The user manual: Flavour Symmetry Embedded - GLoBES (FaSE-GLoBES)

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I. INTRODUCTION

The discovery of neutrino oscillations points out the fact that neutrinos have mass, and provides evidence beyond the Standard Model (BSM). This phenomenon is successfully described by a theoretical framework with the help of three neutrino mixing angles (θ_{12}, θ_{13}, θ_{23}), two mass-square splittings (Δm^2_{21}, Δm^2_{31}), and one Dirac CP phase (δ) [1–4]. Thanks to the great efforts in the past two decades, we almost have a complete understanding of such a neutrino oscillation framework. More data in the neutrino oscillation experiments is needed to determine the sign of Δm^2_{31}, to measure the value of $\sin\theta_{23}$, to discover the potential CP violation in the leptonic sector and even to constrain the size of δ [4]. For these purposes, the on-going long baseline experiments (LBLs), such as the NuMI Off-axis ν_e Appearance experiment (NOνA) [5] and the Tokai-to-Kamioka experiment (T2K) [6], can answer these questions with the statistical significance $\gtrsim 3\sigma$ in most of the parameter space. Based on the analysis with their data, the normal mass ordering ($\Delta m^2_{31} > 0$), the higher θ_{23} octant ($\theta_{23} > 45^\circ$), and $\delta \sim 270^\circ$ are preferred so far [4]. The future LBLs, Deep Underground Neutrino Experiment (DUNE) [7], Tokai to Hyper-Kamiokande (T2HK) [8], and the medium baseline reactor experiment, the Jiangmen Underground Neutrino Observatory (JUNO) [9, 10] will further complete our knowledge of neutrino oscillations.

Flavour symmetry models are used to explain the origin of the neutrino mixing, and to predict the value of oscillation parameters (some of useful review articles are [11–17]). These models are motivated by some interesting features, *such as* $\theta_{12} \sim 33^\circ$, and $\theta_{23} \sim 45^\circ$. Before the discovery of non-zero θ_{13} measurement by Daya Bay experiment [18], the 'tri-bi-maximal' neutrino mixing (TBM) ansatz, which was proposed in 2002 by Horrison, Perkins, and Scott [19], fitted with the experimental data in a good agreement:

$$U_{\rm TBM} = \begin{pmatrix} 2/\sqrt{6} & 1/\sqrt{3} & 0 \\ -1/\sqrt{6} & 1/\sqrt{3} & 1/\sqrt{2} \\ 1/\sqrt{6} & -1/\sqrt{3} & 1/\sqrt{2} \end{pmatrix}.$$

With the fact that $\theta_{13} \approx 8^\circ$, several ways to obtain such non-zero value of θ_{13} are proposed. One of popular proposals is to correct the tri-bi-maximal neutrino mixing such that

$$\sin\theta_{12} = (1+s)/\sqrt{3},\ \sin\theta_{13} = r/\sqrt{2},\ \text{and}\ \sin\theta_{23} = (1+a)/\sqrt{2}.$$

Th neutrino mixing ansatz can be realised by high-energy symmetries G_f. The symmetry

of discrete groups G_f, preserved at the high energy but slightly broken at the lower energy, predicts the neutrino mixing, mass-square splittings, and the CP violation phase (Dirac and Majorana phases), with reduced degrees of freedom. The symmetries need to be broken at the low energy. Otherwise, the flavour of leptons cannot be distinguished. There are several approaches for the symmetry breaking, including the direct, indirect, and semi approaches. Direct approach preserves the residual symmetries of G_f in the charged-lepton or neutrino sector. On the other hand, there is no residual symmetry preserved in neither charged-lepton nor neutrino sector in the indirect approach. In the semi approach, the charged-lepton and neutrino sectors preserve different residual symmetries, respectively. This symmetry is broken by extending the Higgs sector or introducing the flavons. To achieve the δ prediction, many models are based on a discrete family symmetry G_f together with a non-commuting CP symmetry $H_{\rm CP}$. Broken in different approaches, the symmetry $G_f \otimes H_{\rm CP}$ can predict different patterns for the neutrino mixing. For example, in the semi-direct approach, S_4 is preserved in the leading order, and leads the bimaximal (MB) or tri-bimaximal (TB) neutrino mixing, while the higher order terms bring the correction to the neutrino mixing.

Currently, some works discuss on how the future experiments can be used for testing these flavour symmetry models in the phenomenological point of view, *e.g.* Ref. [20–23]. A large number of these works are based on the `c`-library – **G**eneral **Lo**ng **B**aseline **E**xperiment **S**imulator (**GLoBES**) [24, 25], which is a convenient simulation tool to simulate neutrino oscillation experiments via the Abstract Experiment Definition Language (AEDL). Some AEDL files for experiments are also available on **GLoBES** website, while the working group of DUNE experiment also releases their AEDL files [26]. **GLoBES** can be taken as one of popular and useful tools in the community of neutrino oscillation physics. However, it has not yet to be extended for the purpose of analysing flavour symmetry models. In this work, we will present our simulation tool **Fla**vour **S**ymmetry **E**mbedded - **GLoBES** (**FaSE-GloBES**) that is a simulation code based on GLoBES library for studying the flavour symmetry. **FaSE** is a supplemental tool for GLoBES, written in `c/c++` language, and FASE-GLoBES allows the user to assign the flavour symmetry model and analysis how a flavour symmetry model can be excluded by the simulated neutrino oscillation experiment.

II. OVERVIEW FASE-GLOBES

Flavor **S**ymmetry **E**mbedded (**FaSE**) is a supplemental tool for **G**eneral **Lo**ng **B**aseline **E**xperiment **S**imulator (**GLoBES**) in order to analysis how a leptonic flavour symmetry model can be tested in neutrino oscillation experiments. **FaSE** is written in the `c/c++` language, and consist with three codes **FASE_GLoBES.c** and **model-input.c**. The user defines the model in to **model-input.c**, while **FASE_GLoBES.c** does not need to be touched.

