec{B} \rightarrow \vec{f}) \equiv e^{2i\gamma} A(\vec{B} \rightarrow \vec{f})$ , are represented graphically by the two triangles  $OAA'$  and  $OBB'$  in Fig. 1. Here all amplitudes are divided by a common factor  $A$  $\sqrt{2\tilde{r}}_{\mu} |A(B^+\to \pi^+\pi^0)|e^{i(\gamma+\xi)}$ , such that the horizontal line *OI* is of unit length and the radius of the circle is  $\delta_{\text{FW}}$ . Four of the sides of the two triangles are given by

$$
x_{0+} = \frac{1}{\sqrt{2}\tilde{r}_{u}} \frac{|A(B^{+} \to K^{0}\pi^{+})|}{|A(B^{+} \to \pi^{+}\pi^{0})|},
$$
  
\n
$$
x_{+0} = \frac{1}{\tilde{r}_{u}} \frac{|A(B^{+} \to K^{+}\pi^{0})|}{|A(B^{+} \to \pi^{+}\pi^{0})|},
$$
  
\n
$$
\tilde{x}_{0-} = \frac{1}{\sqrt{2}\tilde{r}_{u}} \frac{|A(\bar{B}^{-} \to \bar{K}^{0}\pi^{-})|}{|A(B^{+} \to \pi^{+}\pi^{0})|},
$$
  
\n
$$
\tilde{x}_{-0} = \frac{1}{\tilde{r}_{u}} \frac{|A(\bar{B}^{-} \to K^{-}\pi^{0})|}{|A(B^{+} \to \pi^{+}\pi^{0})|}.
$$
  
\n(5)

The relative orientation of the two triangles depends on  $\gamma$ and is not determined from measurements of the sides alone. Assuming that the rescattering amplitude with weak phase  $\gamma$ *in*  $B^+ \rightarrow K^0 \pi^+$  *can be neglected*, one takes the amplitude (1) to be given approximately by the second (penguin) term [3,7], which implies  $OB = e^{2i\gamma}OA$  in Fig. 1. In this approximation, the weak phase  $\gamma$  is determined by requiring that the angle  $(2\gamma)$  between *OA* and *OB* is equal to the angle  $(2\gamma)$ at the center of the circle  $[7]$ .

In order to study the precision of determining in this way the phase  $\gamma$  as function of the rescattering contribution which is being neglected, let us rewrite Eq.  $(1)$  in the form

$$
A(B^+\to K^0\pi^+) = -V_{cb}\left(1-\frac{\lambda^2}{2}\right)p(1+\epsilon_Ae^{i\phi_A}e^{i\gamma}),
$$
  

$$
p = P_{ct} + P_3^{\text{EW}},
$$
 (6)

where  $\epsilon_A$  measures the magnitude of rescattering effects. In Fig. 1 the magnitude of these effects has a simple geometrical interpretation in terms of the distance of the point *Y* from the origin *O*,  $\epsilon_A = |YO|/|YA|$ , where *YO* and *YA* are the two components in the  $B^+ \rightarrow K^0 \pi^+$  amplitude carrying weak phases  $\gamma$  and zero, respectively,

$$
YO = |\lambda_u^{(s)}| e^{i\gamma} [(A + P_{uc}) - p] / \mathcal{A}, \quad YA = V_{cb} \left( 1 - \frac{\lambda^2}{2} \right) p / \mathcal{A}.
$$

$$
(7)
$$

The rescattering phase  $\phi_A$  is given by  $\phi_A = \text{Arg}(YO/YZ)$ , where *Z* is any point on the line bisecting the angle *AYB*. A second strong phase which affects the determination of  $\gamma$  is  $\phi$ , the relative strong phase between the penguin amplitude  $p$ and the  $I=3/2$  current-current amplitude  $T+C$ . In Fig. 1 this phase is given by  $\phi = \text{Arg}(YZ/OI)$ .

