

Examination of the nature of the N^* and Δ resonances via coupled-channels dynamics

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This work aims at determining the composition of certain N^* and Δ resonances, i.e., whether they are compact states formed directly by quarks and gluons, or hadronic molecules generated from the meson-baryon interaction. The information of the resonance poles is provided by a comprehensive coupled-channels approach, the Jülich-Bonn model. Thirteen states that are significant in this approach are studied. Two criteria for each state are adopted in this paper, the comparison thereof roughly indicates the model uncertainties. It is found that the conclusions for eight resonances are relatively certain: $N(1535)_{\frac{1}{2}}^{-}$, $N(1440)_{\frac{1}{2}}^{+}$, $N(1710)_{\frac{1}{2}}^{+}$, and $N(1520)_{\frac{3}{2}}^{-}$ tend to be composite, whereas $N(1650)_{\frac{1}{2}}^{-}$, $N(1900)_{\frac{3}{2}}^{+}$, $N(1680)_{\frac{5}{2}}^{+}$, and $\Delta(1600)_{\frac{3}{2}}^{+}$ tend to be compact.

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I. INTRODUCTION

Investigating the inner structures of particles is always an important topic in physics. Especially, the study of the composition is a fundamental and difficult task, which usually needs firmly established observations on both the theoretical and the experimental sides, just like the Rutherford scattering experiment which clarified the structure of atoms. In hadron physics, such studies are even more complicated due to the extremely involved dynamics and the fact that most states are unstable.

However, in the 1960s Weinberg's work on the deuteron [1] showed that the information of the composition sometimes could be derived from simple observations. Particularly, if $Z(\in [0, 1])$ is the probability of finding an elementary state in the deuteron (the “elementariness”), then the scattering length a and the effective range r are

$$\begin{aligned} a &= -\frac{2(1-Z)}{2-Z}R + \mathcal{O}(L), \\ r &= -\frac{Z}{1-Z}R + \mathcal{O}(L), \end{aligned} \quad (1)$$

where $R = (2\mu B)^{-1/2} \simeq 4.3$ fm is the binding radius, with μ the two-nucleon reduced mass and $B \simeq 2.22$ MeV the deuteron binding energy. Further, $L = 1/M_\pi \simeq 1.4$ fm, with M_π the pion mass, is the typical interaction range of two nucleons. Experiments give $a = -5.41$ fm and $r = +1.75$ fm, which support a rather small value of Z ; i.e., the deuteron should be composed by two nucleons. The derivation of this criterion is straightforward: one just needs to assume a simple formalism including two-nucleon continuous states $|\alpha\rangle_{2N}$ and possible bare states $|n\rangle$ orthogonal to the former, with the

coupling constant g between the physical deuteron $|d\rangle$ and the two-nucleon states $|\alpha\rangle_{2N}$. Then the scattering can be solved resulting in the projection probability $Z \equiv 1 - \int d\alpha |\langle d|\alpha\rangle_{2N}|^2$. Apart from the original paper [1], Ref. [2] has given a modern explanation of Weinberg's work and shown that it is compatible with the basic conceptions of modern effective field theories (EFTs). For an introduction to EFTs, see [3]. In addition, Ref. [4] interprets Weinberg's criterion via the number operators of the initial or final state particles, in order to make the connections to the EFTs without bare states. Nevertheless, the success of this criterion on the deuteron is somewhat accidental, in the sense that it can hardly be applied to other states without model assumptions or approximations: the particle must be an S -wave stable bound state near a two-body threshold.

Several decades after Weinberg's work, many new hadronic states have been found. Some of them cannot be interpreted by the naive quark model, indicating possibly the mechanism of the molecular states formed by the residual hadron-hadron interactions, e.g., hadron exchanges. For example, in pion-nucleon reactions, there are discussions on the nontrivial structure of the $N^*(1535)$ and $N^*(1440)$ (the Roper resonance); see, e.g., Refs. [5–14]. As for the heavy hadrons, after the discovery of the $X(3872)$ [15], more and more hadron exotic states were found that deviate significantly from the predictions of the conventional quark model, which can likely be interpreted by the picture of hadronic molecules [16]. A criterion of the composition is urgently needed for such states.

¹In general, one is dealing with “dynamically generated” states, which are due to the hadron-hadron (or even three-hadron) interactions. Hadronic molecules are a subclass of this type of states close to one or in between two close-by thresholds.

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Unfortunately, they are all unstable and Weinberg's criterion cannot be applied directly.

There are mainly three ways to establish an extended criterion. The first is the “pole counting rule” proposed in Ref. [17], and applied to various studies of the exotic states; see, e.g., Refs. [18–24]. It abandons the definition of the unstable state and a “probability” as the output, but rather focuses on the pole structure and the dynamics. A typical quantum mechanical potential of an S wave can only produce one near-threshold pole, in contrast to the mechanism of Castillejo, Dalitz, and Dyson [25], which produces two poles near the threshold. This method is model independent, but can only be applied to S -wave near-threshold states.

The second method, called the “spectral density function” approach, was proposed in Ref. [26] and further applied in Refs. [21,24,27–31]. This method also avoids the definition of unphysical states. It has been observed that, when lower channels are switched on, a previously stable state gains a decay width and its elementariness Z disperses into a finite probability distribution $w(z)$ of the physical energy z , called the “spectral density function.” Mathematically, $w(z)$ is just the projection of the physical scattering state (with energy z) on the bare elementary state. The elementariness for the resonance can be obtained just by collecting the function $w(z)$ near the pole mass M_R of the resonance: $Z \simeq \int_{M_R-\Delta E}^{M_R+\Delta E} w(z) dz$, with ΔE a quantity comparable with the pole width. The mathematical formalism can be constructed in principle for any partial wave, and the output is quantitative. The choice of ΔE further generates some uncertainty, and this method does not work well when the resonance is broad or the overlaps of different resonances are large.

The third way is defining a quantity similar to the Z in Weinberg's case: the projection of a Gamow state [32,33] on the bare state. The Gamow states describing resonances are zero-norm [34] and the corresponding Z is complex, without the interpretation as a “probability.” Studies on relevant mathematical properties can be found in, e.g., Refs. [35–37]. Some mathematical transformations or naive measures [4,38–44] are performed to make it a real number between 0 and 1.

In this work, we study the nature of some selected N^* and Δ states, based on the resonance parameters extracted in a recent analysis within the Jülich-Bonn (JüBo) dynamical coupled-channels approach [45]. See Refs. [10,46–60] for earlier works and other recent developments. In the JüBo model, a coupled-channels scattering equation is solved with angular momentum up to $J = 9/2$ and the channel space πN , $\pi\pi N$ (simulated by three effective channels σN , ρN , and $\pi\Delta$), ηN , $K\Lambda$, and $K\Sigma$. Photoproduction reactions are also taken into account. Reference [60] further includes the $\pi N \rightarrow \omega N$ channel. The free parameters are determined by the fit to a large collection of pion- and photon-induced data. The analytical structure is respected and the states are extracted by means of modern pole-searching procedures in the complex energy plane. Such a model provides the basic information one needs to study the elementarinesses of resonances.

Note that many of the N^* and Δ states are not in S wave, and some of them are not narrow or not close to any two-particle thresholds. We have to admit that a systematic and

accurate study of their nature is a very difficult task, and it is impossible to completely get rid of the model dependence. The reason why we still carry out this study is that the JüBo model is data driven, as the parameters are constrained by a tremendous amount of experimental input. Therefore we believe the findings with respect to the elementarinesses make sense. Specifically, the pole counting rule cannot be applied in this study, whereas the spectral density functions approach can be instructive, since these are readily extracted from the s -channel bare states of this model, and require no further assumptions or mathematical transformations. Such functions directly obtained from this model also measure the correlation between the input s -channel states and output poles, helping us to gain a better understanding of the interplay of the different components of the model. We also try our best to further estimate the uncertainties by locally constructing another spectral density function for each state from the pole positions and on-shell residues and comparing it to the function directly given by the model. Additionally, the Gamow states and the complex Z values are also calculated as a further check. This is meaningful since the conception of Gamow states is theoretically very different from the spectral density functions.

The structure of this paper is as follows. In Sec. II a solvable toy model is discussed in order to show the basic concepts and interpretation of the spectral density functions and the Gamow states. In Sec. III we briefly summarize the Jülich-Bonn model and discuss the spectral density functions and Gamow states in this framework. The numerical results are shown in Sec. IV together with pertinent discussions. Finally Sec. V contains the conclusions and provides an outlook.