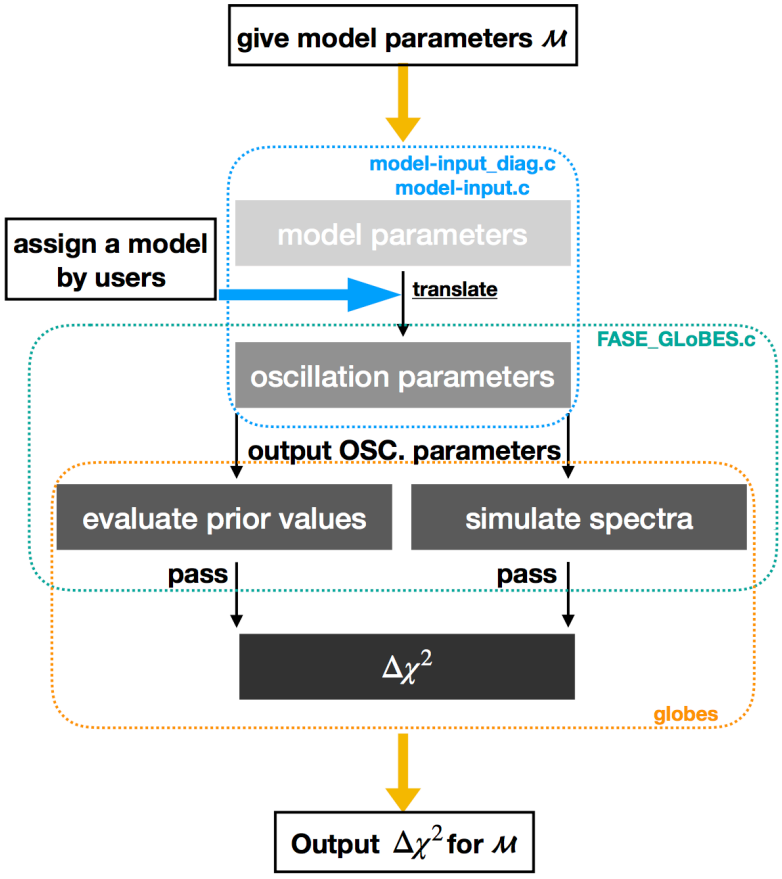

Figure 1: A scheme to correlate the model parameters with standard neutrino oscillation parameters. The error propagation is implemented in the simulation code up to the spectra analysis.

The concept of **FaSE-GLoBES** is shown in Fig. 1, in which three parts are shown: 1. **the parameter translation** (the blue box), 2. **giving oscillation-parameter values** (the green box), and 3. **the χ^2-value calculation** (the orange box). The idea behind this flow chart Fig. 1 is that given a set of value for model parameter, the corresponding values

for oscillation parameters are obtained by a translation, which is assigned by the user in **model-input.c**. And then, through **FASE_GLoBES.c** these oscillation-parameter values are passed in to **GLoBES** library to simulate the event spectrum for evaluating the χ^2 value.

API functions in **FaSE** are listed:

1. `MODEL_init(`N_{para}`)`,
2. `FASE_glb_probability_matrix`,
3. `FASE_glb_set_oscillation_parameters`,
4. `FASE_glb_get_oscillation_parameters`,
5. `FASE_prior_OSC`,
6. `FASE_prior_model`.

The first one is to initialise **FaSE** with the number of input parameters N_{para}. The next three functions need to be included to replace the default **GLoBES** probability engine by the one that can read the output from **model-input.c**, as follows.

```
glbRegisterProbabilityEngine(6,
&FASE_glb_probability_matrix,
&FASE_glb_set_oscillation_parameters,
&FASE_glb_get_oscillation_parameters,
NULL);
```

This probability engine can work with oscillation or model parameters. It can be set by the user with the parameter `PARA`. If `PARA=STAN` (`PARA=MODEL`) the probability engine works with oscillation (model) parameters. The final two items on the API list are prior functions. Once the user gives the prior in oscillation (model) parameters, the user needs to call `FASE_prior_OSC` (`FASE_prior_model`) as follows.

```
glbRegisterPriorFunction(FASE_prior_OSC,NULL,NULL,NULL);
```

or

```
glbRegisterPriorFunction(FASE_prior_model,NULL,NULL,NULL);
```

We note that except for setting the probability engine and the prior function, the other parts in the main code should follow with the GLoBES manual.

In the following, we give more details about using **FASE-GLoBES**. In Sec. II A, we will give the instruction to download and install **GLoBES** and **FASE**. Then, in Sec. II B, we will introduce how to compile and run the code. More complicated is setting the model, and the way to do this will be described in Sec. II C. Finally, we will discuss how to set up the prior in Sec. II D.

A. Download and Installation

download and install GLoBES and FASE-GLoBES

[will fill in after the code ready on Github]

B. Compile and Run

As **FaSE** works based on **GLoBES**, the user should use the **GLoBES** `Makefile`, but include the binary file of **FaSE**. To do so, we include the script in the `Makefile`,

```
{makefile_execution}:  my_program.o FASE_GLoBES.o model-input_diag.o
gcc my_program.o FASE_GLoBES.o model-input_diag.o -o my_executable
$(LDFLAGS) $(local_LDFLAGS)
```

or

```
{makefile_execution}:  my_program.o FASE_GLoBES.o model-input.o
gcc my_program.o FASE_GLoBES.o model-input.o -o my_executable $(LDFLAGS)
$(local_LDFLAGS)
```

where `{compile_execusion}` is the makefile commend for the program `my_program.o` and the execution `my_executable` is the output. After giving this script in the **GLoBES** `Makefile`, the user needs to makefile the program `my_program` on the terminal.

```
makefile {makefile_execution}
```

And, the user can execute the program by typing the following commend on the terminal.

```
./my_executable
```

The user can also compile the program without the **GLoBES** `Makefile` with the following script.

```
gcc my_program.o FASE_GLoBES.o model-input_diag.o -lglobes -LGLB_DIR/lib/
-o my_executable
```

or

```
gcc my_program.o FASE_GLoBES.o model-input.o -lglobes -LGLB_DIR/lib/ -o
my_executable
```

And, the execution is also in the same way: `./my_executable`.

