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# Predictions for leptonic mixing angle correlations and nontrivial Dirac CP violation from $A_5$ with generalized CP symmetry

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Using a discrete flavor group,  $A_5$ , combined with generalized CP, we study the mixing parameter correlations which arise from breaking to residual symmetries in the neutrino  $G_{\nu} = \mathbb{Z}_2 \times CP$ , and the charged lepton sectors  $G_e = \mathbb{Z}_2$ . By focusing on patterns that agree with current experimental data we demonstrate that nontrivial leptonic phases are predicted and discuss a number of distinctive correlations between mixing parameters.

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### I. INTRODUCTION

The mixing structure of the three families of the lepton sector has inspired many to use non-Abelian flavor symmetries to predict the mixing angles and phases of the Pontecorvo-Maki-Nakagawa-Sakata (PMNS) matrix: three mixing angles,  $\theta_{13}$ ,  $\theta_{12}$ ,  $\theta_{23}$ , the *CP*-violating phase  $\delta$  and Majorana phases  $\alpha_{21}$  and  $\alpha_{31}$ . A significant experimental result in 2012 was the measurement of  $\theta_{13}$  [1–3]. This somewhat large value of  $\theta_{13}$  ruled out many classes of flavor models that predicted small or zero  $\theta_{13}$ . These models lead to simple mixing patterns such as bimaximal [4–7], tribimaximal [8] and golden ratio [9] which result from using small flavor groups such as  $A_4$ ,  $A_5$  and  $S_4$  (an indepth review of discrete groups can be found in Ref. [10]).

In order to produce mixing patterns that accommodate experimental data, the flavor model paradigm has shifted to include larger groups such as  $\Delta(96)$  [11,12],  $\Delta(150)$  [13],  $\Delta(600)$  [14] and  $\Delta(1536)$  [15]. These non-Abelian discrete flavor groups cannot be a symmetry at the low-energy scale as leptonic masses are distinct. Therefore the flavor group must be broken into Abelian residual symmetries in the neutrino and charged lepton sectors. The structure of the Abelian residual symmetries is shaped by the larger non-Abelian flavor group and from these low-energy residual symmetries leptonic observables can be predicted. In general, there are two possible implementations of flavor symmetries and they are often referred to as direct and semidirect (e.g. see Ref. [10]). The distinction between the two approaches is the low-energy residual symmetry of the Majorana mass matrix: in the direct approach the Klein group,  $\mathbb{Z}_2 \times \mathbb{Z}_2$ , is a subgroup of the underlying flavor symmetry while in the semidirect approach,  $\mathbb{Z}_2$  emerges as a residual symmetry of the flavor group. In the semidirect models a continuous parameter is introduced, derived from the freedom to rotate in the degenerate subspace of the neutrino residual symmetry, allowing the prediction of a nonzero  $\theta_{13}$ . There are several attractive features of implementing such an approach: first, a UV-complete theory is not necessary in order to predict leptonic observables [16–22]. Second, correlations between the observables can be derived and provide specific signatures which allow the comparison of a range of models to experimental data [18–25]. In contrast with the direct and semidirect models, there have been predictions of leptonic observables using  $Z_2$  residual symmetries which do not inherit structure from a non-Abelian subgroup [26,27].

These models can successfully predict mixing angles consistent with data and a Dirac phase,  $\delta$ . However, due to the constraints imposed on the mass matrices used to construct the PMNS matrix, a number of degrees of freedom cannot be eliminated and therefore these models cannot predict Majorana phases. By extending the flavor group to include generalized CP (gCP) symmetry, the three mixing angles and three phases can be determined using a small number of input parameters [28]. This idea of combining CP with a flavor symmetry is not a recent one and was originally discussed in [29–31] together with a  $\mu - \tau$  symmetry. There have been a number of interesting works on the consistent relation between gCP and flavor symmetry [28,32,33] and many plausible groups have been studied such as  $A_4$  [28,34],  $S_4$  [28,35,36],  $\Delta$ (96) [11,12],  $\Delta(150)$  [13],  $\Delta(600)$  [14],  $\Delta(1536)$  [15],  $\Delta(3n^2)$  [37],  $\Delta(6n^2)$  [37,38] and most recently  $A_5$  [39–41]. In smaller groups such as  $A_4$ ,  $S_4$  and  $A_5$  [28,34,35,39–41], it has been found that the leptonic phases are either trivial or maximal. Moreover, there are often share recurring patterns of predictions such as maximally CP-violating  $\delta$  associated with maximal  $\theta_{23}$ , the origin of which was recently discussed [42,43]. Applying the same framework with a larger flavor group such as  $\Delta(3n^2)$  or  $\Delta(6n^2)$ , leptonic phases are nontrivially dependent upon the continuous parameter and can take values different from  $0, \frac{\pi}{2}, \pi$  and  $\frac{3\pi}{2}$ . In addition to the leptonic observable predictions made

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using non-Abelian flavor groups combined with gCP, there have been interesting recent works which apply an invariant approach to *CP* in flavor symmetries [44–46].

The work presented in this paper is an extension of the study in [41], where a flavor group,  $A_5$ , combined with gCP is broken into residual symmetries in the charged lepton,  $G_e = \{\mathbb{Z}_3, \mathbb{Z}_5, \mathbb{Z}_2 \times \mathbb{Z}_2\}$ , and the neutrino sectors,  $G_{\nu} = \mathbb{Z}_2 \times CP$ . An additional motivation to further explore the predictions of  $A_5$  is, unlike other small groups such as  $A_4$  and  $S_4$ ,  $S_6$  is anomaly safe [47]. In this work, we consider the possibility that the same high-energy symmetry breaks into low-energy residual symmetries  $S_e = \mathbb{Z}_2$  and  $S_e = \mathbb{Z}_2 \times CP$ . By relaxing the possible combination of residual symmetries we find that nontrivial values of the leptonic phases can be accommodated and there are distinctive correlations between observables.

Throughout this work, we assume our low-energy effective theory is the Standard Model augmented by a Majorana mass term and we will use the following  $3\sigma$  global fit data [48]:

$$7.85^{\circ} \le \theta_{13} \le 9.10^{\circ},$$
  $31.29^{\circ} \le \theta_{12} \le 35.91^{\circ},$   $38.2^{\circ} \le \theta_{23} \le 53.3^{\circ}.$ 

The work presented in this paper is structured as follows: in Sec. II we present the assumptions of our theoretical framework; in Sec. III we discuss the construction of the PMNS matrix from a symmetry constraint and the derivation of our results. A number of representative predictions, along with an example, are given in Sec. IV and finally we make concluding remarks in Sec. V.

### II. SYMMETRIES OF THE MODEL

We briefly review the theoretical framework of this study, where we have closely followed the discussion of [41]. We first review the general concepts of flavor and generalized *CP* symmetry and subsequently consider the consistent relations between these two symmetry transformations in preparation for constructing the PMNS matrix.