Let us now investigate the dependence of the error in  $\gamma$ when neglecting rescattering on the relevant hadronic parameters. Our procedure will be as follows. First we generate a set of amplitudes based on the geometry of Fig. 1 and on given values of the parameters  $\gamma$ ,  $\epsilon$ ,  $\epsilon_A$ ,  $\phi_A$ , and  $\phi$ ; then we



solve the equation  $\cos 2\gamma = \cos(BOA)$  and compare the output value of  $\gamma$  with its input value. Here  $\epsilon$  is given in terms of the ratio of charge-averaged branching ratios  $[3,7]$ 

$$
\epsilon \equiv \frac{\lambda}{1 - \lambda^2/2} \frac{f_K}{f_\pi} \sqrt{\frac{2B(B^{\pm} \to \pi^{\pm} \pi^0)}{B(B^{\pm} \to K^0 \pi^{\pm})}},\tag{8}
$$

The geometrical construction in Fig. 1 is described by

$$
YA = \frac{e^{i(\phi - \gamma)}}{\epsilon \sqrt{1 + 2\epsilon_A \cos \phi_A \cos \gamma + \epsilon_A^2}} OI,
$$
  

$$
OY = \epsilon_A e^{i(\phi_A + \gamma)} YA,
$$
 (9)

implying a rate asymmetry between  $B^+ \rightarrow K^0 \pi^+$  and  $B^ \rightarrow \bar{K}^0 \pi^-$ .

For illustration, we take  $\gamma = 76^{\circ}, \epsilon = 0.24$  [5],  $\epsilon_A = 0.1$ (which is a reasonable guess [10,11]), and we vary  $\phi$  and  $\phi_A$ in the range  $0^{\circ} \le \phi \le 180^{\circ}$ ,  $-90^{\circ} \le \phi_A \le 270^{\circ}$ . The results of a search for solutions in the interval  $65^{\circ} \le \gamma \le 90^{\circ}$  are presented in Fig. 2 which displays a twofold ambiguity. Figure 2(a) shows the solution as function of  $\phi_A$  for two values of  $\phi$ ,  $\phi$ =60° and  $\phi$ =90°. Whereas for  $\phi$ <sub>*A*</sub>=90° the solution is very close to the input value, the deviation becomes maximal for  $\phi_A = 0^\circ, 180^\circ$ . This agrees with the geometry of Fig. 1, in which the largest rescattering effects are expected when *Y O* is parallel or antiparallel to the line bisecting the angle *BYA*.

In a second plot, Fig. 2(b), we fix  $\phi_A = 0^\circ$  and vary  $\phi$ over its entire range, which illustrates the maximal rescattering effect. We find two branches of the solution for  $\gamma$ , both of which deviate strongly from the input value  $\gamma=76^{\circ}$  for values of  $\phi$  around 90°. At  $\phi = 90^\circ$  there is no solution for  $\epsilon_A$ =0.1 in the considered interval. We checked that the so-

FIG. 2. The weak phase  $\gamma$  is obtained as the solution to the equation  $cos(2\gamma)=cos(BOA)$ . (a) the dependence of the solution on  $\phi_A$ , for two values of  $\phi$ =60° and  $\phi$ =90°; (b) the dependence of the solution on  $\phi$ , for  $\phi_A = 0^{\circ}$  (both graphs correspond to  $\epsilon_A = 0.1$ ,  $\gamma$  $=76^{\circ}$ ).

lution is restored and approaches the input value as the magnitude of  $\epsilon_A$  decreases to zero, as it should. Thus, the uncertainty in  $\gamma$ , seen both in Fig. 2(a) and Fig. 2(b) at  $\phi_A = 0^\circ$ and around  $\phi=90^{\circ}$ , is about 14°. It can even be worse in the singular cases where no solution for  $\gamma$  can be found.

A variant of this method for determining  $\gamma$ , proposed recently in Ref.  $[8]$ , was formulated in terms of two quantities  $R_*$  and  $\tilde{A}$  defined by

$$
R_{*} = \frac{B(B^{\pm} \to K^{0} \pi^{\pm})}{2B(B^{\pm} \to K^{\pm} \pi^{0})},
$$
  
\n
$$
\tilde{A} = \frac{B(B^{\pm} \to K^{\pm} \pi^{0}) - B(B^{-} \to K^{-} \pi^{0})}{B(B^{\pm} \to K^{0} \pi^{\pm})}
$$
  
\n
$$
-\frac{B(B^{\pm} \to K^{0} \pi^{+}) - B(B^{-} \to \bar{K}^{0} \pi^{-})}{2B(B^{\pm} \to K^{0} \pi^{\pm})}.
$$
 (10)