Finally we need to initialize **FaSE** by giving N_{para} which is the number of model parameters. To do so, we need to include the following script in the main code.
`MODEL_init(`N_{para}`);`

C. Model setting

The function `MtoS` can do the translator from model parameters $\vec{\theta}_{Model}$ to oscillation parameters $\vec{\theta}_{OSC.}$. After the user gives the array $\vec{\theta}_{Model}$ to the function `MtoS`, the output is the corresponding oscillation parameter $\vec{\theta}_{OSC.}$, of which components are θ_{12}, θ_{13}, θ_{23}, δ, Δm^2_{21}, and Δm^2_{31}. For the first four components, values are given in the unit of **rad**, while the other two are in $\mathbf{eV}^2$. These values will be passed in to **FaSE_GLoBES** to simulate the experimental spectra and compute the prior value.

To do the translation from $\vec{\theta}_{Model}$ to $\vec{\theta}_{OSC.}$, the user can assign the relation between the oscillation and model parameter sets, or define the mass matrix in model parameters, which will be diagonalised by the function `ModelTO` to obtain the corresponding oscillation-parameter values. In the way of directly giving the relation between oscillation and model parameter sets, the user needs to provide

$$\vec{\theta}_{Model} = \vec{f}(\vec{\theta}_{OSC.}) \tag{1}$$

in the function `MtoS`.

The oscillation parameters can also be obtained in the way based on

$$U^\dagger \mathcal{M}\mathcal{M}^\dagger U = \mathbf{M}^2, \text{ where } \mathbf{M}^2_{\alpha\beta} = m^2_\alpha \delta_{\alpha\beta}, \tag{2}$$

where $\mathcal{M}$ ($\mathbf{M}$) is the neutrino mass matrix in the flavour (mass) state. The matrix $\mathcal{M}$ is

given by user with model parameters $\vec{\theta}_{Model}$. The matrix U is the neutrino mixing matrix, and can be used for getting mixing angles. The difference between any two diagonal elements of $\mathbf{M}$ ($\mathbf{M}_{ii} - \mathbf{M}_{jj}$) is the mass-squared difference (Δm^2_{ij}). This diagnolisation Eq. 2 can be done with the function `ModelTO`, which needs to be called in `MtoS` and outputs directly the vector $\vec{\theta}_{OSC.}$.

D. Prior setting

Given a set of values for model parameters, **FASE_GLoBES.c** will obtain the corresponding oscillation-parameter values from **model-input.c**, and will pass these values to simulate event spectra and to compute the prior value. Two gaussian prior functions are provided in **FaSE** – `FASE_prior_OSC` and `FASE_prior_model`. These two functions are for different purposes. If the user give the prior in oscillation (model) parameters, the user should register `FASE_prior_OSC` (`FASE_prior_model`) with the **GLoBES** function `glbRegisterPriorFunction`, as we introduced in the beginning of this section. The user also needs to assign the parameter `PARA=STAN` (`PARA=Model`), when the user prefers to give the prior in oscillation (model) parameters. The Gaussian prior is

$$\chi^2_{prior} = \sum_i \frac{(\theta_i - \theta_i^c)^2}{\sigma_i^2}, \tag{3}$$

where θ_i is one of parameters constrained by prior, θ_i^c (σ_i) is the central value (Gaussian width) of the prior for θ_i. We note that θ_i can be either model ($\vec{\theta}_{Model}$) or oscillation parameters ($\vec{\theta}_{OSC.}$). The values of θ_i^c and σ_i need to be given by the user through three arrays: `Central_prior`, `UPPER_prior`, and `LOWER_prior`, in which there are six components. To treat asymmetry of width for upper ($\theta_i > \theta_i^c$) and lower ($\theta_i < \theta_i^c$) Gaussian widths, we give values in two arrays `UPPER_prior`, and `LOWER_prior`, respectively. If the user gives the prior in model parameters, the order of each component follows with the setup of input of the probability engine. While the user gives the prior in oscillation parameters, the six components in order are θ_{12}, θ_{13}, θ_{23}, δ, Δm^2_{21}, and Δm^2_{31}. The first four parameter are in **rad**, and the final two are in $\mathbf{eV}^2$.

Finally, some restrictions are imposed by the studied flavour symmetry model. We set up these restrictions in the function `model_restriction` in `model-input.c`. In the function `model_restriction`, the user needs to *return* 1 once the restriction is broken. For

example, if the normal ordering is imposed, we give " `if (DMS31<0) { return 1;}` " in `model_restriction`, where `DMS31` is the variable for Δm^2_{31}. Then, when the restriction is broken, `model_restriction` returns the value 1 to the prior function `FASE_prior_OSC` or `FASE_prior_model`.

III. EXAMPLES

A. Setup for a model

We now take the littlest seesaw model in the tri-direct approach as an example. In this model, the atmospheric and solar flavon vacuum alignments are $\langle\phi_{\rm atm}\rangle \propto (1,\omega^2,\omega)^T$ and $\langle\phi_{\rm sol}\rangle \propto (1,x,x)^T$, where stands for a cube root of unity and the parameter x is real because of the imposed CP symmetry. Under this model, the light left-handed Majorana neutrino mass matrix is given by

$$m_\nu = m_a \begin{pmatrix} 1 & \omega & \omega^2 \\ \omega & \omega^2 & 1 \\ \omega^2 & 1 & \omega \end{pmatrix} + e^{i\eta} m_s \begin{pmatrix} 1 & x & x \\ x & x^2 & x^2 \\ x & x^2 & x^2 \end{pmatrix} , \tag{4}$$

where x, η, m_a, and the ratio $r \equiv m_s/m_a$ are four parameters and will be constrained by experimental data. We note that from Eq. (4), $m_1 = 0$ and the normal mass ordering are imposed, and will need to be imposed in **FaSE-GLoBES**. Therefore, the restrictions in this model are $m_a > 0$ and $r > 0$.

```
  int MtoS(double OSC_PARAMS[6], double M_para[])
{
double x=M_para[0];
double eta=M_para[1];
double r=M_para[2];
double ma=M_para[3];
double ms=ma*r;

  double complex Mass_Matrix[] = {ma+ms*(cos(eta) + I*sin(eta)),
ma*(cos(6.6666e-1*M_PI) + I*sin(6.6666e-1*M_PI))+x*ms*(cos(eta) +
I*sin(eta)), ma*(cos(6.6666e-1*M_PI) + I*sin(6.6666e-1*M_PI))*(cos(6.6666e-1*M_PI)
```