II. TOY MODEL DISCUSSIONS

A. Basic descriptions

To better understand the formalism employed later, we first consider a toy model. It is nonrelativistic, and the reduced mass of the two particles is denoted as μ . Further, the dispersion relation is $\mathcal{E} = k^2/(2\mu)$ with k the momentum, and the threshold is located at the energy $\mathcal{E} = 0$. The model contains continuous free two-body states $|\psi(\mathbf{k})\rangle$ and one isolated bare state $|\psi_0\rangle$ with the energy E_0 . For simplicity we assume the bare state to be above the threshold, i.e., $E_0 > 0$. These free states form an orthogonal and complete basis:

$$\begin{aligned} \langle\psi_0|\psi(\mathbf{k})\rangle &= \langle\psi(\mathbf{k})|\psi_0\rangle = 0, \\ \langle\psi_0|\psi_0\rangle &= 1, \\ \langle\psi(\mathbf{k}')|\psi(\mathbf{k})\rangle &= (2\pi)^3 \delta^{(3)}(\mathbf{k} - \mathbf{k}'), \\ |\psi_0\rangle\langle\psi_0| + \int \frac{d^3\mathbf{k}}{(2\pi)^3} |\psi(\mathbf{k})\rangle\langle\psi(\mathbf{k})| &= \hat{I}, \end{aligned} \quad (2)$$

where $\hat{I}$ is the identity operator. Any physical state $|\Phi\rangle$ can be expressed as a linear combination of the free states with the coefficients c_0 and $\chi(\mathbf{k})$:

$$|\Phi\rangle = c_0|\psi_0\rangle + \int \frac{d^3\mathbf{k}}{(2\pi)^3} \chi(\mathbf{k})|\psi(\mathbf{k})\rangle. \quad (3)$$

The Hamiltonian consists of a free part $\hat{H}_0$ and an interaction part $\hat{H}_I$; the latter only contains the S -wave coupling between the continuous and the bare states:

$$\hat{H}_0 = E_0 |\psi_0\rangle\langle\psi_0| + \int \frac{d^3\mathbf{k}}{(2\pi)^3} \frac{k^2}{2\mu} |\psi(\mathbf{k})\rangle\langle\psi(\mathbf{k})|, \quad (4)$$

$$\hat{H}_I = \int \frac{d^3\mathbf{k}}{(2\pi)^3} gF(k, \Lambda) [|\psi(\mathbf{k})\rangle\langle\psi_0| + |\psi_0\rangle\langle\psi(\mathbf{k})|];$$

with g the coupling constant and $F(k, \Lambda)$ the regulator,

$$F(k, \Lambda) = \frac{\Lambda^2}{k^2 + \Lambda^2}, \quad (5)$$

and Λ is the cutoff parameter. In the following discussions, Λ is often assumed to be significantly larger than any other quantity with energy dimension 1, so that we can perform expansions. One can also define the free eigenstate with energy $\mathcal{E}$ accordingly:

$$|\psi(\mathcal{E})\rangle \equiv \int \frac{d^3\mathbf{k}}{(2\pi)^3} \sqrt{\frac{2\pi^2}{\mu k}} \delta\left(\frac{k^2}{2\mu} - \mathcal{E}\right) |\psi(\mathbf{k})\rangle, \quad (6)$$

which satisfies $\hat{H}_0 |\psi(\mathcal{E})\rangle = \mathcal{E} |\psi(\mathcal{E})\rangle$. In fact this toy model is equivalent or quite similar to many studies based on the spectral density functions in the literature [21,24,26–31].

B. Physical solutions

The central equation of the toy model is

$$(\hat{H}_0 + \hat{H}_I) |\Phi(\mathcal{E})\rangle = \mathcal{E} |\Phi(\mathcal{E})\rangle, \quad (7)$$

which yields

$$E_0 c_0(\mathcal{E}) + \int \frac{d^3\mathbf{k}}{(2\pi)^3} gF(k, \Lambda) \chi(\mathcal{E}, \mathbf{k}) = \mathcal{E} c_0(\mathcal{E}),$$

$$\frac{k^2}{2\mu} \chi(\mathcal{E}, \mathbf{k}) + gF(k, \Lambda) c_0(\mathcal{E}) = \mathcal{E} \chi(\mathcal{E}, \mathbf{k}). \quad (8)$$

Note that the coefficients depend on the energy eigenvalue $\mathcal{E}$.

1. Bound states

A bound state corresponds to $\mathcal{E} = -B < 0$, with the binding energy B . In this case the coefficient χ can be eliminated by the second line of Eq. (8). Then, the coefficient c_0 in the first line of Eq. (8) is also eliminated, leaving only one equation for the binding energy:

$$-B - E_0 - \Sigma(-B) = 0,$$

$$\Sigma(\mathcal{E}) \equiv \int \frac{d^3\mathbf{k}}{(2\pi)^3} \frac{g^2 F^2(k, \Lambda)}{\mathcal{E} - k^2/(2\mu)}, \quad (9)$$

where $\Sigma(\mathcal{E})$ is the self-energy of the bare state. The energy of the bound state is a pole of the propagator

$$D(\mathcal{E}) \equiv \frac{1}{\mathcal{E} - E_0 - \Sigma(\mathcal{E})}. \quad (10)$$

The bound state wave function is written as

$$|\Phi(-B)\rangle = c_0 \left[|\psi_0\rangle - \int \frac{d^3\mathbf{k}}{(2\pi)^3} \frac{gF(k, \Lambda)}{\frac{k^2}{2\mu} + B} |\psi(\mathbf{k})\rangle \right]. \quad (11)$$

The coefficient c_0 can further be determined by the normalization condition,

$$\langle\Phi(-B)|\Phi(-B)\rangle = 1, \quad (12)$$

which yields the elementariness

$$Z \equiv |c_0|^2 = \frac{1}{1 - \Sigma'(-B)}, \quad (13)$$

where the prime denotes differentiation with respect to energy. Moreover, if one expands to leading order in the inverse of the large cutoff Λ , the elementariness can be written as

$$Z = \left(1 + \frac{g^2 \mu^2}{2\pi \sqrt{2\mu B}}\right)^{-1} + \mathcal{O}(\Lambda^{-1}), \quad (14)$$

which exhibits the typical feature $g^2 \propto (Z^{-1} - 1)$ as discussed in Ref. [16]. The compositeness is further defined as

$$X \equiv \int \frac{d^3\mathbf{k}}{(2\pi)^3} |\chi(-B, \mathbf{k})|^2, \quad (15)$$

so that, according to the normalization condition (12), $Z + X = 1$.

2. Scattering states

For the scattering states, any $\mathcal{E} > 0$ can be an eigenvalue. In this case one should give the energy an infinitesimal imaginary part: $\mathcal{E} \rightarrow \mathcal{E} + i0^+$. It is convenient to eliminate the coefficient c_0 by the first line of Eq. (8), and then the second line becomes

$$\int \frac{d^3\mathbf{p}}{(2\pi)^3} \left[\frac{p^2}{2\mu} (2\pi)^3 \delta^{(3)}(\mathbf{p} - \mathbf{k}) + \frac{g^2 F(k, \Lambda) F(p, \Lambda)}{\mathcal{E} - E_0 + i0^+} \right] \chi(\mathcal{E}, \mathbf{p}) = \mathcal{E} \chi(\mathcal{E}, \mathbf{k}). \quad (16)$$

Here, χ can be understood as the wave function in momentum space. The corresponding potential operator and the Hamiltonians are

$$\hat{V} = \hat{H}_I \frac{|\psi_0\rangle\langle\psi_0|}{\mathcal{E} - E_0 + i0^+} \hat{H}_I,$$

$$\hat{h}_0 = \int \frac{d^3\mathbf{k}}{(2\pi)^3} \frac{k^2}{2\mu} |\psi(\mathbf{k})\rangle\langle\psi(\mathbf{k})|,$$

$$\hat{h} \equiv \hat{h}_0 + \hat{V}. \quad (17)$$

The scattering amplitude can be worked out with the help of the Lippmann-Schwinger equation,

$$\hat{T} = \hat{V} + \hat{V} \hat{G} \hat{T}, \quad \hat{G} \equiv (\mathcal{E} - \hat{h}_0 + i0^+)^{-1}. \quad (18)$$

The scattering amplitude is found as

$$\langle\psi(\mathbf{k})|\hat{T}|\psi(\mathbf{p})\rangle \equiv T(k, p, \mathcal{E}) = \frac{g^2 F(k, \Lambda) F(p, \Lambda)}{\mathcal{E} - E_0 - \Sigma(\mathcal{E} + i0^+)}, \quad (19)$$

which corresponds to an s -channel Feynman diagram with the dressed propagator. The physical (on-shell) scattering amplitude should be $T(q_\varepsilon, q_\varepsilon, \mathcal{E})$, shortened as $T(\mathcal{E})$, with $q_\varepsilon = \sqrt{2\mu\mathcal{E}}$ the on-shell momentum.