These quantities do not contain  $\mathcal{O}(\epsilon_A)$  terms; their dependence on the rescattering parameter  $\epsilon_A$  appears only at order  $\mathcal{O}(\epsilon \epsilon_A)$ . Therefore, it was argued in Ref. [8], the determination of  $\gamma$ , by setting  $\epsilon_A = 0$  in the expressions for  $R_*$  and  $\tilde{A}$ , is insensitive to rescattering effects. This procedure gives two equations for  $\gamma$  and  $\phi$  which can be solved simultaneously from  $R_*$  and  $\tilde{A}$ . Using two pairs of input values for (*R*\* ,*<sup>A</sup> ˜*) <sup>~</sup>corresponding to a restricted range for  $\phi_A$  and  $\phi$ ) seemed to indicate that the error in  $\gamma$  for  $\epsilon_A$  $=0.08$  is only about 5°. (The relations between the parameters used in Ref. [8] and ours are  $\phi = -\phi$ ,  $\eta = \phi_A$  $+\pi$ ,  $\bar{\epsilon}_{3/2} = \epsilon$ , and  $\epsilon_a = \epsilon_A$ .)

In Fig. 3 we show the results of such an analysis carried out for the entire parameter space of  $\phi_A$  and  $\phi$ . Whereas the

> FIG. 3. (a) the weak phase  $\gamma$ extracted from the method using the parameters  $(R_{*}, \tilde{A})$ , as a function of the strong phase  $\phi$  for several values of  $\phi_A(\epsilon_A=0.1)$ . The horizontal line shows the assumed physical value of  $\gamma=76^{\circ}$ . (b) the strong phase  $\phi$  can be reconstructed using the  $(R_*, \tilde{A})$  data.





FIG. 4. Geometric construction for the method described in Sec. III. The lines *OC* and *OD* denote the amplitudes  $A(B^{\pm} \rightarrow K^{\pm} K^0)$ , normalized as described in text. The positions of the points *C* and *D* are found as intersection points of the lines *AY* and *BY* with the two circles of radii given by  $|A(B^{\pm} \rightarrow K^{\pm} K^0)|$ .

angle  $\phi$  can be recovered with small errors, the results for  $\gamma$ show the same large rescattering effects for values of  $\phi$ around  $90^{\circ}$  as in Fig. 2. (A slight improvement is the absence of a discrete ambiguity in the value of  $\gamma$ .) These results show that the large deviation of  $\gamma$  from its physical value for  $\phi = 90^\circ$  is a general phenomenon, common to all variants of this method. Some information about the size of the expected error can be obtained by first determining  $\phi$ . Values not too close to 90° would be an indication for a small error.

### **III.** ELIMINATING RESCATTERING BY  $B^{\pm} \rightarrow K^{\pm} K^0$

The amplitude for  $B^+ \to K^+ \overline{K}^0$  is obtained from  $A(B^+)$  $\rightarrow$ *K*<sup>0</sup> $\pi$ <sup>+</sup>) in Eq. (1) by a *U*-spin rotation [10]

$$
A(B^+\rightarrow K^+\overline{K}^0) = |\lambda_u^{(d)}|e^{i\gamma}(A+P_{uc})
$$
  
 
$$
+ |\lambda_t^{(d)}|e^{-i\beta}(P_{ct}+P_3^{\text{EW}}).
$$
 (11)

In the limit of  $SU(3)$  symmetry the amplitudes in Eq.  $(11)$ are exactly the same as those appearing in Eq.  $(1)$ . In Fig. 4  $A(B^+\rightarrow K^+\overline{K}^0)$ , scaled by the factor  $\lambda/(1-\lambda^2/2)$  (and divided by  $A$  as in Fig. 1), is given by the line  $OC$  and its charge-conjugate is given by *OD*. We have shown in Ref. [16] that knowledge of these two amplitudes allows one to completely eliminate the rescattering contribution  $A + P_{uc}$ from the determination of  $\gamma$ . This is achieved by effectively replacing in the GLRN method the origin *O* by the intersection *Y* of the lines *AC* and *BD*.  $\gamma$  is determined by requiring that the angle  $(2\gamma)$  between *YA* and *YB* is equal to the angle  $(2\gamma)$  at the center of the circle.