### A. Flavor symmetry

We assume there exists a finite, discrete flavor symmetry,  $G_f$ , at the high-energy scale. The purpose of this symmetry is to unify the three flavors of leptonic doublets into a single mathematical object: a three-dimensional irreducible representation of the flavor group,  $\Psi$ . The flavor group acts on  $\Psi$  such that

$$\Psi \to \rho(g)\Psi,$$
 (1)

where  $\rho(g)$  is a three-dimensional unitary representation of group element  $g \in G_f$ . The non-Abelian flavor symmetry must be broken at the low-energy scale as leptonic masses

are distinct. This implies that if a flavor symmetry is operational in the high-energy regime then only its Abelian residual symmetries would be observable at the scale of mass generation. Therefore, we assume that the non-Abelian flavor symmetry is broken into Abelian residual symmetries in the charged lepton sector,  $G_e$ , and the neutrino sector,  $G_{\nu}$ . For group elements  $g_e \in G_e$  and  $g_{\nu} \in G_{\nu}$ , the charged lepton and neutrino fields transform under the residual symmetries according to

$$e_{\rm L} \to \rho(g_e)e_{\rm L}$$
 and  $\nu_{\rm L} \to \rho(g_{\nu})\nu_{\rm L}$ , (2)

where generational indices have been suppressed. The transformations of Eq. (2) enforce constraints on the charged lepton and neutrino mass matrices,

$$\rho(g_e)^{\dagger}(m_e m_e^{\dagger})\rho(g_e) = m_e m_e^{\dagger}, \tag{3}$$

$$\rho(g_{\nu})^{\mathrm{T}} m_{\nu} \rho(g_{\nu}) = m_{\nu}. \tag{4}$$

To deduce the possible forms of the residual symmetries, we must consider the largest symmetry of each sector and the structure inherited from the larger non-Abelian flavor group. On the basis in which the charged lepton mass matrix is diagonal and the masses are distinct, the largest symmetry of this sector is  $U(1)^3$ . This is derived from the freedom to rephase the fields of each generation of the charged leptons. The most general discrete residual symmetry of this sector must be a subgroup of  $U(1)^3$  and is therefore a direct product of the cyclic groups,  $\mathbb{Z}_n$ . The Abelian subgroups of  $A_5$  that satisfy this condition are  $\mathbb{Z}_5$ ,  $\mathbb{Z}_3$ ,  $\mathbb{Z}_2 \times \mathbb{Z}_2$  and  $\mathbb{Z}_2$ , where the cases of  $\mathbb{Z}_3$ ,  $\mathbb{Z}_5$  and  $\mathbb{Z}_2 \times$  $\mathbb{Z}_2$  have been studied in the analysis of [41]. As we assume that the neutrinos are Majorana in nature, rather than Diractype particles, their mass matrix is always invariant under a Klein symmetry,  $\mathbb{Z}_2 \times \mathbb{Z}_2$ . Therefore the residual symmetry of the neutrino sector is the Klein group or a subgroup thereof.

### B. Generalized *CP* symmetry

In addition to the non-Abelian flavor symmetry operational at the high-energy scale, we assume there also exists a gCP symmetry. This symmetry parity transforms and charge conjugates the field, as well as acting on its generational indices [49]. The gCP transformation acts on the multiplet of fields,  $\Psi$ , as

$$\Psi \to X \Psi^C,$$
 (5)

where X is a unitary, symmetric matrix and  $\Psi^C$  denotes the CP conjugate of  $\Psi$ . We have chosen gCP to be an involutory, meaning that two gCP transformations are equivalent to the identity:  $XX^* = 1$ . If gCP remains a symmetry of the charged lepton or neutrino sector, it must leave the mass terms invariant:

$$X^{\mathrm{T}} m_{\nu} X = m_{\nu}^*, \tag{6}$$

$$X^{\dagger}(m_e m_e^{\dagger}) X = (m_e m_e^{\dagger})^*. \tag{7}$$

It has been demonstrated [35,50] that if gCP remains unbroken at the low-energy scale in both the charged lepton and neutrino sectors, then Eqs. (6) and (7) are satisfied and consequently no *CP*-violating effects will be observed. Therefore, in this paper we assume that gCP is broken in the charged lepton sector and remains a preserved symmetry of the neutrino sector.

In summary, the residual flavor and gCP symmetries place a series of constraints on the charged lepton and neutrino mass matrices. These constraints shape the form of the diagonalizing matrices,  $U_e$  and  $U_\nu$ , which in turn constrain the form of the PMNS matrix. The consistent interaction between the non-Abelian flavor group and gCP must be considered in order to determine which gCP transformations are physical.

### C. Combining flavor and gCP symmetries

To illustrate the consistency between the flavor and gCP symmetries, consider a multiplet of fields transforming under a gCP, flavor and subsequent gCP transformation,

$$\Psi \to X \rho(g)^* X^* \Psi \equiv \rho(g') \Psi. \tag{8}$$

In [28,32], the authors showed that in order to combine flavor and gCP symmetries consistently, gCP must act as an automorphism on the flavor group,  $G_f$  [in Eq. (8), gCP maps group element g to another element g such that the identity and group multiplication is respected]. This idea was further developed by [33], which pointed out that for a physical CP transformation to occur, gCP should be an outer automorphism that maps the representations of the fields to their conjugate representations. Moreover, in a generic setting, these outer automorphisms of the group must be class-inverting automorphisms, implying that Eq. (8) becomes

$$X\rho(g)^*X^*\Psi \equiv \rho(h)\rho(g^{-1})\rho(h)^{-1}\Psi, \tag{9}$$

for  $h \in G_f$ . The detailed derivation of X is discussed fully in [41], however in this work we will briefly summarize the group theoretic concepts that were considered. In order to find the forms of X that constitute physical gCP transformations, the outer automorphism of the flavor group must be known. In general, the outer automorphism group for  $A_n$ , where  $n \le 5$ , is  $\mathbb{Z}_2$  (for more group theory insights see Ref. [51]). This implies there is only one nontrivial outer automorphism of  $A_5$  and this maps elements of one conjugacy classes of order 5 to the other. In addition to

finding the nontrivial outer automorphism of the flavor group,  $A_5$  has a special property that simplifies the derivation of X.  $A_5$  is an ambivalent group, meaning each element is conjugate to its inverse. Applying this property of the group and choosing to work in a real representation, Eq. (9) can be significantly simplified. Henceforth, a series of deductions can be made and it can be concluded that the forms of X that act on  $A_5$  as class-inverting, involutory automorphism are the Klein group.

### III. METHODOLOGY

We first discuss the construction of the PMNS matrix from the symmetry constraints and subsequently describe the method used to derive the correlations between observables.