It is meaningful to relate χ to the amplitude. The definition of the $\hat{\chi}$ operator is

$$\chi(\mathcal{E}, \mathbf{k}) \equiv \langle \psi(\mathbf{k}) | \hat{\chi} | \psi(\mathcal{E}) \rangle, \quad (20)$$

then Eq. (16) is translated as

$$\hat{V} \hat{\chi} | \psi(\mathcal{E}) \rangle = \hat{G}^{-1} \hat{\chi} | \psi(\mathcal{E}) \rangle. \quad (21)$$

One may also rewrite the scattering equation (18), considering $\hat{G}^{-1} | \psi(\mathcal{E}) \rangle = 0$:

$$\hat{G}^{-1} (\hat{I} + \hat{G} \hat{T}) | \psi(\mathcal{E}) \rangle = \hat{V} (\hat{I} + \hat{G} \hat{T}) | \psi(\mathcal{E}) \rangle;$$

then

$$\hat{\chi} | \psi(\mathcal{E}) \rangle = \mathcal{N}(\mathcal{E}) (\hat{I} + \hat{G} \hat{T}) | \psi(\mathcal{E}) \rangle, \quad (22)$$

with $\mathcal{N}$ a normalization factor that might depend on $\mathcal{E}$. Finally the solution of χ is

$$\begin{aligned} \chi(\mathcal{E}, \mathbf{k}) = \mathcal{N}(\mathcal{E}) & \left[\sqrt{\frac{2\pi^2}{\mu k}} \delta\left(\frac{k^2}{2\mu} - \mathcal{E}\right) \right. \\ & \left. + \sqrt{\frac{\mu q_\varepsilon}{2\pi^2}} \frac{T(k, q_\varepsilon, \mathcal{E})}{\mathcal{E} - k^2/(2\mu) + i0^+} \right]. \end{aligned} \quad (23)$$

According to Eq. (8),

$$c_0(\mathcal{E}) = \mathcal{N}(\mathcal{E}) \sqrt{\frac{\mu q_\varepsilon}{2\pi^2}} \frac{T(k, q_\varepsilon, \mathcal{E})}{gF(k, \Lambda)}. \quad (24)$$

The normalization factor is determined as $\mathcal{N}(\mathcal{E}) = 1$ by the normalization condition of the scattering states,

$$\langle \Phi(\mathcal{E}') | \Phi(\mathcal{E}) \rangle = \delta(\mathcal{E} - \mathcal{E}'). \quad (25)$$

The spectral density function is defined as the probability density of finding the bare state in the scattering states:

$$\begin{aligned} w(\mathcal{E}) &= |c_0(\mathcal{E})|^2 \\ &= \rho \left| \frac{T(k, q_\varepsilon, \mathcal{E})}{gF(k, \Lambda)} \right|^2 = \rho \left| \frac{gF(q_\varepsilon, \Lambda)}{\mathcal{E} - E_0 - \Sigma(\mathcal{E})} \right|^2, \end{aligned} \quad (26)$$

with the factor $\rho = \mu q_\varepsilon / (2\pi^2)$.

Note that the effective range expansion can be performed on Eq. (19). Combining with Eq. (14), Weinberg's criterion Eq. (1) can be reproduced by this toy model.

C. Poles and criteria

1. Pole trajectories

Equations (9) and (19) indicate that the bound states are just the poles of the physical scattering amplitude in the energy region $\mathcal{E} < 0$. Besides, the dispersion relation $q_\varepsilon = \sqrt{2\mu\mathcal{E}}$ divides the complex energy plane into two Riemann sheets by the branch cut $\mathcal{E} \in [0, +\infty)$, with the physical sheet featured as $\text{Im } q_\varepsilon > 0$. The unphysical states appear as poles on the second sheet with $\text{Im } q_\varepsilon < 0$. Among those poles, virtual states lie on the real energy axis below the threshold, while the resonances have nonzero imaginary parts.

We can expand the denominator of the amplitude in Eq. (19) to leading order in the inverse of the large cutoff Λ ,

which gives a result similar to the Flatté parametrization:

$$T(\mathcal{E}) = \frac{g^2}{\mathcal{E} - E_0 + \frac{g^2 \mu \Lambda}{4\pi} + i \frac{g^2 \mu}{2\pi} q_\varepsilon} + \mathcal{O}(\Lambda^{-1}). \quad (27)$$

Defining

$$G \equiv g^2 \mu^2 / (2\pi), \quad (28)$$

the trajectories of the poles² can be plotted; see Fig. 1. For similar discussions see, e.g., Refs. [16,61]. Under the condition $E_0 > 0$, the bound states exists only if the following condition holds:

$$G > G_3 \equiv \frac{2\mu E_0}{\Lambda}. \quad (29)$$

2. Spectral density function and poles

The physical solutions of Eq. (7) also form a complete and orthogonal basis:

$$|\Phi(-B)\rangle \langle \Phi(-B)| + \int_0^{+\infty} d\mathcal{E} |\Phi(\mathcal{E})\rangle \langle \Phi(\mathcal{E})| = \hat{I}, \quad (30)$$

which, by applying to the bare state $\langle \psi_0 | \cdots | \psi_0 \rangle$, leads to

$$Z + \int_0^\infty d\mathcal{E} w(\mathcal{E}) = 1, \quad (31)$$

where Z is the elementariness of the bound state. This is the spectral density function sum rule, which indicates that the free bare state will surely be found if one explores every physical solution. It is also worth mentioning that the spectral density function is proportional to the imaginary part of the propagator in Eq. (10):

$$w(\mathcal{E}) = -\frac{1}{\pi} \text{Im } D(\mathcal{E}). \quad (32)$$

Therefore the sum rule Eq. (31) can also be understood as the Källén-Lehmann spectral representation of an isolated state.

To make closer contact to the physics, we investigate the spectral density function near the pole energy of a narrow resonance. Taking a small value of g in Eq. (27), and suppressing the quantities with g^4 and Λ^{-1} , the pair of resonance poles is located at

$$\begin{aligned} \mathcal{E}_\pm &= \left(E_0 - \frac{G\Lambda}{2\mu} \right) \mp \frac{iG\sqrt{2\mu E_0}}{\mu} \\ &\equiv \mathcal{E}_R \mp \frac{i}{2} \Gamma_R, \end{aligned} \quad (33)$$

where $\mathcal{E}_R$ and Γ_R denote the pole energy and pole width, respectively. According to Eq. (26), the spectral density function near the pole energy is

$$w(\mathcal{E} \simeq \mathcal{E}_R) = \frac{1}{\pi} \frac{\Gamma_R/2}{(\mathcal{E} - \mathcal{E}_R)^2 + \Gamma_R^2/4}. \quad (34)$$

²Note that $G \rightarrow \infty$ is not consistent with the expansion in $1/\Lambda$. However, we tentatively keep a finite value of Λ to finish the discussion, which can be regarded as another model based on the Flatté parametrization.

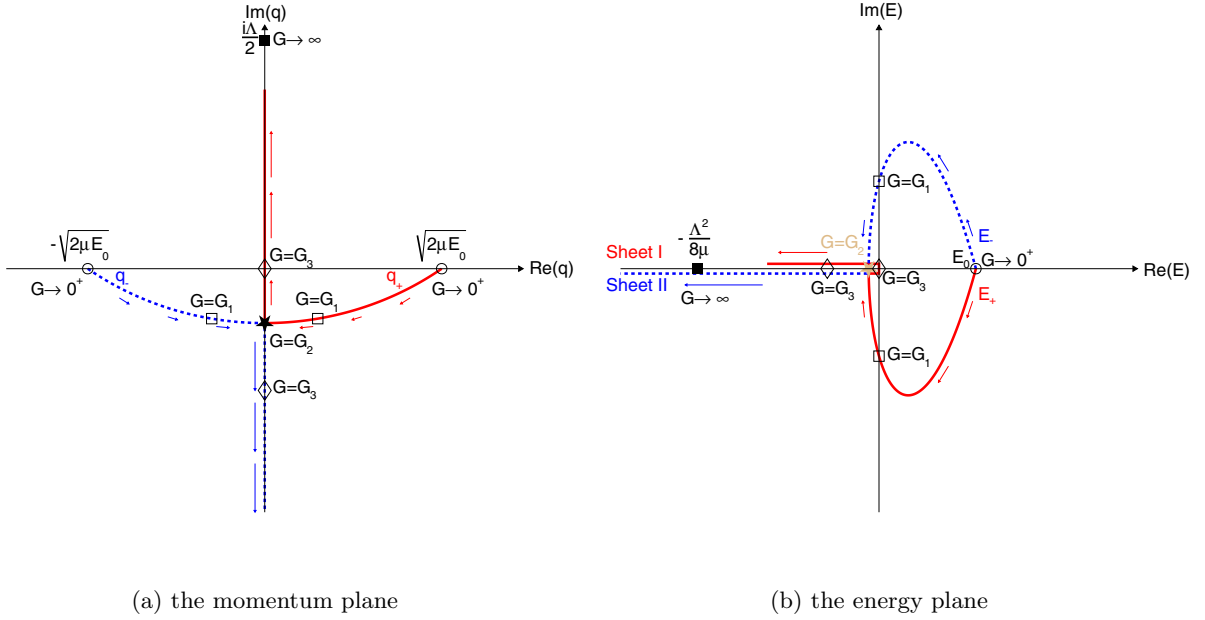


FIG. 1. The schematic plot of the pole trajectories in the toy model. The red solid (blue dashed) line denotes the trajectory of the first (second) pole. The critical points are as follows. G_1 : a resonance pole with the zero width. G_2 : a resonance pole and the conjugate pole collide, forming a second-order virtual state pole. G_3 : a bound state is produced.