The amplitude  $(11)$  can be decomposed into two terms carrying definite weak phases in form very similar to Eq.  $(6)$ ,

$$
\frac{\lambda}{1 - \lambda^2/2} A(B^+ \to K^+ \overline{K}^0) = -V_{cb} \left(1 - \frac{\lambda^2}{2}\right) p
$$

$$
\times \left(-\frac{\lambda^2}{(1 - \lambda^2/2)^2} + \epsilon_A e^{i\phi_A} e^{i\gamma}\right),\tag{12}
$$

The ratio  $|CY|/|AY| = \lambda^2/(1-\lambda^2/2)^2$  implies that the triangle *AYB* is about 25 times larger than the triangle *CYD*. This will result in a large uncertainty in  $\gamma$  also when the equality between the corresponding terms in  $B^+ \to K^0 \pi^+$ and  $B^+ \rightarrow K^+ \overline{K}^0$  amplitudes involves relatively small SU(3) violation.

The geometrical construction by which rescattering amplitudes can be completely eliminated in the  $SU(3)$  limit consists of three steps. (See Fig. 4. For an alternative suggestion, see Ref. [12].)

 $(a)$  Determine the position of the point *Y* as a function of the variable angle  $2\gamma$  and the decay rates of  $B^{\pm} \rightarrow K \pi$  and  $B^+\rightarrow \pi^+\pi^0$ . The point *Y* is chosen on the midperpendicular of *AB* such that the equality of the angles marked  $2\gamma$  is preserved for any value of  $\gamma$ .

(b) Draw two circles of radii  $\lambda/(1-\lambda^2/2)|A(B^{\pm})|$  $\rightarrow$ *K*<sup>0</sup>*K*<sup> $\pm$ </sup>) centered at the origin *O* (dashed-dotted circles in Fig. 4). The intersections of the lines *AY* and *BY* with these circles determine  $C$  and  $D$ , respectively (up to a twofold ambiguity), again as functions of  $\gamma$ .

(c) The physical value of  $\gamma$  is determined by the requirement  $|AC| = |BD|$  [16]. This condition on  $\gamma$  can be formulated in an algebraic form, showing that only the chargeaveraged rate of  $B^{\pm} \rightarrow K^{\pm} K^{0}$  is needed. The condition is given by Eq.  $(A1)$  in the Appendix.

Let us examine the precision of this method for  $\epsilon_A = 0.1$  at  $\phi \approx 90^{\circ}$ , for which the simpler method of Sec. II receives large rescattering corrections. In Fig.  $5(a)$  we show the lefthand side of Eq. (A1) as a function of variable  $\gamma$  at  $\phi$ =90° for several values of  $\phi_A$ . The value of  $\gamma$  is obtained from the condition that the left-hand side of this equation vanishes. In the absence of  $SU(3)$  breaking this method re-



FIG. 5. (a) The left-hand of Eq.  $(A1)$  as a function of variable  $\gamma$  for  $\phi=90^\circ$  and for different values of  $\phi_A$ . All these curves intersect at  $\gamma=76^{\circ}$ , which is the assumed physical value.  $(b)$  SU $(3)$ breaking effects introduce an error on the extracted value of  $\gamma$ , here shown as function of  $\phi_A$  at  $\phi$  $=90^\circ$ .

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produces precisely the physical value of  $\gamma$  ( $\gamma$ =76°) for all values of  $\phi_A$ . However, SU(3) breaking effects can become important, to the point of completely spoiling this method. We simulate these effects by taking the amplitudes *p* and *a*  $\equiv$ *A* + *P<sub>uc</sub>* - *p* in  $B^{\pm}$   $\rightarrow$   $K^{\pm}$   $K^0$  [Eq. (11)] to differ by at most 30% from those in  $B^{\pm} \rightarrow K^0 \pi^{\pm}$  [Eq. (1)]. This expands the lines of Fig.  $5(a)$  into bands of finite width, which give a range for the output value of  $\gamma$ .

In Fig.  $5(b)$  we show the effects of  $SU(3)$  breaking on the determination of  $\gamma$  as a function of  $\phi_A$  for  $\phi=90^\circ$ . We see that for values of  $|\phi_A|$  larger than about 25° the error on  $\gamma$  is quite large. Thus, we conclude that for certain values of the strong phases the determination of  $\gamma$  using this method is unstable under SU(3) breaking in the relation between  $B^+$  $\rightarrow$ *K*<sup>0</sup> $\pi$ <sup>+</sup> and *B*<sup>+</sup> $\rightarrow$ *K*<sup>+</sup> $\overline{K}$ <sup>0</sup>.