# A. Constructing the PMNS matrix from symmetry considerations

The flavor and gCP symmetry constrain the form of the neutrino and charged lepton mass matrices. From these constraints the form of their diagonalizing matrices,  $U_{\nu}$  and  $U_{e}$ , may be deduced and thus the PMNS matrix can be constructed:  $U_{\rm PMNS} = U_{e}^{\dagger}U_{\nu}$ . Let us consider how to derive  $U_{e}$  from the symmetry constraints. First, Eq. (3) can be reexpressed in the form of a commutator:  $[\rho(g_{e}), (m_{e}m_{e}^{\dagger})] = 0$ . As the unitary representation  $\rho(g_{e})$  commutes with the Hermitian matrix  $m_{e}m_{e}^{\dagger}$ , there exists a unitary matrix,  $U_{e}$ , that simultaneously diagonalizes both. In the case where  $\rho(g_{e})$  has degenerate eigenvalues, there is not a unique diagonalizing matrix of  $\rho(g_{e})$  but rather an additional complex rotation can be performed in the degenerate subspace of  $\rho(g_{e})$ . Therefore, the most general form of the diagonalizing matrix of  $\rho(g_{e})$  and  $(m_{e}m_{e}^{\dagger})$  is

$$U_e = U_l R(\omega, \gamma), \tag{10}$$

where  $U_l$  diagonalizes  $\rho(g_e)$  and  $R(\omega, \gamma)$  is an SU(2) transformation in the degenerate eigenspace. It is worth stressing that we allow for the existence of this complex rotation by permitting  $G_e = \mathbb{Z}_2$ . This differs from the analysis of [41], as their choice of  $G_e$  had no such degenerate subspace.

In order to deduce the form of  $U_{\nu}$ , we consider constraints from the flavor residual symmetry, gCP and the logical relation between the two symmetries. The action of gCP on the neutrino residual symmetry,  $\rho(g_{\nu})$ , maps these elements to their inverse:

$$X\rho(g_{\nu})^*X^* = \rho(g_{\nu})^{-1} = \rho(g_{\nu}),$$
 (11)

where in the final step we have used the fact  $\mathbb{Z}_2$  elements are self-inverse. Equation (11) can equivalently be viewed as forming a direct product between  $\mathbb{Z}_2$  and gCP. In [28] they showed that it is always possible to make a convenient basis change  $\Omega$  ( $X = \Omega\Omega^T$ ) such that

 $<sup>^{1}</sup>A_{5}$  contains five conjugacy classes: one for order 1, 2 and 3 elements and two for order 5 elements.

$$(\Omega^T m_{\nu} \Omega) = (\Omega^T m_{\nu} \Omega)^*. \tag{12}$$

Therefore, this basis transformation ensures that  $m_{\nu}$  is real valued. Moreover, from Eq. (4), it can be seen that the diagonal from of  $\rho(g_{\nu})$  commutes with  $(\Omega^T m_{\nu}\Omega)$ , which implies that  $m_{\nu}$  must be block diagonal. To fully diagonalize the matrix of Eq. (12), an additional real rotation,  $R(\theta)$ , must be performed. From these considerations,  $U_{\nu}$  may be written as

$$U_{\nu} = \Omega R(\theta). \tag{13}$$

Using symmetry constraints alone, the PMNS matrix may be written as

$$U_{\rm PMNS} = R(\omega, \gamma) U_l^{\dagger} \Omega R(\theta). \tag{14}$$

For our chosen representation of  $A_5$ ,  $\Omega$  can take three possible forms,

$$\Omega_{12} = \begin{pmatrix} i & 0 & 0 \\ 0 & i & 0 \\ 0 & 0 & 1 \end{pmatrix}, \qquad \Omega_{13} = \begin{pmatrix} i & 0 & 0 \\ 0 & 1 & 0 \\ 0 & 0 & i \end{pmatrix} \quad \text{and} \quad \Omega_{23} = \begin{pmatrix} 1 & 0 & 0 \\ 0 & i & 0 \\ 0 & 0 & i \end{pmatrix}, \tag{15}$$

which have been fully derived in [41].

## B. Derivation of predictions for leptonic mixing parameters

In the work of [41], the combination of Abelian residual symmetries studied were  $G_e = \{\mathbb{Z}_3, \mathbb{Z}_5, \mathbb{Z}_2 \times \mathbb{Z}_2\}$  and  $G_{\nu} = \mathbb{Z}_2 \times CP$ . The observables are a function of one continuous parameter, derived from the freedom to make a real rotation in the degenerate subspace of the  $\mathbb{Z}_2$  residual symmetry in the neutrino sector. Moreover, the authors of [41] considered the combination of  $G_{\nu} = \mathbb{Z}_2 \times \mathbb{Z}_2 \times CP$ and  $G_e = \mathbb{Z}_2$ . In this case, each observable would be a function of two input parameters which is obtained from the ability to make an SU(2) transformation in the degenerate eigenspace of the  $\mathbb{Z}_2$  charged lepton residual symmetry. This was fully explored and there were no predictions that agreed with data at the  $3\sigma$  level. A further consideration of  $G_e = \mathbb{Z}_2$  and  $G_{\nu} = \mathbb{Z}_2 \times CP$  was proposed in [41]. Moreover, predictions that agreed to a  $3\sigma$ level with global fit data were found, however these correlations were not analyzed further.

In this work, we aim to develop on this combination of residual symmetries that was originally described in [41] and explore the nontrivial correlations between leptonic observables. This consideration allows for two additional continuous parameters and therefore each observable is a function of three input parameters:  $\theta$ ,  $\omega$  and  $\gamma$ .

For each of the 15 charged lepton residual symmetries,  $\rho(g_e)$ , we construct a diagonalizing matrix,  $U_e = U_l R(\omega, \gamma)$ , where we have fixed the degenerate subspace,  $R(\omega, \gamma)$ . To construct  $U_{\nu}$  we have explored all possible combinations of  $\Omega R(\theta)$  and from this, we can construct  $U_{PMNS}$ . For each constructed PMNS matrix, we considered the arbitrary ordering of the diagonalizing matrices and therefore appropriately permuted the rows and columns. For each PMNS matrix, the three continuous parameters  $(\theta, \omega, \gamma)$  are randomly scanned over the range  $[0,\pi]$  and the three mixing angles are calculated for each point in the phase space. Subsequently, only points that simultaneously agree to a  $3\sigma$  level with global fit data [48] are retained and from these points the phases are calculated. We have chosen to present our results in terms of correlations between the leptonic phases  $\theta_{13}$ ,  $\theta_{12}$  as a function of  $\theta_{23}$ .