In other words,

$$\lim_{g \rightarrow 0} w(\mathcal{E} \simeq \mathcal{E}_R) = \delta(\mathcal{E} - \mathcal{E}_R). \quad (35)$$

On the one hand, as $g \rightarrow 0$ the bare state decouples from the continuous states, and the Hamiltonian becomes free, so then the state at $\mathcal{E}_R = E_0$ has elementariness $Z = 1$. On the other hand, the spectral density function is concentrated totally on a single energy point. Since a finite g just makes the δ function disperse like Eq. (34), we can still assume the spectral density function near $\mathcal{E} = \mathcal{E}_R$ to be a measure of the elementariness of the resonance, which is the central idea for applying this method. Nevertheless, one also finds that the broader the resonance is, the more ambiguous this method will be. In fact, when the resonance is too broad even Eq. (34) becomes problematic.

3. Gamow states

Resonances correspond to the poles on the second Riemann sheet,

$$E_R - E_0 - \Sigma^{\text{II}}(E_R) = 0, \quad (36)$$

where the self-energy on the second sheet is evaluated by the deformed integral contour C_+ :

$$\begin{aligned} \Sigma^{\text{II}}(E_R) &= \int_{C_+} \frac{k^2 dk}{2\pi^2} \frac{g^2 F^2(k, \Lambda)}{E_R - k^2/(2\mu)} \\ &\equiv \int \frac{k^2 dk}{2\pi^2} \frac{g^2 F^2(k, \Lambda)}{(E_R - \frac{k^2}{2\mu})_+}. \end{aligned} \quad (37)$$

To reach the second sheet, the energy E_R moves across the cut $\mathcal{E} \in [0, \infty)$, and meanwhile the pole of the integrand $k_R = \sqrt{2\mu E_R}$ hits the original contour $k \in [0, \infty)$ and deforms it. The deformed contour is topologically the original contour

plus a residue term at $k_R = \sqrt{2\mu E_R}$; see Fig. 2. Then the Gamow states can be defined accordingly [see Eq. (11)]:

$$\begin{aligned} |\Phi(E_R)\rangle &\equiv c_0 \left[|\psi_0\rangle + \int \frac{d^3\mathbf{k}}{(2\pi)^3} \frac{gF(k, \Lambda)}{(E_R - \frac{k^2}{2\mu})_+} |\psi(\mathbf{k})\rangle \right], \\ |\Phi(E_R^*)\rangle &\equiv c_0^* \left[|\psi_0\rangle + \int \frac{d^3\mathbf{k}}{(2\pi)^3} \frac{gF(k, \Lambda)}{(E_R^* - \frac{k^2}{2\mu})_-} |\psi(\mathbf{k})\rangle \right], \end{aligned} \quad (38)$$

where E_R^* is the conjugate pole and the “ $-$ ” refers to the corresponding deformed contour. Note that the contour deformation in Eq. (38) does not introduce new $|\psi(\mathbf{k})\rangle$'s with complex $\mathbf{k}$. Instead it only changes the rule of the inner product: whenever another state $\langle\phi|$ does the inner product with the Gamow state, namely $\langle\phi|\Phi(E_R)\rangle$, the integral contour should be deformed. Equation (36) ensures the Gamow state is an eigenstate of the Hamiltonian, but since the Hamiltonian is Hermitian, the Gamow state cannot be normalized. The

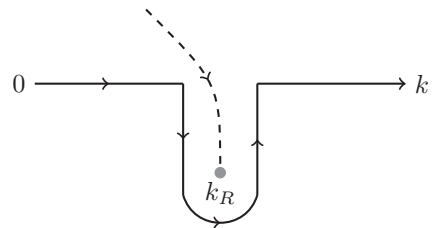


FIG. 2. Reaching the second sheet by contour deformation. The solid line stands for the integral contour, while the dashed line represents the move of the singularity $k = k_R$ when the energy moves across the cut. See the text for explanations.

normalization should be defined via the conjugate part:

$$(\Phi(E_R^*)|\Phi(E_R)) = 1. \quad (39)$$

From this condition one can only define the c_0^2 , instead of $|c_0|^2$, as the elementariness:

$$Z_R \equiv c_0^2 = \left[1 - \frac{d}{d\mathcal{E}} \Sigma^{\text{II}}(\mathcal{E} = E_R) \right]^{-1}, \quad (40)$$

which is a complex quantity without the interpretation as a probability. Then the compositeness can also be defined as $X_R \equiv 1 - Z_R$.

The compositeness of the Gamow state can also be related to the scattering amplitude. On the one hand, equivalently to Eq. (18), the scattering operator $\hat{T}$ can be rewritten as

$$\hat{T} = \hat{V} + \hat{V} \hat{\mathcal{G}} \hat{V}, \quad \hat{\mathcal{G}} \equiv (\mathcal{E} - \hat{h} + i0^+)^{-1}, \quad (41)$$

where $\hat{\mathcal{G}}$ is the full Green's operator of the Hamiltonian $\hat{h}$; see Eq. (17). On the other hand, when there is a resonance, the completeness condition (30) can be modified

$$|\Phi(E_R)\rangle(\Phi(E_R^*)| + \int_0^{+\infty} d\mathcal{E} |\tilde{\Phi}(\mathcal{E})\rangle\langle\Phi(\mathcal{E})| = \hat{I}, \quad (42)$$

where $|\tilde{\Phi}(\mathcal{E})\rangle$ is the modified scattering state, for details see Ref. [62]. Since $\hat{\mathcal{G}}|\Phi(E_R)\rangle = (\mathcal{E} - E_R)^{-1}|\Phi(E_R)\rangle$, the Laurent expansion of the T amplitude on the second sheet can be obtained from the modified completeness condition:

$$T^{\text{II}}(k, p, \mathcal{E}) = \frac{\mathfrak{r}(k)\mathfrak{r}(p)}{\mathcal{E} - E_R} + \dots, \quad (43)$$

with the off-shell residue

$$\mathfrak{r}(k) \equiv \langle\psi(\mathbf{k})|\hat{V}|\Phi(E_R)\rangle. \quad (44)$$

Moreover, $\langle\psi(\mathbf{k})|\hat{V}|\Phi(E_R)\rangle = \langle\psi(\mathbf{k})|(\hat{h} - \hat{h}_0)|\Phi(E_R)\rangle = [E_R - k^2/(2\mu)]_+ \langle\psi(\mathbf{k})|\Phi(E_R)\rangle$, where “+” denotes the deformed contour; one obtains the compositeness from the off-shell residue,

$$\begin{aligned} X(E_R) &= \int \frac{d^3\mathbf{k}}{(2\pi)^3} \langle\psi(\mathbf{k})|\Phi(E_R)\rangle^2 \\ &= \int \frac{d^3\mathbf{k}}{(2\pi)^3} \frac{\mathfrak{r}^2(k)}{[E_R - k^2/(2\mu)]_+^2}. \end{aligned} \quad (45)$$

According to Eq. (19), the off-shell residue in this toy model is $\mathfrak{r}(k) = gF(k, \Lambda)[1 - \Sigma^{\text{II}}]^{-1/2}$, so Eq. (45) gives exactly the same result as Eq. (40).