## **IV.** THE USE OF  $B^{\pm} \rightarrow \pi^{\pm} \eta_8$

In Ref. [16] we proposed to use in addition to  $B^+$  $\rightarrow$ *K*<sup>+</sup> $\bar{K}$ <sup>0</sup> also *B*<sup>+</sup> $\rightarrow$  $\pi$ <sup>+</sup> $\eta_8$  and their charge conjugates. Writing

$$
A(B^{+} \to \pi^{+} \eta_{8}) = |\lambda_{u}^{(d)}| e^{i\gamma} (-T - C - 2A - 2P_{uc})
$$
  
 
$$
+ |\lambda_{t}^{(d)}| e^{-i\beta} (-P_{ct} + P_{5}^{EW}), \qquad (13)
$$

we find the triangle relation

$$
A(B^{+} \to K^{+} \bar{K}^{0}) + \sqrt{\frac{3}{2}} A(B^{+} \to \pi^{+} \eta_{8})
$$
  
=  $\frac{1}{\sqrt{2}} A(B^{+} \to \pi^{+} \pi^{0}).$  (14)

This relation and its charge conjugate provide another condition which determines the positions of the points *C* and *D*. As in Sec. III, the phase  $\gamma$  is determined by the equation  $cos(BYA) = cos 2\gamma$ , where the point *Y* is fixed by the intersection of the lines *AC* and *BD*. General considerations, based on the relative sizes of the amplitudes involved, suggest that this method is relatively insensitive to  $SU(3)$  breaking effects  $\lceil 16 \rceil$ .

We illustrate this in Fig. 6 where we show on the same plot the two sides of the equation  $cos(BYA) = cos 2\gamma$  as functions of the variable  $\gamma$ . As in the method of Sec. III, SU(3) breaking is simulated by taking the penguin  $(p)$  and annihilation (*a*) amplitudes in  $B^{\pm} \rightarrow K^{\pm} K^0$  to differ by at most 30% (separately for their real and imaginary parts) from those in  $B^{\pm} \rightarrow K^{0} \pi^{\pm}$ . The latter are used to construct the positions of the points *C* and *D*. In the example of Fig. 6 we take  $\epsilon_A$  $=0.1,\phi=90^\circ,\phi_A=45^\circ$ , for which the two methods described in Secs. II and III were shown to lead to large errors in  $\gamma$ . For an input value  $\gamma=76^{\circ}$ , the output is given by the range  $74^{\circ} < \gamma < 78^{\circ}$ , obtained by the intersection of the solid line with the band formed by the diamond points. We see that the error in  $\gamma$  due to SU(3) breaking is less than  $\pm 2^{\circ}$ , which confirms the general arguments of Ref.  $[16]$ . This scheme, or rather its analogous version using  $B^0$  and  $B_s$  decay [16], may prove useful for a determination of  $\gamma$  in case



FIG. 6. Numerical results for the method of Sec. IV. The two sides of the equation  $cos(BYA)=2\gamma$  as function of variable  $\gamma$ , including 30% SU(3) breaking effects in the  $p$  and  $a$  amplitudes. The physical value of  $\gamma$  is determined by the intersection of the solid line with the wide band. The strong phases are taken as  $(\phi, \phi_A)$  $= (90^{\circ}, 45^{\circ}).$ 

that the strong phases  $(\phi, \phi_A)$  turn out to have values which preclude the use of the two simpler methods.

## **V. CONCLUSION**

We compared three ways of dealing with rescattering effects in  $B^{\pm} \rightarrow K^0 \pi^{\pm}$ , in order to achieve a precise determination for the weak phase  $\gamma$  from these processes and  $B^{\pm}$  $\rightarrow K^{\pm} \pi^{0}$ . In the simplest GLRN method which neglects rescattering we find that large errors in  $\gamma$  are possible for a particular region of the strong phases,  $\phi \sim 90^{\circ}$ , even when the rescattering term is only at a level of 10%. Limits on rescattering at this level are attainable from  $B^{\pm} \rightarrow K^{\pm} K^{0}$ .  $B^{\pm}$ and  $B^-$  decay rate measurements into  $K\pi$  are expected to provide rather precise information on  $\phi$ . In the likely case that  $\phi$  turns out to be far away from 90° small errors in  $\gamma$ would be implied.

On the other hand, in the less likely case that values of  $\phi$ are measured near 90° one may try to eliminate the rescattering effects by using also the charge-averaged  $B^{\pm}$  $\rightarrow$ *K*<sup> $\pm$ </sup>*K*<sup>0</sup> rate. This method suffers from a sizable uncertainty due to  $SU(3)$  breaking. These uncertainties could be resolved by also measuring  $B^{\pm} \rightarrow \pi^{\pm} \eta_8$ . The effect of  $\eta - \eta'$  mixing requires further study. Alternatively, to avoid this effect, one can apply the same method using  $B^0$  and  $B_s$  decays.