### IV. RESULTS

There is a varied program of currently running and planned experiments that aim to increase precision in the measurement of a number of the oscillation parameters. In the near term, accelerator long-baseline experiments such as T2K [52] and NO $\nu$ A [53] aim at improving the current measurements of parameters such as  $\theta_{23}$ ,  $\delta$  and  $\Delta m_{32}{}^2$ . In the longer term, accelerator facilities such as DUNE [54] and T2HK [55] hope to further increase the sensitivity to these oscillation parameters. T2HK will have the ability to resolve  $\delta$  to a  $1\sigma$  uncertainty of 19° for all allowed values (using an integrated beam power of 7.5 MW seconds of exposure with  $1.56 \times 10^{22}$  protons on target). Moreover, using a 10 kt detector and expected knowledge from T2K and NO $\nu$ A would allow DUNE to achieve a  $3\sigma$  sensitivity for detecting CP violation in 50% of  $\delta$  values.

In conjunction, future medium baseline reactor experiments such as RENO-50 [56] and JUNO [57] aim to better the measurement of  $\theta_{12}$ . These experiments utilize the survival probability of electron antineutrinos, which are copiously produced in fission reactors, to determine the mass ordering and make subpercent measurements of  $\theta_{12}$ .

The determination of the nature of the neutrino remains of fundamental importance and neutrinoless double beta decay  $(\nu 0\beta\beta)$  experiments such as GERDA, CUORICINO, EXO-200 and KamLAND-Zen hope to explore *CP*-conserving upper boundary of the inverted ordering region. The decay rate of this rare process is proportional to the effective Majorana mass,  $m_{ee}$  (see e.g. [58–64]), and the values of  $m_{ee}$  are influenced by the combinations of phases,  $e^{i\alpha_{21}}$  and  $e^{i(\alpha_{31}-2\delta)}$ . We will comment on some specific predictions which have particularly relevant consequences for  $\nu 0\beta\beta$ . However, due to the ambitious plans to improve the measurement of  $\delta$ ,  $\theta_{12}$  and  $\theta_{23}$  by a range of complementary neutrino oscillation experiments, we will mainly focus on the mixing angle and  $\delta - \theta_{23}$  correlations.

Using the symmetry construction of Sec. II, each PMNS matrix is a function of three continuous parameters and therefore we find a large number of cases that agree to a  $3\sigma$  level with global fit data and we have in the order of 50 different predictions. For illustrative purposes we provide an explicit example of one such prediction in Sec. IV A. We group the remainder of our selected predictions into categories according to the octant of the  $\theta_{23}$ : the lower octant cases are discussed in Sec. IV B, upper octant in Sec. IV C and finally cases that span both octants are discussed in Sec. IV D.

### A. An example

We utilize the group representations of Refs. [65] and [66]. To construct  $U_e$ , let us first consider a  $\mathbb{Z}_2$  group element of  $A_5$  in the three-dimensional real representation

$$\mathbb{Z}_{2} = \frac{1}{2} \begin{pmatrix} -1 & \phi & \frac{-1}{\phi} \\ \phi & \frac{1}{\phi} & -1 \\ \frac{-1}{\phi} & -1 & -\phi \end{pmatrix}, \tag{16}$$

where  $\phi = \frac{(1+\sqrt{5})}{2}$  is the golden ratio. A diagonalizing matrix of Eq. (16) is

$$U_{l} \sim \begin{pmatrix} 0.665 & -0.555 & -\frac{1}{2} \\ -0.58 & -0.025 & -\frac{\phi}{2} \\ 0.461 & 0.832 & -\frac{1}{2\phi} \end{pmatrix}, \tag{17}$$

where the degenerate eigenvalues of the matrix of Eq. (16) are in the 12-plane. Therefore,  $U_e^{\dagger}$  takes the form

$$U_e^{\dagger} \sim \begin{pmatrix} c_{\omega} & e^{i\gamma} s_{\omega} & 0\\ -e^{-i\gamma} s_{\omega} & c_{\omega} & 0\\ 0 & 0 & 1 \end{pmatrix} \times \begin{pmatrix} 0.665 & -0.58 & 0.461\\ -0.555 & -0.025 & 0.832\\ -\frac{1}{2} & -\frac{\phi}{2} & -\frac{1}{2\phi} \end{pmatrix}, \quad (18)$$

where  $c_{\omega} \equiv \cos(\omega)$  and  $s_{\omega} \equiv \sin(\omega)$ . Not all combinations of  $\Omega$  and  $R(\theta)$  produce predictions within  $3\sigma$  of the global fit data, however, one such combination that does is

$$U_{\nu} = \begin{pmatrix} 1 & 0 & 0 \\ 0 & i & 0 \\ 0 & 0 & i \end{pmatrix} \begin{pmatrix} 1 & 0 & 0 \\ 0 & c_{\theta} & s_{\theta} \\ 0 & -s_{\theta} & c_{\theta} \end{pmatrix}. \tag{19}$$

Combining Eqs. (18) and (19), we construct the PMNS matrix and perform a random scan over the three continuous parameters  $(\theta, \omega, \gamma)$  in the interval  $[0, \pi]$ . The points of this parameter space that agree to a  $3\sigma$  level with global fit

data are retained and the phases are calculated. Throughout this work, the Particle Data Group parametrization has been applied to find the mixing angles and phases [67]. In Fig. 1, we plot the three leptonic phases,  $\theta_{12}$  and  $\theta_{13}$  as a function of  $\theta_{23}$ . This PMNS matrix requires  $44.2^{\circ} \le \theta_{23} \le 53.3^{\circ}$  and the  $\delta$  phase reaches a maximal value of 69° for large values of  $\theta_{23}$  (53°). A *CP*-conserving value of  $\delta$  is possible for all viable  $\theta_{23}$ 's and for the maximal  $\theta_{23}$ 's,  $-20^{\circ} \le \delta \le 20^{\circ}$ . From Fig. 1, there appears to be no preferred  $\delta$  phase within the viable parameter space. The values of the Majorana phases range from  $\alpha_{21} \sim \pm 25^{\circ}$  to  $\alpha_{31} \sim \pm 80^{\circ}$ . In the case of  $\alpha_{31}$ , small values (< 15°) are strongly preferred over the whole range of  $\theta_{23}$ . However for  $\theta_{23} > 49^{\circ}$ ,  $\alpha_{31}$  can take large values. The consequences of this prediction on neutrinoless double beta  $(\nu 0\beta\beta)$  decay would be interesting to explore. The magnitudes of  $\alpha_{21}$  and  $(\alpha_{31} - 2\delta)$  of this prediction can be small and this results in little cancellation between the mass terms of  $m_{ee}$ . This would imply the prediction for  $m_{ee}$  can be close to the CP-conserving upper boundary of the inverted ordering region, which experiments hope to explore. Therefore, it would be feasible to use  $\nu 0\beta\beta$  decay to study this particular prediction.