Table I: A summary of the relation between oscillation parameters and TDLS model parameters [27]. Two requirements are imposed by TDLS: the smallest mass state $m_1 = 0$ and the normal mass ordering. The sign of $\sin\delta$ depends on the sign of $x\cos\psi$: "+" ("−") is for $x\cos\psi > 0$ (< 0).

model parameters	x, η, r, m_a
combinations of model parameters	$y = \frac{5x^2+2x+2}{2(x^2+x+1)}(m_a + e^{i\eta}m_s)$, $z = -\frac{\sqrt{5x^2+2x+2}}{2(x^2+x+1)}\left[(x+2)m_a - x(2x+1)e^{i\eta}m_s\right]$, $w = \frac{1}{2(x^2+x+1)}\left[(x+2)^2 m_a + x^2\,(2x+1)^2\,e^{i\eta}m_s\right]$, $\sin\psi = \frac{\Im(y^*z+wz^*)}{\lvert y^*z+wz^*\rvert}$, $\quad \cos\psi = \frac{\Re(y^*z+wz^*)}{\lvert y^*z+wz^*\rvert}$. $\sin 2\theta = \frac{2\lvert y^*z+wz^*\rvert}{\sqrt{(\lvert w\rvert^2-\lvert y\rvert^2)^2+4\lvert y^*z+wz^*\rvert^2}}$, $\cos 2\theta = \frac{\lvert w\rvert^2-\lvert y\rvert^2}{\sqrt{(\lvert w\rvert^2-\lvert y\rvert^2)^2+4\lvert y^*z+wz^*\rvert^2}}$.
oscillation parameters	$\Delta m^2_{21} = m^2_2 = \frac{1}{2}\left[\lvert y\rvert^2 + \lvert w\rvert^2 + 2\lvert z\rvert^2 - \frac{\lvert w\rvert^2-\lvert y\rvert^2}{\cos\theta}\right]$, $\Delta m^2_{31} = m^2_3 = \frac{1}{2}\left[\lvert y\rvert^2 + \lvert w\rvert^2 + 2\lvert z\rvert^2 + \frac{\lvert w\rvert^2-\lvert y\rvert^2}{\cos\theta}\right]$, $\sin^2\theta_{12} = 1 - \frac{3x^2}{3x^2+2(x^2+x+1)\cos^2\theta}$, $\sin^2\theta_{13} = \frac{2\left(x^2+x+1\right)\sin^2\theta}{5x^2+2x+2}$, $\sin^2\theta_{23} = \frac{1}{2} + \frac{x\sqrt{3(5x^2+2x+2)}\sin 2\theta\sin\psi}{2[3x^2+2(x^2+x+1)\cos^2\theta]}$, $\cos\delta = \frac{\cot 2\theta_{23}\left[3x^2-\left(4x^2+x+1\right)\cos^2\theta_{13}\right]}{\sqrt{3}\lvert x\rvert\sin\theta_{13}\sqrt{(5x^2+2x+2)\cos^2\theta_{13}-3x^2}}$, $\sin\delta = \pm\csc 2\theta_{23}\sqrt{1 + \frac{(x^2+x+1)^2\cot^2\theta_{13}\cos^2 2\theta_{23}}{3x^2[3x^2\tan^2\theta_{13}-2(x^2+x+1)]}}$.

```
+ I*sin(6.6666e-1*M_PI))+x*ms*(cos(eta) + I*sin(eta)),
ma*(cos(6.6666e-1*M_PI) + I*sin(6.6666e-1*M_PI))+x*ms*(cos(eta) +
I*sin(eta)), ma*(cos(6.6666e-1*M_PI) + I*sin(6.6666e-1*M_PI))*(cos(6.6666e-1*M_PI)
+ I*sin(6.6666e-1*M_PI))+x*x*ms*(cos(eta) + I*sin(eta)),
ma+x*x*ms*(cos(eta) + I*sin(eta)), ma*(cos(6.6666e-1*M_PI) +
I*sin(6.6666e-1*M_PI))*(cos(6.6666e-1*M_PI) + I*sin(6.6666e-1*M_PI))+x*ms*(cos(eta)
+ I*sin(eta)), ma+x*x*ms*(cos(eta) + I*sin(eta)), ma*(cos(6.6666e-1*M_PI) +
I*sin(6.6666e-1*M_PI))+x*x*ms*(cos(eta) + I*sin(eta))};

  STAN_OSC(Mass_Matrix,OSC_PARAMS);

  return 0;
```

```
}
   double MtoS(double osc_para[6], double M_para[])
{
/*example:  tri-direct*/

double x_in = M_para[0];
double eta_in = M_para[1];
double r_in = M_para[2];
double ma_in = M_para[3];

osc_para[0]=TDth12(x_in,eta_in,r_in, ma_in);
osc_para[1]=TDth13(x_in,eta_in,r_in, ma_in);
osc_para[2]=TDth23(x_in,eta_in,r_in, ma_in);
osc_para[3]=TDdCP(x_in,eta_in,r_in, ma_in);
osc_para[4]=TDdm21(x_in,eta_in,r_in, ma_in);
osc_para[5]=TDdm31(x_in,eta_in,r_in, ma_in);

return 0;
}
```

The functions `TDth12`, `TDth13`, `TDth23`, `TDdCP`, `TDdm21`, and `TDdm31` are included for obtaining the value of θ_{12}, θ_{13}, θ_{23}, δ, Δm^2_{21}, and Δm^2_{31} according to Table I, respectively.