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#### **APPENDIX**

The weak angle  $\gamma$  is fixed in the method described in Sec. III by the condition  $|AC| = |BD|$ , or equivalently  $|YC|$   $=$ [*YD*]. Explicitly, this can be written after some algebra as an equation in  $\gamma$ 

$$
2(1-x_0)^2 \vec{Y}^2 + 2x_0(1-x_0) \vec{Y} \cdot (\vec{A} + \vec{B}) + x_0^2 (x_0^2 + \tilde{x}_{0-}^2)
$$
  
 
$$
-(y_{+0}^2 + \tilde{y}_{-0}^2) = 0,
$$
 (A1)

where  $x_0$  is defined as the ratio of two  $CP$  rate differences

$$
x_0 \equiv (y_{+0}^2 - \tilde{y}_{-0}^2)/(x_{0+}^2 - \tilde{x}_{0-}^2) \to -\lambda^2/(1 - \lambda^2/2)^2.
$$
\n(A2)

Here

$$
y_{+0} = \frac{1}{\sqrt{2}} \frac{f_{\pi}}{f_K} \frac{|A(\bar{B}^+ \to K^+ \bar{K}^0)|}{|A(B^+ \to \pi^+ \pi^0)|},
$$
  

$$
\tilde{y}_{-0} = \frac{1}{\sqrt{2}} \frac{f_{\pi}}{f_K} \frac{|A(\bar{B}^- \to K^- K^0)|}{|A(B^+ \to \pi^+ \pi^0)|}
$$
(A3)

obey an SU(3) relation with the amplitudes (5) of  $B^{\pm}$  $\rightarrow$ *K* $\pi$ <sup> $\pm$ </sup>

$$
y_{+0}^2 - \tilde{y}_{-0}^2 = -\frac{\lambda^2}{(1 - \lambda^2/2)^2} (x_{0+}^2 - \tilde{x}_{0-}^2).
$$
 (A4)

This implies that *CP* rate differences in  $B^{\pm} \rightarrow K^{0} \pi^{\pm}$  and  $B^{\pm} \rightarrow K^{\pm} K^{0}$  are equal and of opposite sign [12]. We see that in the  $SU(3)$  limit the condition  $(A1)$ , which eliminates rescattering effects, requires only a measurement of the charge-averaged rate of  $B^{\pm} \rightarrow K^{\pm} K^{0}$  and not the *CP* asymmetry in these processes  $[12]$ .

To prove Eq. (A1), let us consider two lines *AY* and *BY* cutting two circles of radii  $R_1$ ,  $R_2$  (centered at the origin) at points *C* and *D*, respectively. The intersection points can be written as  $\vec{C} = \vec{Y} + x_1(\vec{A} - \vec{Y})$  and  $\vec{D} = \vec{Y} + x_2(\vec{B} - \vec{Y})$ , where  $\rightarrow$  $\rightarrow$  $\rightarrow$  $\rightarrow$  $\rightarrow$  $x_1, x_2$  are solutions of the equations

$$
(\vec{A} - \vec{Y})^2 x_1^2 + 2x_1 \vec{Y} \cdot (\vec{A} - \vec{Y}) + (\vec{Y}^2 - R_1^2) = 0,
$$
 (A5)

$$
(\vec{B} - \vec{Y})^2 x_2^2 + 2x_2 \vec{Y} \cdot (\vec{B} - \vec{Y}) + (\vec{Y}^2 - R_2^2) = 0.
$$
 (A6)

The condition  $|YC| = |YD|$  is equivalent to requiring that these two equations have a common solution  $x_1 = x_2$ . Obviously, if such a solution exists, it is given by

$$
x_0 = \frac{R_1^2 - R_2^2}{2\vec{Y} \cdot (\vec{A} - \vec{B})} = \frac{R_1^2 - R_2^2}{\vec{A}^2 - \vec{B}^2},
$$
 (A7)

where we used the equality  $(\vec{A} - \vec{Y})^2 = (\vec{B} - \vec{Y})^2$ . Taking the sum of Eqs.  $(A5)$  and  $(A6)$  with the value  $(A7)$  for *x* leads immediately to the condition  $(A1)$ .

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