As the  $3\sigma$  range of  $\theta_{13}$  is highly constrained compared with the other mixing angles, there is little discernible structure in the  $\theta_{13} - \theta_{23}$  correlation, however the  $\theta_{12} - \theta_{23}$  dependence has greater predictivity. For near maximal values of  $\theta_{23}$ ,  $\theta_{12}$  is predicted to be at the very upper boundary of its  $3\sigma$  range (~36°). For larger values of  $\theta_{23}$  close to the upper  $3\sigma$  boundary, the range of predicted  $\theta_{12}$  increases (31.6°–35.9°). Although for most viable values of  $\theta_{23}$  there are a range of  $\theta_{12}$  predictions, the density of solutions clusters near the boundary of the viable region of the  $\delta - \theta_{23}$  parameter space.

### **B.** Lower octant predictions

The chosen lower octant predictions are presented in Fig. 2 and it can be seen that the possible range of  $\theta_{23}$ values differs between the various cases: lower octants (LOs) 1–3 have viable predictions for the entire lower octant (38.8°-45°) while LO 4 is somewhat more constrained, as  $\theta_{23}$  spans only 3° (40°-43°). LO 5 is the most highly constrained and therefore most easily testable, with  $38^{\circ} \le \theta_{23} \le 38.8^{\circ}$ . LOs 1–3 share the same  $\delta - \theta_{23}$  correlation, which attains a maximal  $\delta$  (85°) for a  $\theta_{23}$  close to the lower  $3\sigma$  allowed region. The *CP*-conserving values of  $\delta$ require  $40.6^{\circ} \le \theta_{23} \le 44.2^{\circ}$ . In the case of LO 4, although the  $\delta - \theta_{23}$  correlation structure is similar to that of LOs 1–3, the maximal value of  $\delta$  is slightly greater reaching 90° and solutions tend to cluster at these points. In contrast to LOs 1–4, the  $\delta$  value of LO 5 is close to zero, however it reaches a maximum of 26°.

LOs 1–3 have a common  $\theta_{12} - \theta_{23}$  dependence: for values of  $\theta_{23}$  close to the lower  $3\sigma$  boundary, all values of  $\theta_{12}$  are allowed. For a near maximal  $\theta_{23}$ , the  $\theta_{12}$  prediction becomes increasing constrained: for example for the

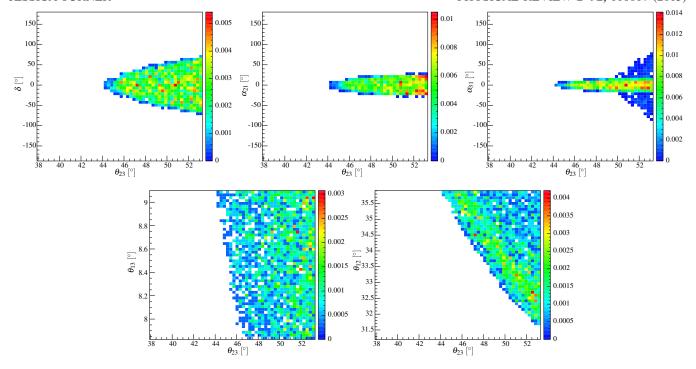


FIG. 1 (color online). Two-dimensional histograms showing the phases and mixing angles as a function of  $\theta_{23}$  for predictions that agree to a  $3\sigma$  level with global fit data [48]. The color chart shows relative frequencies of solutions where red (dark blue) represents a higher (lower) frequency.

current best fit value of  $\theta_{23}$  (42.3°) [48], only values of  $33.5^{\circ} \le \theta_{12} \le 35.91^{\circ}$  are predicted. In the case of LO 4, for  $38.8^{\circ} \le \theta_{23} \le 42.3^{\circ}$ , the predicted  $\theta_{12}$  spans the  $3\sigma$  range of  $\theta_{12}$ . Similar to LOs 1–3, the range of predicted  $\theta_{12}$  becomes more constrained for a near maximal  $\theta_{23}$  (smaller  $\theta_{12}$  is preferred). Using the  $\theta_{12} - \theta_{23}$  correlation as a means of differentiating between LOs 1–3 and LO 4 would be problematic in the regions  $\theta_{23} \le 41^{\circ}$ , as the predictions are indistinguishable. As the viable parameter space of LO 5 is significantly smaller than that of the previous four predictions, there is no discernible correlation between  $\theta_{12}$  and  $\theta_{23}$ ; in this regard its most discriminating feature is that  $\theta_{12}$  can only range between  $31^{\circ} - 33^{\circ}$ .

The Majorana phases are the only observables that differ amongst LOs 1–3. It is worth noting that LO 1 is the only lower octant prediction of this sample that has a CP-conserving value of  $\alpha_{21}$ . It would be an interesting future study to investigate the effect that this would have on  $\nu 0\beta \beta$  decay and the feasibility of discriminating between predictions.

In summary, there are several general features which are shared amongst these cases; the most striking of these is the prediction of nontrivial leptonic phases. Moreover, the  $\delta$  and  $\alpha_{21}$  phases are bound between  $\pm 90^\circ$ . If  $\delta$  is measured to be maximally CP violating, as hinted at by T2K [68], the only remaining viable prediction is LO 4. Some predictions cannot be discriminated between by using  $\delta$  and  $\theta_{12}$  alone and access to the Majorana phases is necessary. Moreover, the ability to discriminate between LO 4 and LOs 1–3 is

highly dependent upon the value of  $\theta_{23}$ : in the scenario of a maximal or near maximal  $\theta_{23}$ , this is possible. Of the cases presented, LO 5 is the most easily testable, as its  $\theta_{23}$  values are highly constrained and lie at the extreme lower boundary of the  $3\sigma$  range.

### C. Upper octant predictions

Similar to the lower octant results, we have chosen three cases [upper octants (UOs) 1–3], presented in Fig. 3, for which the mixing angle and  $\delta$  phase correlations are indistinguishable and only the Majorana phases differ. UOs 1–3 share the feature of viable predictions over the entire upper octant (45°–53.3°). The  $\theta_{23}$  allowed range UO 4 is slightly more constrained, with  $46.3^{\circ} \leq \theta_{23} \leq 53.3^{\circ}$ . UO 5 is an analogous case to LO 5, where its  $\theta_{23}$  prediction span is small and occurs at the very upper limit of the  $3\sigma$  boundary,  $51.2^{\circ} \leq \theta_{23} \leq 53.3^{\circ}$ .