The restrictions $m_a > 0$ and $r > 0$ need to be setup in `model_restriction` as follows.

```
double model_restriction(double model [])
{
double x=model[0];
double eta=model[1];
double r=model[2];
double ma=model[3];

if(ma<0) {return 1;}
```

```
if(r<0) {return 1;}


return 0;
}
```

If there is no any restrictions, we simply *return* 0 in `model_restriction` as follows.

```
double model_restriction(double model []){ return 0;}
```

B. Initialise the code

In the beginning of main code, we need to initialise GLoBES and FASE, and include the considered experiment (here we consider MOMENT experiment) as follows.

```
glbInit(argv[0]);
glbInitExperiment("expMOMENT_FIX_FLUX_150KM_addATM_NC.glb",&glb_experiment_list[0],
&glb_num_of_exps);
MODEL_init(4);
```

Then we register the probability engine.

```
glb_init_probability_engine();
glbRegisterProbabilityEngine(6,
&FASE_glb_probability_matrix,
&FASE_glb_set_oscillation_parameters,
&FASE_glb_get_oscillation_parameters,
NULL);
```

In the following we define the model-parameter values, and set the true value in model parameters.

```
float degree = M_PI/180;
float x_true,eta_true,r_true,ma_true;
x_true=-3.65029; eta_true=1.13067*M_PI; r_true=0.511325;
ma_true=3.71199e-3;
glb_params true_values = glbAllocParams();
```

```
glb_params test_values = glbAllocParams();
glb_params input_errors = glbAllocParams();
glb_params centers = glbAllocParams();
glbDefineParams(true_values,x_true,eta_true,r_true,ma_true,0,0);
glbSetDensityParams(true_values,1.0,GLB_ALL);
glbCopyParams(true_values,test_values);
PARA=MODEL;
glbSetOscillationParameters(true_values);
glbSetRates();
```

We then finally set up the projection. The projection `free` is to set all four parameters free to find the minimal value of χ^2, which `project` is used for studying the constraint on x and η.

```
glb_projection projection = glbAllocProjection();
glb_projection free = glbAllocProjection();
glbDefineProjection(free,GLB_FREE,GLB_FREE,GLB_FREE,GLB_FREE,GLB_FIXED,GLB_FIXED);
glbSetDensityProjectionFlag(free, GLB_FREE, GLB_ALL);
glbDefineProjection(projection,GLB_FIXED,GLB_FIXED,GLB_FREE,GLB_FREE,GLB_FIXED,GLB_FIXED
glbSetDensityProjectionFlag(projection, GLB_FREE, GLB_ALL);
```

C. Prior setup

To set up the prior, we need to register the prior function. As mentioned, when we use the prior in oscillation and model parameters, we register `FASE_prior_OSC` and `FASE_prior_model`, respectively. Here we present an example setting up priors in oscillation parameters according to NUFIT4.0 Table II.

Table II: The best fit and 3σ uncertainty, in the results of NuFit4.0 [4].

Parameter	$\theta_{12}/^\circ$	$\theta_{13}/^\circ$	$\theta_{23}/^\circ$	$\delta/^\circ$	$\Delta m^2_{21}/10^{-5}\text{eV}^2$	$\Delta m^2_{31}/10^{-3}\text{eV}^2$
best fit	33.82	8.61	49.6	215	7.39	2.525
3σ Range	$31.61 - 36.27$	$8.22 - 8.99$	$40.3 - 52.4$	$125 - 392$	$6.79 - 8.01$	$2.47 - 2.625$

We therefore register `FASE_prior_OSC` for the prior function as follows.

```
glbRegisterPriorFunction(FASE_prior_OSC,NULL,NULL,NULL);
```

And, then we give the value for the central values, and upper and lower Gaussian widths as follows.

```
UPPER_prior[0]=36.27*degree; LOWER_prior[0]=31.61*degree;
Central_prior[0]=33.82*degree;
UPPER_prior[1]=8.99*degree; LOWER_prior[1]=8.22*degree;
Central_prior[1]=8.61*degree;
UPPER_prior[2]=52.4*degree; LOWER_prior[2]=40.3*degree;
Central_prior[2]=49.6*degree;
UPPER_prior[3]=392*degree; LOWER_prior[3]=125*degree;
Central_prior[3]=215*degree;
UPPER_prior[4]=8.01e-5; LOWER_prior[4]=6.79e-5; Central_prior[4]=7.39e-5;
UPPER_prior[5]=2.625e-3; LOWER_prior[5]=2.427e-3; Central_prior[5]=2.525e-3;


int i;
for (i=0;i<6;i++) {UPPER_prior[i]=fabs(UPPER_prior[i]-Central_prior[i])/3;
LOWER_prior[i]=fabs(LOWER_prior[i]-Central_prior[i])/3;}
for (i=0; i<6; i++) glbSetOscParams(centers,0,i);
glbSetDensityParams(centers,1.0,GLB_ALL); glbCopyParams(centers,input_errors);
glbSetCentralValues(centers); glbSetInputErrors(input_errors);
```

The parameter `degree` is defined as $\pi/180$.