D. Coupled-channels extension

Here, we slightly extend the toy model to a coupled-channels situation. We keep the channel discussed above as channel “1”, with all the notations remaining the same. In addition, there is a lower channel with index “0”. The threshold of the lower channel is $-\Delta E < 0$. The free continuous states of channel 0 are labeled as $|\psi^0(\mathbf{k})\rangle$, which are orthogonal to the $|\psi(\mathbf{k})\rangle$. For a given energy $\mathcal{E}$, the on-shell momentum of the channel 1 remains the same, i.e., $q_\mathcal{E} = \sqrt{2\mu\mathcal{E}}$, while for channel 0 it is $q_{0\mathcal{E}} = \sqrt{2\mu_0(\mathcal{E} + \Delta E)}$. Equivalently to Eq. (6),

we define another energy eigenstate for the lower channel 0:

$$\begin{aligned} |\psi^0(\mathcal{E})\rangle &= \int \frac{d^3\mathbf{k}_0}{(2\pi)^3} \sqrt{\frac{2\pi^2}{\mu_0 k_0}} \\ &\times \delta\left(\frac{k_0^2}{2\mu_0} - \mathcal{E} - \Delta E\right) |\psi^0(\mathbf{k}_0)\rangle. \end{aligned} \quad (46)$$

Besides, there is still only one bare state $|\psi_0\rangle$.

The Hamiltonian now is³

$$\begin{aligned} \hat{H}_0 &= E_0 |\psi_0\rangle\langle\psi_0| + \int \frac{d^3\mathbf{k}}{(2\pi)^3} \frac{k^2}{2\mu} |\psi(\mathbf{k})\rangle\langle\psi(\mathbf{k})| \\ &+ \int \frac{d^3\mathbf{k}_0}{(2\pi)^3} \left(\frac{k_0^2}{2\mu_0} - \Delta E\right) |\psi^0(\mathbf{k}_0)\rangle\langle\psi^0(\mathbf{k}_0)|, \\ \hat{H}_I &= \int \frac{d^3\mathbf{k}}{(2\pi)^3} gF(k, \Lambda) |\psi(\mathbf{k})\rangle\langle\psi_0| \\ &+ \int \frac{d^3\mathbf{k}_0}{(2\pi)^3} g_0 F(k_0, \Lambda) |\psi^0(\mathbf{k}_0)\rangle\langle\psi_0| + \text{H.c.}, \end{aligned} \quad (47)$$

where “H.c.” refers to the Hermitian conjugation. The physical scattering state is

$$\begin{aligned} |\Phi(\mathcal{E})\rangle &= c_0(\mathcal{E}) |\psi_0\rangle + \int \frac{d^3\mathbf{k}}{(2\pi)^3} \chi(\mathcal{E}, \mathbf{k}) |\psi(\mathbf{k})\rangle \\ &+ \int \frac{d^3\mathbf{k}_0}{(2\pi)^3} \chi_0(\mathcal{E}, \mathbf{k}_0) |\psi^0(\mathbf{k}_0)\rangle. \end{aligned} \quad (48)$$

The eigenstates can still be solved from Eq. (7). In this case, the amplitude is

$$T_{ij}(k, p, \mathcal{E}) = \frac{g_i g_j F(k, \Lambda) F(p, \Lambda)}{\mathcal{E} - E_0 - \tilde{\Sigma}(\mathcal{E} + i0^+)}, \quad (49)$$

with $i, j = 0, 1$ as the channel indices, and $g_1 = g$. Meanwhile the propagator is

$$D(\mathcal{E}) = \frac{1}{\mathcal{E} - E_0 - \tilde{\Sigma}(\mathcal{E} + i0^+)}, \quad (50)$$

with the two-channel self-energy

$$\tilde{\Sigma}(\mathcal{E}) = \int \frac{d^3\mathbf{k}}{(2\pi)^3} \left[\frac{g_0^2 F^2(k, \Lambda)}{\mathcal{E} - \frac{k^2}{2\mu_0} + \Delta E + i0^+} + \frac{g^2 F^2(k, \Lambda)}{\mathcal{E} - \frac{k^2}{2\mu} + i0^+} \right]. \quad (51)$$

Due to the same reason as in the single channel case, for both $i = 0, 1$ ($|\psi^i(\mathcal{E})\rangle = |\psi(\mathcal{E})\rangle$)

$$\hat{\chi} |\psi^i(\mathcal{E})\rangle = \mathcal{N}_i(\mathcal{E}) (\hat{I} + \hat{G}\hat{T}) |\psi^i(\mathcal{E})\rangle. \quad (52)$$

The correspondence of the function χ to the $\hat{\chi}$ operator is not unique in the coupled-channels system, and the normalization factor $\mathcal{N}$ is no longer trivially 1. Fortunately the spectral density function is easily obtained from Eq. (32):

$$w(\mathcal{E}) = \frac{g_0^2 \mu_0 q_{0\mathcal{E}} F^2(q_{0\mathcal{E}}, \Lambda) + g^2 \mu q_\mathcal{E} F^2(q_\mathcal{E}, \Lambda) \Theta(\mathcal{E})}{2\pi^2 |\mathcal{E} - E_0 - \tilde{\Sigma}(\mathcal{E})|^2}, \quad (53)$$

³The cutoff Λ is chosen to be the same for the two channels.

where $\Theta(\mathcal{E})$ is the Heaviside step function. It is easy to verify that, when $g_0 = 0$, Eq. (26) is recovered.

At last we investigate a “quasibound state” (qb) in the two-channel case. We assume when channel 0 is switched off, the channel 1 forms a bound state at $\mathcal{E} = -B$, with the binding energy $B \ll \Delta E$. As the coupling g_0 grows from zero but is still small, the bound state becomes a narrow resonance. Such a pole lies on the Riemann sheet defined by $\text{Im } q_{0\epsilon} < 0$, $\text{Im } q_\epsilon > 0$, i.e., the physical sheet for channel 1 but the unphysical sheet for channel 0. Expanding Eq. (51) and suppressing the $\mathcal{O}(g_0^4)$ and $\mathcal{O}(\Lambda^{-1})$ terms, the location of the pole is (the conjugate one is omitted)

$$\begin{aligned}\mathcal{E}_+ &= E_{qb} - \frac{i}{2}\Gamma_{qb}, \\ E_{qb} &= -B - \frac{g_0^2 \mu_0 \Lambda}{4\pi} Z_B, \\ \Gamma_{qb} &= \frac{g_0^2 \mu_0}{\pi} \sqrt{2\mu_0(\Delta E - B)} Z_B,\end{aligned}\quad (54)$$

where $Z_B = [1 - \Sigma'(-B)]^{-1}$ is the elementariness of the bound state when the channel 0 is switched off. Moreover, the spectral density function near the pole energy is

$$w(\mathcal{E} \simeq E_{qb}) = \frac{Z_B}{\pi} \frac{\Gamma_{qb}/2}{(\mathcal{E} - E_{qb})^2 + (\Gamma_{qb}/2)^2}. \quad (55)$$

Hence in the narrow resonance limit,

$$\lim_{g_0 \rightarrow 0} w(\mathcal{E} \simeq E_{qb}) = Z_B \delta(\mathcal{E} + B). \quad (56)$$

This indicates again that the spectral density function near the pole energy carries the information of the elementariness, which is fully compatible with the elementariness of the bound state (Weinberg’s criterion). When g_0 is finite, the delta function in Eq. (56) disperses and becomes a finite distribution.

The elementariness of a quasibound state with finite width can be evaluated as $\int_{E_{qb}-\Delta E}^{E_{qb}+\Delta E} w(\mathcal{E}) d\mathcal{E}$. Actually the choice of the integral interval (ΔE) is ambiguous. Note that the Z_B in Eq. (55) is the crucial quantity we are interested in. So for the pole at $\mathcal{E} = E_R - \frac{i}{2}\Gamma_R$, we define the “Breit-Wigner” spectral density function only according to the pole position:

$$\text{BW}(\mathcal{E}) \equiv \frac{1}{\pi} \frac{\Gamma_R/2}{(\mathcal{E} - E_R)^2 + (\Gamma_R/2)^2}. \quad (57)$$

Then the elementariness is modified as

$$Z \simeq \frac{\int_{E_R-\Delta E}^{E_R+\Delta E} w(\mathcal{E}) d\mathcal{E}}{\int_{E_R-\Delta E}^{E_R+\Delta E} \text{BW}(\mathcal{E}) d\mathcal{E}}. \quad (58)$$

Since, in the narrow resonance limit of Eq. (55), Eq. (58) gives exactly the quantity Z_B no matter what value ΔE is, it is expected that the dependence of the result on ΔE is weakened. However, again, when the resonance is too broad, Eq. (55) does not hold and this evaluation is also ambiguous.

The Gamow states for coupled channels can also be discussed in a totally similar manner to the single channel case. We skip the details here.

III. APPLICATION TO THE JÜLICH-BONN MODEL

A. Short description of the model

The Jülich-Bonn model is a comprehensive coupled-channels model, which currently contains the hadronic channels πN , $\pi\pi N$, ηN , $K\Lambda$, $K\Sigma$, and ωN . The $\pi\pi N$ system is simulated by three effective channels, namely $\pi\Delta$, σN , and ρN . The thresholds are shown in Fig. 3.