Maximal CP violation is possible in UOs 1–3 and UO 5, however the  $\delta-\theta_{23}$  correlation structure differs between cases. UOs 1–3 share the same pattern where the maximal  $\delta$  value (90°) occurs for large  $\theta_{23}$  values and CP-conserving values of  $\delta$  are associated with 45.6°  $\leq \theta_{23} \leq$  48.4°. The  $\delta$  correlation of UO 5 differs significantly from UOs 1–3 as CP-conserving values of  $\delta$  are not predicted and maximal  $\delta$  favored. In the case of UO 4, the correlation structures are particularly distinctive and unlike the previously discussed cases, the maximal  $\delta$  value (55°) is much smaller. A unique aspect of LO 4 is that there are two distinct regions of  $\theta_{23}$ 

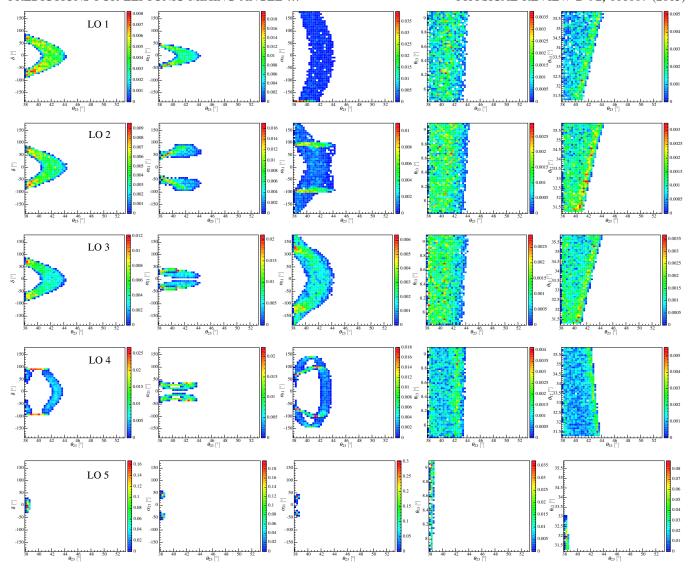


FIG. 2 (color online). Two-dimensional histograms showing the phases and mixing angles as a function of  $\theta_{23}$  for predictions that agree to a  $3\sigma$  level with global fit data [48]. The color chart shows relative frequencies of solutions where red (dark blue) represents a higher (lower) frequency. Each prediction is labeled lower octant (LO) 1–5.

where *CP*-conserving values of  $\delta$  can occur:  $47.4^{\circ} \le \theta_{23} \le 49.2^{\circ}$  and  $51^{\circ} \le \theta_{23} \le 52^{\circ}$ .

In regards to the  $\theta_{12}-\theta_{23}$  correlation of UOs 1–3, all regions of the  $3\sigma$  range of  $\theta_{12}$  are allowed for  $49^{\circ} \leq \theta_{23} \leq 53.3^{\circ}$ . Larger values of  $\theta_{12}$  are favored for a near maximal  $\theta_{23}$ . It is worth noting this dependence (large  $\theta_{12}$  associated with near maximal values of  $\theta_{23}$ ) is similar to the lower octant predictions LOs 1–3. For  $49^{\circ} \leq \theta_{23} \leq 52^{\circ}$ , the  $\theta_{12}$  predictions of UO 4 are indistinguishable from UOs 1–3. In spite of this, for certain  $\theta_{23}$ 's, these cases can be differentiated. For example,  $\theta_{23} > 52.6^{\circ}$  UO 4 predicts large valued  $\theta_{12}$  (~36°), whereas the  $\theta_{12}$  of UOs 1–3 can attain any value in the  $3\sigma$  range. Moreover, at near maximal values of  $\theta_{23}$  (46°), UO 4 predicts smaller  $\theta_{12}$  values (31.3°) than UOs 1–3. Discrimination between UOs 1–3 and UO 5 is not possible using  $\theta_{12}-\theta_{23}$  correlations (as there is

complete overlap in the predictions) and therefore a combination of  $\theta_{23}$  and  $\delta$  measurements in conjunction with  $\nu 0\beta\beta$  decay study would be required to disentangle these predictions.

In summary,  $\delta$  and  $\alpha_{21}$  are bounded between  $\pm 90^\circ$ . Moreover, the ability to discriminate between predictions is often dependent upon the value of  $\theta_{23}$  and in certain cases predictions are only differentiable with knowledge of the Majorana phases.

### D. Predictions spanning both octants

We have chosen five representative cases that span both the upper and lower octants of  $\theta_{23}$ . The predicted regions of  $\theta_{23}$  vary amongst these cases: both octants (BO) 1 has the greatest viable range, which fully covers the  $3\sigma$  region of

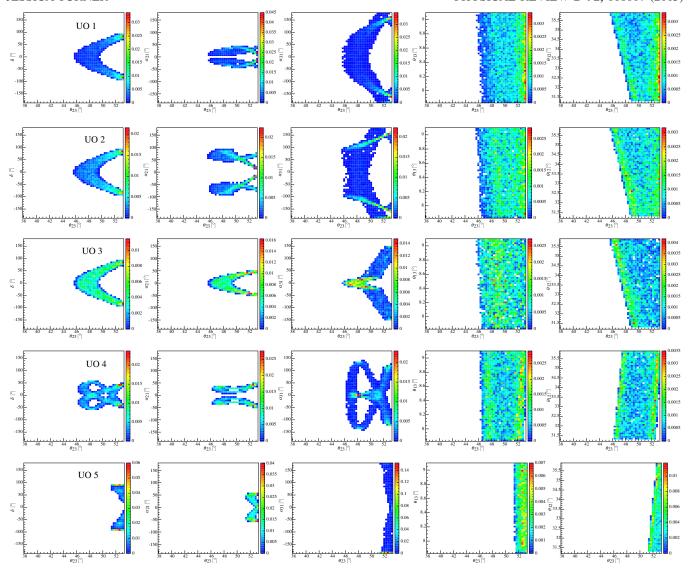


FIG. 3 (color online). Two-dimensional histograms showing the phases and mixing angles as a function of  $\theta_{23}$  for predictions that agree to a  $3\sigma$  level with global fit data [48]. The color chart shows relative frequencies of solutions where red (dark blue) represents a higher (lower) frequency. Each prediction is labeled upper octant (UO) 1–5.

 $\theta_{23}$ . BO 2 and BO 5 also have a wide range of  $\theta_{23}$ :  $38.2^{\circ} \le \theta_{23} \le 49^{\circ}$  and  $38.2^{\circ} \le \theta_{23} \le 51^{\circ}$  respectively. BO 3 and BO 4 have the smallest viable range of  $\theta_{23}$  with  $44.3^{\circ} \le \theta_{23} \le 53.3^{\circ}$  and  $38.2^{\circ} \le \theta_{23} \le 45.9^{\circ}$  respectively.