D. Constraint of model parameters

The user can use **FaSE-GLoBES** to study the constraint of model parameters. The expression of χ^2 is used as the default **GLoBES** setting. In more detail, the χ^2 function is constructed based on a log-likelihood ratio,

$$\chi^2(\vec{\theta},\xi_s,\xi_b) = 2\sum_i \left(\eta_i(\vec{\theta},\xi_s,\xi_b) - n_i + n_i \ln \frac{n_i}{\eta_i(\vec{\theta},\xi_s,\xi_b)}\right) + p(\xi_s,\sigma_s) + p(\xi_b,\sigma_b) + \chi^2_{prior}, \quad (5)$$

where i runs over the number of bins, $\eta_i(\vec{\theta}, \xi_s, \xi_b)$ is the hypothesis event rate for bin i and E_i is the central bin energy. The vector $\vec{\theta}$ consists of test model or oscillation parameters. The parameters ξ_s and ξ_b are introduced to account for the systematic uncertainty of normalisation for the signal (subscript $_s$) and background (subscript $_b$) components for the event rate, and are allowed to vary in the fit as nuisance parameters. For a given hypothesised set of parameters $\vec{\theta}$, the event rate for bin i is calculated as

$$\eta_i(\vec{\theta}, \xi_s, \xi_b) = (1 + \xi_s) \times n_i + (1 + \xi_b) \times b_i, \tag{6}$$

where n_i and b_i are the expected number of signal and background events in bin i, respectively. The nuisance parameters are constrained by the Gaussian prior $p(\xi, \sigma) = \xi^2/\sigma^2$ with corresponding uncertainties σ_s and σ_b for the signal and background, respectively. Finally, χ^2_{prior} is a set of Gaussian priors for hypothesis, and is expressed as Eq. 3. After doing all minimisations, the user obtain the χ^2 value for a specific hypothesis $\vec{\theta}^{hyp.}$, $\chi^2(\vec{\theta}^{hyp.})$.

The user of **FaSE-GLoBES** is able to study how model parameters can be constrained by the simulated experiment. To do so, the user needs to define the true event spectrum n_i with a set of model or oscillation parameters, *i.e.* $n_i(\vec{\theta}^{true}_{Model})$ or $n_i(\vec{\theta}^{true}_{OSC.})$. The hypothesis $\vec{\theta}^{hyp.}_{Model}$ predicts the tested event spectrum $\eta_i(\vec{\theta}^{hyp.}_{Model}, \xi_s, \xi_b)$. With the default setting for χ^2 function Eq. 3, from **FaSE-GLoBES** the user computes the statistical quantity,

$$\chi^2(\vec{\theta}^{hyp.}_{Model}), \text{ with } n_i(\vec{\theta}^{true}_{Model}) \text{ or } n_i(\vec{\theta}^{true}_{OSC.}). \tag{7}$$

We note that the minimum of χ^2 in the whole parameter space ($\chi^2_{min.}$) may not be 0. Therefore, to study the precision of model parameters, the user should use the value $\Delta\chi^2(\vec{\theta}^{hyp.}_{Model}) \equiv \chi^2(\vec{\theta}^{hyp.}_{Model}) - \chi^2_{min.}$, instead of χ^2 itself. By varying different hypothesis $\vec{\theta}^{hyp.}_{Model}$, the user can obtain the allowed region of model parameters with the statistical quantity $\Delta\chi^2(\vec{\theta}^{hyp.}_{Model})$.

To get the constraint of x and η, we obtain the minimum of χ^2 value,

```
glbSetProjection(free);
float res0=glbChiNP(true_values,NULL,GLB_ALL);
```

Then, we set two loops to get the $\Delta\chi^2$ value for different x and η

```
float x,eta,r,ma,dx,deta;
```

```
float lower_x,upper_x,lower_eta,upper_eta;
FILE* File=fopen("data/constraint_x_eta.dat", "w");
lower_x=-9; upper_x=-3; lower_eta=0.8*M_PI; upper_eta=2*M_PI;
dx=(upper_x-lower_x)/100; deta=(upper_eta-lower_eta)/100;
glbSetProjection(projection);
for (x=lower_x;x<=upper_x;x=x+dx){
for (eta=lower_eta;eta<=upper_eta;eta=eta+deta){
glbSetOscParams(test_values,x,0); glbSetOscParams(test_values,eta,1);
float res=glbChiNP(test_values,NULL,GLB_ALL);
fprintf(File,"%f %f %f \n",x,eta/M_PI,res-res0);
} fprintf(File,"\n");}
```

We use the same code to obtain the constraint of any other combination of two parameters. Then, we can have the result as in Fig. 2.

E. Model testing

The user can also study how well a flavour symmetry model explains the computed data as predicting how the simulated experiment can exclude this model. In the other word, the user studies the minimum of χ^2 for the flavour symmetry model $\vec{\theta}_{Model}$ as a hypothesis, assuming different true oscillation values, *i.e.* different $\vec{\theta}^{true}_{OSC}$. To do so, we compute the same statistical quantity Eq. 7, while the true spectrum is varied with different true values $\vec{\theta}^{true}_{OSC}$. All model parameters are allowed to be varied with the user-defined prior. Finally, the user needs to adopt Wilk's theorem [28]. When comparing nested models, the $\Delta\chi^2$ test statistics is a random variable asymptotically distributed according to the χ^2-distribution with the number of degrees of freedom, which is equal to the difference in the number of free model parameters `dof`.

We can also study on excluding the model, assuming different true values for oscillation parameters. In this example, we present testing the tri-driect littlest seesaw model in various θ_{23} and δ. To do so, we set the true value in oscillation parameters. We change values of θ_{23} and δ of the true theory, and compute the minimal χ^2 value for the tested model with all four model parameters, free to be varied. And, the studied statistics function is exactly