In this model the reactions are studied through the following scattering equation:

$$\begin{aligned}T_{\mu\nu}(p'', p', z) &= V_{\mu\nu}(p'', p', z) + \sum_{\kappa} \int_0^{\infty} p^2 dp V_{\mu\kappa}(p'', p, z) \\ &\quad \times G_{\kappa}(p, z) T_{\kappa\nu}(p, p', z),\end{aligned}\quad (59)$$

where T denotes the scattering amplitude, V denotes the interaction kernel (potential), p' and p'' are the three-momenta of the initial and final states in the center-of-mass frame, respectively, and z is the center-of-mass energy. The channel labels μ , ν , and κ denote the meson-baryon system with specific isospin (I), angular momentum (J , up to $9/2$), spin (S), and orbital angular momentum L . $G_{\kappa}(p, z)$ is the propagator of the intermediate channel:

$$G_{\kappa}(z, p) = [z - E_{\kappa} - \omega_{\kappa} - \Sigma_{\kappa}(z, p) + i0^+]^{-1}, \quad (60)$$

where Σ_{κ} is the self-energy of the unstable particle (ρ , σ , or Δ) in the effective channel κ . When κ is not an effective channel, $\Sigma_{\kappa} = 0$. Further, E_{κ} , ω_{κ} denote the energies of the baryon and the meson in channel κ , respectively, with the relativistic dispersion relation, e.g., $E_{\kappa} = \sqrt{p^2 + M_{\kappa}^2}$. The whole formalism is established in the framework of time-ordered perturbation theory [63], which, together with an expansion in the partial-wave basis, ensures the integral in Eq. (59) is only one-dimensional.

To simplify the calculation, we further perform the following separation:

$$T = T^{NP} + T^P, \quad (61)$$

where T^P and T^{NP} are the pole part and the nonpole part, respectively. The nonpole part T^{NP} is generated only by the potentials without s -channel bare states (V^{NP}):

$$\begin{aligned}T_{\mu\nu}^{NP}(p'', p', z) &= V_{\mu\nu}^{NP}(p'', p', z) + \sum_{\kappa} \int_0^{\infty} p^2 dp V_{\mu\kappa}^{NP} \\ &\quad \times (p'', p, z) G_{\kappa}(p, z) T_{\kappa\nu}^{NP}(p, p', z),\end{aligned}\quad (62)$$

whereas the pole part is

$$T_{\mu\nu}^P(p'', p', z) = \sum_{i,j} \Gamma_{\mu,i}^a(p'') D_{ij}(z) \Gamma_{\nu,j}^c(p'). \quad (63)$$

Here, i, j are the indices of the s -channel bare states in a given partial wave, and $\Gamma_{\mu,i}^{a(c)}$ is the dressed vertex function describing the annihilation (creation) of the i th state to channel μ . The dressed propagator of the s -channel state is denoted as D ,

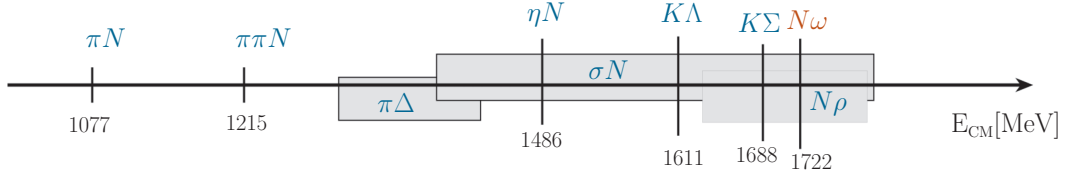


FIG. 3. Thresholds of the scattering channels currently considered in the Jülich-Bonn model as function of the center-of-mass energy. The ωN channel is not considered in this work; see the discussions in the beginning of Sec. IV.

which is related to the self-energies Σ_{ij} :

$$\begin{aligned}\Gamma_{\mu,i}^a(p'') &= \gamma_{\mu,i}^a(p'') \\ &+ \sum_{\kappa} \int_0^\infty p^2 dp T_{\mu\kappa}^{NP}(p'', p, z) G_{\kappa}(p, z) \gamma_{\kappa,i}^a(p), \\ \Gamma_{\nu,j}^c(p') &= \gamma_{\nu,j}^c(p') \\ &+ \sum_{\kappa} \int_0^\infty p^2 dp \gamma_{\kappa,j}^c(p) G_{\kappa}(p, z) T_{\kappa\nu}^{NP}(p, p', z), \\ D_{ij}^{-1}(z) &= \delta_{ij}(z - m_i^b) - \Sigma_{ij}(z), \\ \Sigma_{ij}(z) &= \sum_{\kappa} \int_0^\infty p^2 dp \gamma_{\kappa,i}^c(p) G_{\kappa}(p, z) \Gamma_{\kappa,j}^a(p).\end{aligned}\quad (64)$$

Here, the γ 's are bare vertices and m_i^b is the bare mass of the i th bare state. All the relevant expressions can be found in Ref. [60] and its Supplemental Material. Photoproduction reactions are described in a semiphenomenological approach with $T_{\mu\nu}$ as the hadronic final-state interaction; see Ref. [53] for details. The free parameters in this model are determined by a global fit to almost all available data. The resonances are extracted by scanning the complex energy plane on the unphysical Riemann sheet of the full T matrix; for details see Ref. [50]. Note that the decomposition of the scattering amplitude into a pole and a nonpole part in Eq. (61) is ambiguous since resonance poles can be produced not only by the s -channel bare states but can also be dynamically generated in the nonpole part in Eq. (62). However, due to the complicated multichannel space and strong dressing effects inherent in the model, a clear interpretation of a certain observed pole as an s -channel or “genuine” state in contrast to a dynamical one is often very difficult.

B. Spectral density functions of the model

The Jülich-Bonn model can be rewritten in terms of the states in the Hilbert space. The free continuous states are denoted as $|\mu, p\rangle$, which satisfy

$$\langle \nu, q | \mu, p \rangle = p^{-2} \delta_{\mu\nu} \delta(p - q), \quad (65)$$

whereas the s -channel bare states are $|\Phi_n\rangle$, with

$$\langle \Phi_m | \Phi_n \rangle = \delta_{mn}. \quad (66)$$

Further conditions are

$$\langle \Phi_n | \mu, p \rangle = \langle \mu, p | \Phi_n \rangle = 0, \quad (67)$$

and

$$\sum_n |\Phi_n\rangle \langle \Phi_n| + \sum_{\mu} \int_0^\infty p^2 dp |\mu, p\rangle \langle \mu, p| = \hat{1}. \quad (68)$$

The Hamiltonian is written as

$$\hat{H} = \hat{H}_0 + \hat{H}_I, \quad (69)$$

with the free part

$$\hat{H}_0 = \sum_j m_j^b |\Phi_j\rangle \langle \Phi_j| + \sum_{\mu} \int p^2 dp (E_{\mu} + \omega_{\mu}) |\mu, p\rangle \langle \mu, p| \quad (70)$$

and the interaction part

$$\begin{aligned}\hat{H}_I &= \sum_i \sum_{\mu} \int p'^2 dp' [\gamma_{\mu,i}^a(p'') |\mu, p''\rangle \langle \Phi_i| + \text{H.c.}] \\ &+ \sum_{\mu,\nu} \iint dp' dp'' (p' p'')^2 V_{\mu\nu}^{NP}(p'', p') |\mu, p''\rangle \langle \nu, p'|.\end{aligned}\quad (71)$$

The scattering equation (59) is fully equivalent to the eigenequation of the energy z :

$$(\hat{H}_0 + \hat{H}_I) |z\rangle = z |z\rangle, \quad (72)$$

where a physical scattering state is denoted as $|z\rangle$. The expansion is very similar to Eq. (48):

$$|z\rangle = \sum_i c_i(z) |\Phi_i\rangle + \sum_{\alpha} \int k^2 dk \chi_{\alpha}(z, k) |\alpha, k\rangle. \quad (73)$$

The spectral density function of the scattering state on the i th s -channel bare state can be defined as

$$w_i(z) \equiv |\langle \Phi_i | z \rangle|^2 = |c_i(z)|^2. \quad (74)$$

The sum rule for every i is

$$\mathcal{Z}_{B,i} + \int_{m_{\pi} + m_N}^{\infty} dz w_i(z) = 1, \quad (75)$$

where $\mathcal{Z}_{B,i}$ is the partial elementariness of a possible bound state on the i th s -channel bare state. Note that in the Jülich-Bonn model the only bound state is the nucleon itself in the P_{11} wave of πN scattering.