There is little structure in the  $\delta-\theta_{23}$  correlation of BO 1:  $\delta$  can attain any value in the range  $\pm 90^\circ$  and there is no dependence on  $\theta_{23}$ . BO 2 has a similar correlation structure to the lower octant predictions: the maximal  $\delta$  value (73°) is correlated to smaller  $\theta_{23}$  values and CP-conserving  $\delta$  spans  $42^\circ \leq \theta_{23} \leq 48.2^\circ$ . BO 3 and BO 4 have comparable  $\delta-\theta_{23}$  dependence; the maximal  $\delta$ , 69° and 61° respectively, occurs at the extreme upper and lower  $3\sigma$  limit of  $\theta_{23}$ . In comparison with BO 1–4, BO 5 has a highly constrained  $\delta$  with a maximal value of  $14^\circ$  for  $49^\circ \leq \theta_{23} \leq 50^\circ$ . In the scenario where  $\delta$  is maximally CP violating, the only viable prediction of this set is BO 1. Interestingly, in spite of BO 1

lacking predictivity in regards to parameters  $\delta$ ,  $\alpha_{31}$  and  $\theta_{13}$ , its  $\alpha_{21}$  and  $\theta_{12}$  predictions attain very specific values ( $\alpha_{21}=0^\circ$  and  $34.8^\circ \le \theta_{12} \le 35.2^\circ$ ). BO 1 would be of particular interest in  $\nu 0\beta\beta$  decay studies as it has a single  $\alpha_{21}$  value and exceptionally narrow  $\theta_{12}$  range.

The  $\theta_{12} - \theta_{23}$  dependence of BO 3 and BO 4 are similar: for near maximal  $\theta_{23}$ , there is a very limited range of  $\theta_{12}$  values (~36°) and for  $\theta_{23}$  close to the upper or lower  $3\sigma$  boundary, the possible  $\theta_{12}$ 's become less constrained. This appears to be a common theme of many of the predictions: near maximal  $\theta_{23}$ 's have very specific  $\theta_{12}$  predictions. In the case of BO 2,  $\theta_{12}$  can attain any value in the  $3\sigma$  range for  $\theta_{23} \le 43.5^\circ$  and for larger  $\theta_{23}$ , smaller values of  $\theta_{12}$  are preferable. There is significant overlap in  $\theta_{12}$  predictions for BO 2 and BO 5 and only in the scenario  $\theta_{23} \ge 47^\circ$  do their predictions differ. A special feature of BO 5, akin to

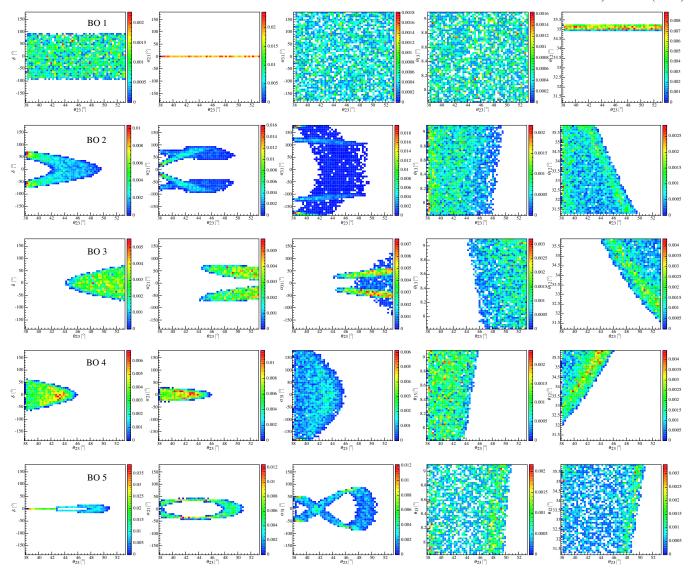


FIG. 4 (color online). Two-dimensional histograms showing the phases and mixing angles as a function of  $\theta_{23}$  for predictions that agree to a  $3\sigma$  level with global fit data [48]. The color chart shows relative frequencies of solutions where red (dark blue) represents a higher (lower) frequency. Each prediction is labeled both octants (BO) 1–5.

BO 1, is that certain observables are more constrained than others; for instance predictions of  $\theta_{13}$  and  $\theta_{12}$  range widely, whereas the leptonic phases are more highly constrained ( $|\delta| \le 14^{\circ}$ ,  $|\alpha_{21}| \le 40^{\circ}$ ,  $|\alpha_{31}| \le 90^{\circ}$ ).

In summary,  $\delta$  and  $\alpha_{21}$  can only attain values  $\pm 90^\circ$ . Furthermore, there are two examples (BO 1 and BO 5), in which certain observables are highly unconstrained but in balance other parameters can only attain very specific values. Therefore in spite of a lack of predictivity in certain observables, these cases still remain testable by upcoming long- and medium baseline experiments.

### V. CONCLUSIONS

In this article, we studied the correlations of leptonic observables that result from a flavor symmetry,  $A_5$ ,

combined with gCP breaking into residual symmetries  $G_{\nu}=\mathbb{Z}_2\times CP$  and  $G_e=\mathbb{Z}_2$ . This combination of residual symmetries introduces three continuous parameters and, unsurprisingly, we obtain a wider range of predictions than in studies that use only one input parameter. The flavor symmetry studies that implement one input parameter and are of low order such as  $A_4$  [28,34],  $S_4$  [28,35] and  $A_5$  [39–41] share common predictions such as  $|\sin\delta|=1$ ,  $|\sin\alpha_{21}|=|\sin\alpha_{31}|=0$  and maximally CP-violating  $\delta$  associated with maximal  $\theta_{23}$ . We find that the addition of two continuous parameters allows for more possibilities in correlations and predictions of nontrivial leptonic phases differing from 0,  $\frac{\pi}{2}$ ,  $\pi$  and  $\frac{3\pi}{2}$ . Using a number of example cases we have shown that certain predictions are indistinguishable using

oscillation parameters  $\delta$ ,  $\theta_{12}$  and  $\theta_{23}$  alone and therefore input from  $\nu 0\beta\beta$  decay experiments is necessary. We find that, in general, the ability to discriminate between predictions is improved for a near maximal  $\theta_{23}$  and that even in specific cases in which there is no predictivity for one parameter (e.g. BO 1 and BO 5), other leptonic observables may be highly constrained and may provide testable predictions. In spite of a greater number of predictions, all of our cases share the feature of  $\delta$  and  $\alpha_{21}$  phases being bounded by  $\pm 90^\circ$ , the former of which is testable by long-baseline oscillation experiments.

In conclusion, we find that relaxing the possible combinations of low-energy residual symmetries permits a wider range of predictions with more complex correlations between leptonic observables which have the potential to be tested at upcoming neutrino oscillation and  $\nu 0\beta\beta$  experiments.