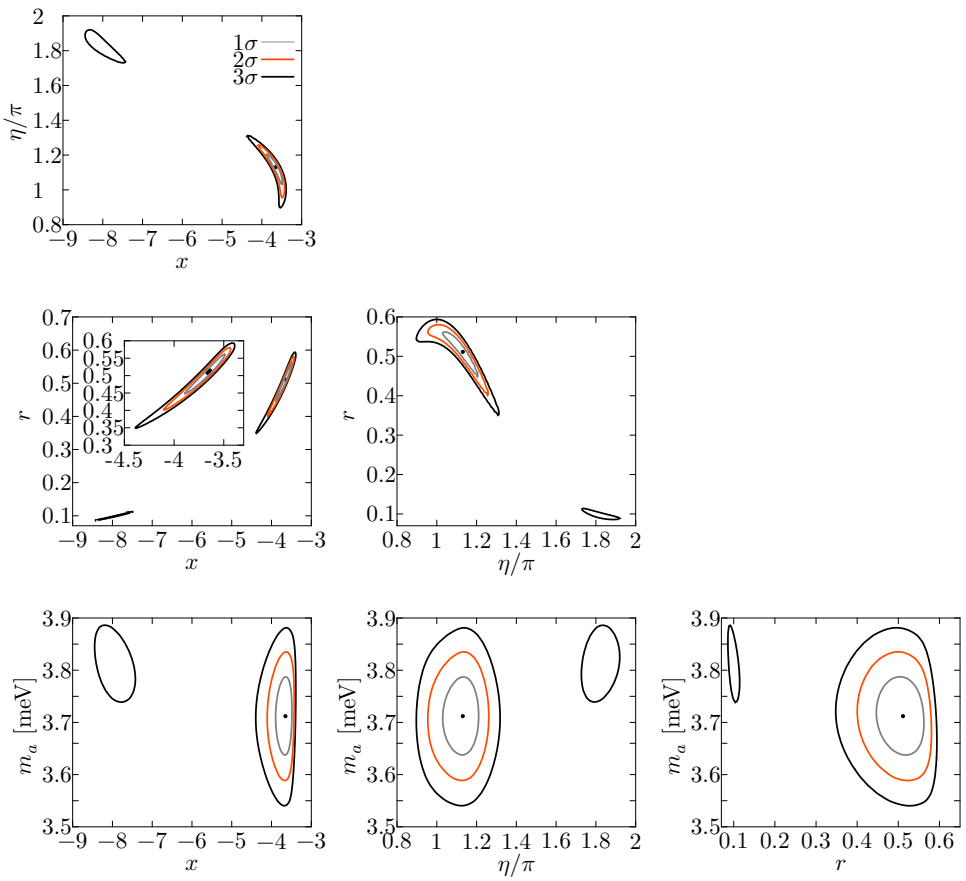

Figure 2: Precision measurements of any two model parameters at 3σ confidence level in the framework of three neutrino oscillations taking uncertainties of the current global fit results, for MOMENT, at 1σ (gray), 2σ (orange), 3σ (black). True values for the model parameters are used $(x,\ \eta,\ r,\ M_a) = (-3.65,\ 1.13\pi,\ 0.511,\ 3.71\ \text{meV})$.

Eq. 5, but the true event rate n_i is predicted by a set of oscillation parameters, which will be varied in the code. And all four model parameter can be varied with the prior given by Eq. 3.

The example code is given below.

```
{ ...  initalize the code ...}
{ ...  register the probability engine ...}

float degree = M_PI/180;
float x_true,eta_true,r_true,ma_true;
x_true=-3.65029; eta_true=1.13067*M_PI; r_true=0.511325;
```

```
ma_true=3.71199e-3;
double M_para[4],OSC_para[6];
M_para[0]=x_true; M_para[1]=eta_true; M_para[2]=r_true; M_para[3]=ma_true;
MtoS(OSC_para, M_para);
double th12_true=OSC_para[0];
double th13_true=OSC_para[1];
double th23_true=OSC_para[2];
double dCP_true=OSC_para[3];
double DM21_true=OSC_para[4];
double DM31_true=OSC_para[5];
glb_params true_values = glbAllocParams();
glb_params test_values = glbAllocParams();
glb_params input_errors = glbAllocParams();
glb_params centers = glbAllocParams();

glbDefineParams(test_values,x_true,eta_true,r_true,ma_true,0,0);
glbSetDensityParams(test_values,1.0,GLB_ALL);
glbDefineParams(true_values,th12_true,th13_true,th23_true,dCP_true,DM21_true,DM31_true);
glbSetDensityParams(true_values,1.0,GLB_ALL);
{ ...  set up for prior ...}
float th23,dCP,dth23,ddCP,lower_th23,upper_th23,lower_dCP,upper_dCP;
FILE* File=fopen("data/ModelTest_th23_dCP_test.dat", "w");
lower_th23=40; upper_th23=50; lower_dCP=125; upper_dCP=392;
dth23=(upper_th23-lower_th23)/100; ddCP=(upper_dCP-lower_dCP)/100;
glbSetProjection(free);
for (th23=lower_th23;th23<=upper_th23;th23=th23+dth23){
for (dCP=lower_dCP;dCP<=upper_dCP;dCP=dCP+ddCP){
glbSetOscParams(true_values,th23*degree,GLB_THETA_23);
glbSetOscParams(true_values,dCP*degree,GLB_DELTA_CP);
PARA=STAN;
glbSetOscillationParameters(true_values);
```

```
glbSetRates();
PARA=MODEL;
float res=glbChiNP(test_values,NULL,GLB_ALL);
fprintf(File,"%f %f %f \n",th23,dCP,gsl_cdf_chisq_Qinv(gsl_cdf_chisq_Q(fabs(res),dof),1)
} fprintf(File,"\n");
}
{ ...  destroy the pointer ...}
```

Finally, in the code we adopt Wilk's theorem [28]. When comparing nested models, the $\Delta\chi^2$ test statistics is a random variable asymptotically distributed according to the χ^2-distribution with the number of degrees of freedom, which is equal to the difference in the number of free model parameters `dof`. Here the number `dof` is 2. And that is our output: `gsl_cdf_chisq_Qinv(gsl_cdf_chisq_Q(fabs(res),dof),1)`.

We use the same code to analysis the exclusion ability for other oscillation-parameter combinations. Then, we have the results shown in Fig. 3.