The Jülich-Bonn model is much more complicated than the toy model discussed previously, so the analytical expressions cannot be solved explicitly. Nevertheless the spectral density functions in Eq. (74) can still be obtained by the imaginary part of the propagator in Eq. (64),

$$w_i(z) = -\frac{1}{\pi} \text{Im} D_{ii}(z), \quad (76)$$

and the propagator can be calculated numerically. For the resonance pole at $z = M_R - i\Gamma_R/2$, the elementariness corresponding to the i th bare state, denoted as $\mathcal{Z}_i$, is evaluated as

$$\mathcal{Z}_i \simeq \frac{\int_{M_R - \Gamma_R}^{M_R + \Gamma_R} w_i(z) dz}{\int_{M_R - \Gamma_R}^{M_R + \Gamma_R} BW(z) dz}. \quad (77)$$

Note that the ΔE in Eq. (58) is chosen as the pole width of the resonance here. To avoid the uncertainties of the overlap between two resonances in the same partial wave, the value of ΔE should not be too large. At last, when there are more than one bare states in the same partial wave, the “total elementariness” can be estimated as

$$\mathcal{Z} = 1 - \prod_i (1 - \mathcal{Z}_i). \quad (78)$$

C. Estimation of the model uncertainties

We emphasize that the number of resonance poles does not correspond to the number of s -channel bare states. The free parameters of the model can move the pole of an s -channel state to distant regions in the complex plane beyond the reach of the pole searching, while the T^{NP} part in Eq. (61) can also contain “dynamically generated poles.” Actually, the separation of the amplitude and the number of s -channel states are model dependent, which also holds for Eq. (77) that does depend on those, and hence it can only be regarded as a naive indication from the model. Despite the difficulties of getting rid of these model dependences, we will try to give some rough estimation for the uncertainties.

1. Locally constructed spectral density functions

The first proposal is to locally construct (lc) a spectral density function for every individual resonance, using only the pole positions and the residues. We can first construct the on-shell T amplitude near the pole:

$$T_{\alpha\beta}^{\text{lc}}(z) = \frac{c g_\alpha g_\beta f_\alpha^a(q_{\alpha z}) f_\beta^c(q_{\beta z})}{z - M_0 - \sum_\kappa g_\kappa^2 L_\kappa(z)} + \dots, \quad (79)$$

where $q_{\alpha z}$ is the on-shell momentum for energy z in channel α , the g 's are real coupling constants, $f_\alpha^{a(c)}$ is the vertex function of the resonance annihilation (creation) without the coupling constant, M_0 is a mass parameter to be determined by the pole position, “ $\dots$ ” refers to the other contributions apart from the pole, and L_κ is the loop function of channel κ with respect to the vertex functions:

$$L_\kappa(z) \equiv \int_0^\infty p^2 dp G_\kappa(p, z) f_\alpha^a(q_{\kappa z}) f_\alpha^c(q_{\kappa z}). \quad (80)$$

Equation (79) can be regarded as a generalization of the Laurent expansion. For simplicity we just choose the form of the vertex function f the same as the bare vertices in Jülich-Bonn model (together with the regulator, with the cutoff parameters set as the pole masses). The σN channel does not have bare vertices in this model. Instead, we take $f_{\sigma N}^{a,c}(q) \sim q^l$ with l the angular momentum. The number c is an extra compensation factor. Specifically, as we already have the residue r_α at the pole $z = M_R - i\Gamma_R/2$, all the parameters can be given (here,

the subscript “1” refers to the πN channel):

$$\begin{aligned} h_\kappa &\equiv \frac{g_\kappa^2}{g_1^2} = \left| \frac{r_\kappa f_1^a}{r_1 f_\kappa^a} \right|^2, \\ g_1^2 &= -\frac{\Gamma_R}{2 \sum_\kappa h_\kappa \text{Im}(L_\kappa^{\text{II}})}, \\ M_0 &= M_R - g_1^2 \sum_\kappa h_\kappa \text{Re}(L_\kappa^{\text{II}}), \\ c &= \frac{r_1^2}{g_1^2 f_1^a f_1^c} \left(1 - g_1^2 \sum_\kappa h_\kappa \frac{d}{dz} L_\kappa^{\text{II}} \Big|_{z=M_R - i\Gamma_R/2} \right). \end{aligned} \quad (81)$$

This parametrization cannot reproduce the phase of the residues. However, the phases mainly affect the interference behavior of a resonance with the others. Here we construct the amplitude only for every individual state with the energy near the pole mass. Finally, the locally constructed spectral density function is

$$w^{\text{lc}}(z) = -\frac{1}{\pi} \text{Im} \left[z - M_0 - \sum_\kappa g_\kappa^2 L_\kappa(z) \right]^{-1}. \quad (82)$$

According to the asymptotic behavior of L_κ and the analytic properties, the sum rule in Eq. (31) still holds for this construction.

Note that Eq. (79) is still too simple to simulate the amplitude in Jülich-Bonn model. This formalism fails if $\sum_\kappa h_\kappa \text{Im} L_\kappa^{\text{II}} > 0$, which does not happen when the state is narrow enough. In this case the uncertainty is of course large. We can employ a “plan B” and directly assign the couplings in Eq. (82) as the absolute values of the dimensionless normalized residues defined by the Particle Data Group [64]:

$$g_\kappa^2 \rightarrow \left| \sqrt{\frac{2\pi \rho_\kappa}{\Gamma_R}} r_\kappa \right|^2, \quad (83)$$

where $\rho_\kappa = q_{\kappa z} E_\kappa \omega_\kappa / z$ is a kinematic factor. The fixed g'_κ s do not automatically lead to the correct pole position. In fact, we have to give M_0 an imaginary part compensating for the other implicit effects. This destroys the sum rule in Eq. (31) and sometimes makes the spectral density function negative. Before applying the plan B, one has to check if such bad features are significant.

2. Complex compositenesses of the Gamow states

The second proposal is to adopt the Gamow states and their compositenesses. Totally similar to Eq. (45), the compositeness for channel κ is

$$X_\kappa \equiv \int_{C_+} p^2 dp \mathfrak{r}_\kappa^2(p) G_\kappa^2(p, z_{\text{pole}}), \quad (84)$$

where $\mathfrak{r}_\kappa$ is the off-shell residue from the Jülich-Bonn model, and C_+ is the deformed contour to ensure the correct Riemann sheet. That means when the pole is lower than the threshold of channel κ , the integral contour is $[0, +\infty]$. When the pole is higher than the threshold, the extra contributions from the

TABLE I. Partial waves, the number of s -channel bare states (N_s), and the states to be studied. The $L_{2I\ 2J}$ notation is only for the πN channel.

$I(J^P)$	$L_{2I\ 2J}$	N_s	States to be studied
$\frac{1}{2}(\frac{1}{2}^-)$	S_{11}	2	$N(1535), N(1650)$
$\frac{3}{2}(\frac{1}{2}^-)$	S_{31}	1	$\Delta(1620)$
$\frac{1}{2}(\frac{1}{2}^+)$	P_{11}	2	$N(1440), N(1710)$
$\frac{1}{2}(\frac{3}{2}^+)$	P_{13}	2	$N(1720), N(1900)$
$\frac{3}{2}(\frac{3}{2}^+)$	P_{33}	2	$\Delta(1232), \Delta(1600)$
$\frac{1}{2}(\frac{3}{2}^-)$	D_{13}	1	$N(1520)$
$\frac{3}{2}(\frac{3}{2}^-)$	D_{33}	1	$\Delta(1700)$
$\frac{1}{2}(\frac{5}{2}^-)$	D_{15}	1	$N(1675)$
$\frac{1}{2}(\frac{5}{2}^+)$	F_{15}	1	$N(1680)$

deformed contour are estimated by the following extrapolation:

$$\tau_\kappa(p) \rightarrow \frac{r_\kappa}{f_\kappa(p_{\text{pole}})} f_\kappa^a(p), \quad (85)$$

with f the bare vertex function just used in Eq. (79). This extrapolation ensures that the on-shell residue r_κ is correctly reproduced. The complex elementariness of a state is

$$Z = 1 - \sum_\kappa X_\kappa. \quad (86)$$

Last, to make a comparison to the probabilities from the spectral density functions, we can also use the naive measure proposed in Ref. [42] to get the rates between 0 and 1:

$$\tilde{X}_\kappa \equiv \frac{|X_\kappa|}{\sum_\alpha |X_\alpha| + |Z|}, \quad \tilde{Z} \equiv \frac{|Z|}{\sum_\alpha |X_\alpha| + |Z|}. \quad (87)$$

Note that the basic philosophy of this method is very different from that of the spectral density functions, so we do not expect perfect matches of those results. However, it is the difference in nature that makes the adoption of the Gamow states somehow instructive as the complex quantities only depend on the pole parameters, hence they are free from the ambiguities of choosing the integral interval, or the overlaps of the resonances. Another advantage is that the complex compositenesses directly show which channel dominates the composition.