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### **APPENDIX**

In this appendix we will provide the form of the matrices that we used to derive predictions in the lower, upper and both octant results. The notation that will be used to denote real rotations is

$$R_{12} = \begin{pmatrix} c_{\theta} & s_{\theta} & 0 \\ -s_{\theta} & c_{\theta} & 0 \\ 0 & 0 & 1 \end{pmatrix}, \qquad R_{13} = \begin{pmatrix} c_{\theta} & 0 & s_{\theta} \\ 0 & 1 & 0 \\ -s_{\theta} & 0 & c_{\theta} \end{pmatrix} \quad \text{and}$$

$$R_{23} = \begin{pmatrix} 1 & 0 & 0 \\ 0 & c_{\theta} & s_{\theta} \\ 0 & -s_{\theta} & c_{\theta} \end{pmatrix}. \tag{A1}$$

In regards to the complex rotations they will be written as

$$R_{12}C = \begin{pmatrix} c_{\omega} & s_{\omega}e^{i\gamma} & 0\\ -s_{\omega}e^{-i\gamma} & c_{\omega} & 0\\ 0 & 0 & 1 \end{pmatrix},$$

$$R_{13}C = \begin{pmatrix} c_{\omega} & 0 & s_{\omega}e^{i\gamma}\\ 0 & 1 & 0\\ -s_{\omega}e^{-i\gamma} & 0 & c_{\omega} \end{pmatrix} \text{ and }$$

$$R_{23}C = \begin{pmatrix} 1 & 0 & 0\\ 0 & c_{\omega} & s_{\omega}e^{i\gamma}\\ 0 & -s_{\omega}e^{-i\gamma} & c_{\omega} \end{pmatrix}.$$
(A2)

The  $\mathbb{Z}_2$  elements that give distinct results are

$$\mathbb{Z}_{2}1 = \frac{1}{2} \begin{pmatrix} \phi & -1 & \phi \\ -1 & -\phi & -\frac{1}{\phi} \\ \phi & -\frac{1}{\phi} & -1 \end{pmatrix}, 
\varphi = \frac{1}{\phi} \begin{pmatrix} -\phi & -\frac{1}{\phi} & -1 \\ -\frac{1}{\phi} & -1 & \phi \\ -1 & \phi & \frac{1}{\phi} \end{pmatrix}, 
\mathbb{Z}_{2}2 = \frac{1}{2} \begin{pmatrix} -1 & -\phi & -\frac{1}{\phi} \\ -\phi & \frac{1}{\phi} & 1 \\ -\frac{1}{\phi} & 1 & -\phi \end{pmatrix}, 
\mathbb{Z}_{2}3 = \frac{1}{2} \begin{pmatrix} \frac{1}{\phi} & -1 & -\phi \\ -1 & -\phi & \frac{1}{\phi} \\ -\phi & \frac{1}{\phi} & -1 \end{pmatrix}, 
\mathbb{Z}_{2}4 = \frac{1}{2} \begin{pmatrix} -\phi & -\frac{1}{\phi} & 1 \\ -\frac{1}{\phi} & -1 & -\phi \\ 1 & -\phi & \frac{1}{\phi} \end{pmatrix}, 
\mathbb{Z}_{2}5 = \frac{1}{2} \begin{pmatrix} -1 & \phi & \frac{1}{\phi} \\ \phi & \frac{1}{\phi} & -1 \\ \frac{1}{\phi} & 1 & -\phi \end{pmatrix}, 
\mathbb{Z}_{2}6 = \frac{1}{2} \begin{pmatrix} -1 & \phi & \frac{-1}{\phi} \\ \phi & \frac{1}{\phi} & -1 \\ \frac{1}{\phi} & -1 & -\phi \end{pmatrix}. \tag{A3}$$

The diagonalizing matrix of  $\mathbb{Z}_2$ i will be denoted by  $U_i$  for  $i \in 1...6$ . The permutations that have been applied to account for the arbitrariness of ordering of the eigenvectors will be denoted by  $p_1...p_6$ ,

PREDICTIONS FOR LEPTONIC MIXING ANGLE ...

# $p_{1} = \begin{pmatrix} 1 & 0 & 0 \\ 0 & 1 & 0 \\ 0 & 0 & 1 \end{pmatrix}, \qquad p_{2} = \begin{pmatrix} 1 & 0 & 0 \\ 0 & 0 & 1 \\ 0 & 1 & 0 \end{pmatrix},$ $p_{3} = \begin{pmatrix} 0 & 1 & 0 \\ 1 & 0 & 0 \\ 0 & 0 & 1 \end{pmatrix}, \qquad p_{4} = \begin{pmatrix} 0 & 1 & 0 \\ 0 & 0 & 1 \\ 1 & 0 & 0 \end{pmatrix},$ $p_{5} = \begin{pmatrix} 0 & 0 & 1 \\ 1 & 0 & 0 \\ 0 & 1 & 0 \end{pmatrix}, \qquad p_{6} = \begin{pmatrix} 0 & 0 & 1 \\ 0 & 1 & 0 \\ 1 & 0 & 0 \end{pmatrix}. \quad (A4)$

### 2. Upper octant matrices

Result	$U_{ m PMNS}$
UO 1	$p_1 R_{13} C U_2^{\dagger} \Omega_{12} R_{13} p_1$
UO 2	$p_1 R_{13} C U_4^{\dagger} \Omega_{12} R_{13} p_3$
UO 3	$p_4 R_{12} C U_3^{\dagger} \Omega_{12} R_{23} p_4$
UO 4	$p_3 R_{23} C U_5^{\dagger} \Omega_{12} R_{13} p_3$
UO 5	$p_3 R_{23} C U_1^{\dagger} \Omega_{12} R_{12} p_2$

### 1. Lower octant matrices

Result	$U_{ m PMNS}$
LO 1	$p_4 R_{23} C U_1^{\dagger} \Omega_{12} R_{12} p_2$
LO 2	$p_2 R_{13} C U_2^{\dagger} \Omega_{12} R_{13} p_3$
LO 3	$p_1 R_{12} C U_3^{\dagger} \Omega_{12} R_{23} p_4$
LO 4	$p_2 R_{13} C U_2^{\dagger} \Omega_{12} R_{13} p_1$
LO 5	$p_4 R_{23} C U_1^{\dagger} \Omega_{12} R_{13} p_3$

### 3. Both octant matrices

Result	$U_{ m PMNS}$
BO 1	$p_1 R_{23} C U_5^{\dagger} \Omega_{12} R_{12} p_4$
BO 2	$p_1 R_{12} C U_6^{\dagger} \Omega_{13} R_{12} p_4$
BO 3	$p_4 R_{23} C U_1^{\dagger} \Omega_{23} R_{12} p_6$
BO 4	$p_4 R_{12} C U_7^{\dagger} \Omega_{23} R_{23} p_1$
BO 5	$p_3 R_{23} C U_5^{\dagger} \Omega_{13} R_{13} p_3$

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