IV. SUMMARY AND CONCLUSIONS

We have presented **FaSE**, which is a supplemental simulation tool for **GLoBES** to study the flavour symmetry with neutrino oscillation experiments. **FaSE** provides c-codes: **model-input.c** and **FASE_GLoBES.c**. Shown in Fig. 1, **FASE_GLoBES**, which calls functions in **model-input.c**, plays a role as a bridge between **FaSE** and **GLoBES** to simulate the expected spectra and compute the prior value. It can be left to be untouched by users. However, all inputs of the user needs to be given in the code **model-input.c**. Given a set of model parameters $\vec{M}$, with **GLoBES**, the output can be the χ^2 value for the hypothesis $\vec{M}$. `Makefile` is easy to include these two binary files (**model-input** and **FASE_GLoBES**) in the `makefile` script for **GLoBES**.

We also present two examples for **FaSE-GLoBES** with the flavour symmetry model – tri-direct littlest seesaw (TDLS) – and the future neutrino oscillaiton experiment – MOMENT. We show the model can be assign in two ways: the relation between model and oscillation parameters or the form of mass matrix in model parameters. The input of the oscillation probability and prior value can be in model or oscillation parameters. We fur-

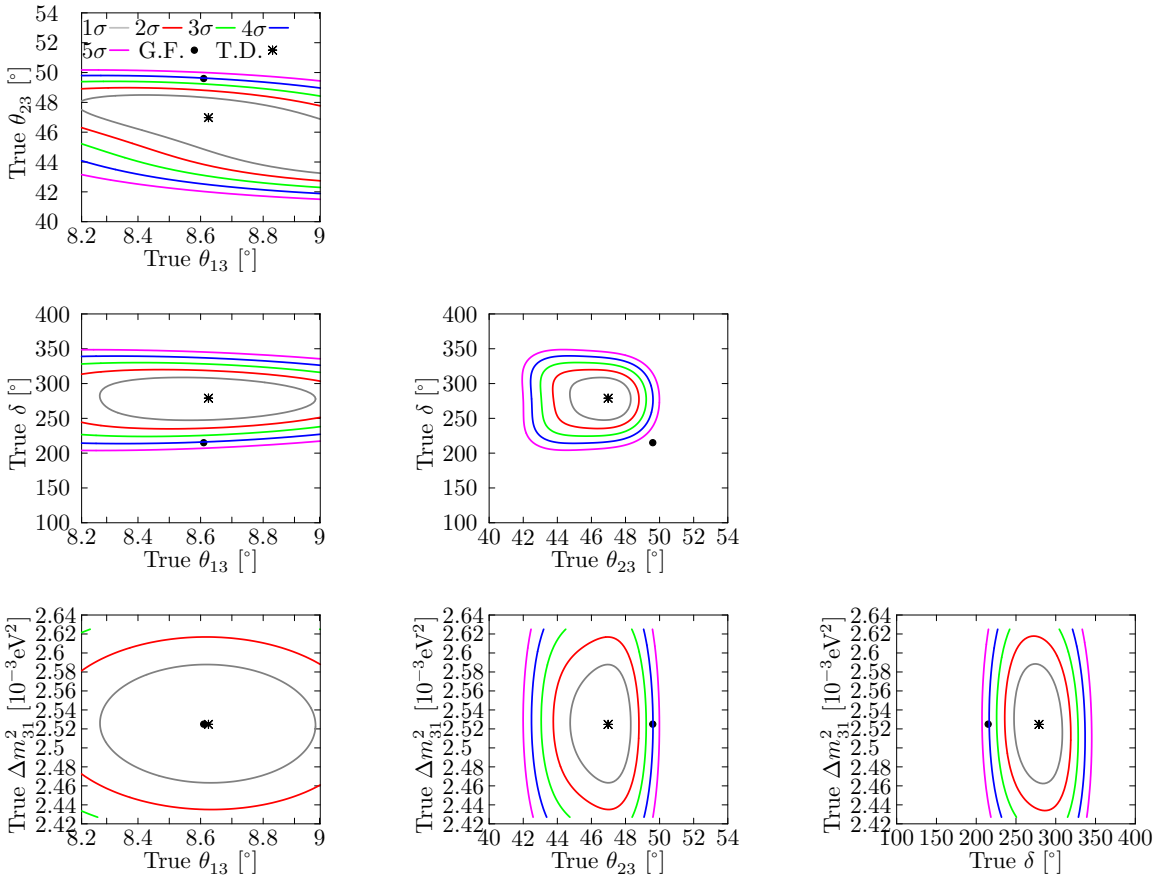

Figure 3: The 2-D exclusion contour for tri-direct littlest seesaw model on the plane of any two true standard parameters, from 1σ to 5σ. The range for each parameter is taken according to the 3σ uncertainty in NuFit4.0 results. The black dot denotes the best fit of NuFit4.0 results $((\theta_{12},\ \theta_{13},\ \theta_{23},\ \delta,\ \Delta m^2_{21},\ \Delta m^2_{31}) = (33.82°,\ 8.61°,\ 49.6°,\ 215°,\ 7.39\times10^{-5}\ \text{eV}^2,\ 2.525\times10^{-3}\ \text{eV}^2))$, while the star is the prediction by the tri-direct littlest seesaw model with NuFit4.0 results $((\theta_{12},\ \theta_{13},\ \theta_{23},\ \delta,\ \Delta m^2_{21},\ \Delta m^2_{31}) \sim (36.25°,\ 8.63°,\ 47°,\ 279°,\ 7.39\times10^{-5}\ \text{eV}^2,\ 2.525\times10^{-3}\ \text{eV}^2))$.

ther demonstrate how to use **FaSE-GLoBES** to obtain the constraint of any two of model parameters, and to study the ability to TDLS by MOMENT experiment.

Finally, **GLoBES** is a popular and powerful simulation tool to analysis the neutrino oscillation experiments in a simple language (AEDL), without losing too much detail. Considering the success of the flavour symmetry theory to explain the neutrino oscillations, **FaSE-GLoBES** should benefit model builders of leptonic flavour symmetry and phenomenologists for neutrino oscillation physics. We leave the flexibility for the user, and some other improvements and extensions might be done in the future.

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