IV. NUMERICAL RESULTS

A. Selection of the resonances

There are two recent results of the Jülich-Bonn model: the “JüBo_omegaN” solution [60], which is based on the study of purely hadronic, pion-induced reactions, with the channel space extended to ωN , and the “JüBo2022” solution [45], which includes also photoproduction reactions with $\gamma p \rightarrow K \Sigma$ newly considered, while the ωN channel is absent. In this work we use the output from the latter. Actually some states are not significant in the hadronic part, e.g., $N(1900) \frac{3}{2}^+$, but play an important role in photoproduction. Moreover, the

TABLE II. Residues of the selected N^* states in the σN channel given by the “JüBo2022” solution.

State	$r_{\sigma N} (10^{-6} \text{ MeV}^{-1/2})$
$N(1535) \frac{1}{2}^-$	$0.294 - 0.207i$
$N(1650) \frac{1}{2}^-$	$-0.163 + 0.296i$
$N(1440) \frac{1}{2}^+$	$3.948 - 8.295i$
$N(1710) \frac{1}{2}^+$	$1.303 - 14.334i$
$N(1720) \frac{3}{2}^+$	$-0.564 - 0.212i$
$N(1900) \frac{3}{2}^+$	$2.707 - 5.367i$
$N(1520) \frac{3}{2}^-$	$0.935 + 3.844i$
$N(1675) \frac{5}{2}^-$	$-0.083 - 0.032i$
$N(1680) \frac{5}{2}^+$	$0.198 - 0.091i$

quality and quantity of the available photoproduction data supersede the ones of the pion-induced data by far. We therefore consider the resonance parameters extracted in JüBo2022 to be more reliable. Note, however, that we cannot exclude the possibility that some of the results might change when the ωN channel (or further channels) is (are) considered in the future.

The results for higher partial waves are always less stable [45]. Therefore we only study the $J \leq 5/2$ N^* states and the $J \leq 3/2$ Δ states. Although there are two Δ states with $J = 5/2$ listed by the Particle Data Group [64], the $\Delta(1905) \frac{5}{2}^+$ and $\Delta(1930) \frac{5}{2}^-$, we exclude those states since they proved to be rather unstable in recent JüBo analyses; cf. the discussion in Ref. [45] for details. This may be related to the fact that the database for the $I = 3/2$ channels is still smaller than for the $I = 1/2$ channels. In addition, the states with widths $\Gamma_R > 300$ MeV are not considered, since the spectral density functions does not work well for broad states. Specifically, in this model the $\Delta(1910) \frac{3}{2}^+$ state is always not significant and very broad ($\Gamma_R \simeq 550$ MeV). It is thus excluded.

Moreover, we expect that the uncertainties of the resonance parameters estimated in Ref. [45] are negligible compared to the systematic uncertainties from the criteria of the elementariness. Therefore we do not use the uncertainties of the resonance parameters in the current calculations.

In summary, 13 states out of 25 in Ref. [45] are selected and summarized in Table I. In the P_{11} wave there are two bare s -channel states, of which the first is the bare nucleon. By applying the sum rule Eq. (75), the nucleon wave function renormalization constant is $Z_N \simeq 0.54$.⁴ Note that we also use the bare nucleon to measure the elementarinesses of the other states in the P_{11} wave.

Last, as mentioned in the last section, the estimations of the uncertainties require the residues as input. The residues of the σN channel were not published in Ref. [45], and are listed here in Table II. Most of the values are much smaller (of order $10^{-6} \text{ MeV}^{-1/2}$) than the residues of the other channels (of order $10^{-3} \text{ MeV}^{-1/2}$), since in the JüBo2022 model the

⁴In our model the bare mass and bare couplings of the nucleon are always adjusted such that its physical mass is 938 MeV.

TABLE III. The elementarinesses of the selected N^* states. The label “(F)” means that the local construction of Eq. (81) has failed and Eq. (83) is used instead.

State	Pole position (MeV)	$\mathcal{Z}_1$	$\mathcal{Z}_2$	$\mathcal{Z}_{\text{tot}}$	$\mathcal{Z}^{\text{lc}}$	$\tilde{Z}$
$N(1535) \frac{1}{2}^-$	$1504 - 37i$	24.8%	5.6%	29.0%	50.8%	39.4%
$N(1650) \frac{1}{2}^-$	$1678 - 64i$	13.4%	91.7%	92.8%	70.5%	8.5%
$N(1440) \frac{1}{2}^+$	$1353 - 102i$	48.7%	1.7%	49.5%	31.5%	36.9%
$N(1710) \frac{1}{2}^+$	$1605 - 58i$	11.5%	10.3%	20.6%	10.2%	40.3%
$N(1720) \frac{3}{2}^+$	$1726 - 93i$	34.1%	68.5%	79.3%	62.5%	41.4%
$N(1900) \frac{3}{2}^+$	$1905 - 47i$	19.9%	100%	100%	99.9%	38.5%
$N(1520) \frac{3}{2}^-$	$1482 - 63i$	29.4%		29.4%	7.2%	40.4%
$N(1675) \frac{5}{2}^-$	$1652 - 60i$	16.6%		16.6%	100% (F)	61.8%
$N(1680) \frac{5}{2}^+$	$1657 - 60i$	67.9%		67.9%	69.9%	55.0%

bare resonance vertices do not couple to σN . The σN residues receive contributions from coupled-channels and nonpole effects and are thus extremely small.

B. Analyses of the N^* states

For each selected state, the elementarinesses from three methods will be shown, namely $\mathcal{Z}_i$ ($i = 1, 2, \text{tot}$) from the naive indication of the model [Eq. (76)], $\mathcal{Z}^{\text{lc}}$ from the local construction [Eq. (82)], and the naive measure $\tilde{Z}$ of the Gamow states [Eq. (87)]. The first is further labeled by subscripts: $\mathcal{Z}_1$ ($\mathcal{Z}_2$) for the partial elementariness on the first (second) bare s -channel state, and $\mathcal{Z}_{\text{tot}}$ for the “total” elementariness in Eq. (78). Note that in a complex coupled-channels environment with strong dressing effects as in the JüBo model, in cases with two states per partial wave it is often difficult to identify a pole as unambiguously induced by a specific bare s -channel state. Therefore we list $\mathcal{Z}_1$ and $\mathcal{Z}_2$ for each pole.

The elementarinesses of the selected N^* states are summarized in Table III. The complex quantities of the Gamow states are listed in Tables IV and V. Since the σN residues are always very small, we skip the discussion of the compositenesses in the σN channel. The spectral density functions are plotted in Fig. 4. The results are explained and discussed one by one in what follows.

1. $N(1535) \frac{1}{2}^-$

As shown in Table III, all results from the three different methods indicate that the elementariness of $N^*(1535)$ is not large. As shown in Fig. 4, the peak structure near the pole position only shows up in the first spectral density function, which means that the $N^*(1535)$ pole is highly related to the first s -channel state in this model, and the mixture between the two bare states is weak. The locally constructed spectral density function is bigger than that of the model. However, it is still significantly smaller than the Breit-Wigner peak. Note that the integral intervals $[M_R - \Gamma_R, M_R + \Gamma_R]$ for $N^*(1535)$ and $N^*(1650)$ have an overlap, which increases the uncertainties. On the other hand, Table IV shows a large complex compositeness of the Gamow state in the ηN channel, with the naive measure of 35.8% for the composition. It is suggested in

this work that the $N^*(1535)$ tends to be composite. Actually in some other approaches the $N^*(1535)$ resonance can be dynamically generated; see, e.g., Refs. [5,9,12,65].

In fact, in the JüBo_omegaN solution [60] the coupling of the $N^*(1535)$ to the ωN channel is rather large, even though the pole position is not close to the ωN threshold. This large coupling, as well as the ωN composition, will be investigated in the future when studying ωN photoproduction.

2. $N(1650) \frac{1}{2}^-$

Just like the $N^*(1535)$, in some models, e.g., Refs [9,12,65], the $N^*(1650)$ can also be dynamically generated. But here in our model it seems to be highly related to the second bare state. The spectral density function of this state, no matter if directly from the model or locally constructed, suggests a high elementariness; see Fig. 4.

However, the compositenesses of the Gamow state lead to another result. In JüBo2022 the on-shell residues are $r_{\pi N} = (9.31 - 0.90i) \times 10^{-3} \text{ MeV}^{-1/2}$ and $r_{\eta N} = (0.25 - 3.97i) \times 10^{-3} \text{ MeV}^{-1/2}$; in contrast the off-shell residue of the ηN channel is larger than that of πN , so the result $X_{\eta N}$ is also large in Table IV. Consequently the elementariness is small, with the naive measure of only 8.5%. We should emphasize again that the naive measure is not mathematically a probability, it is understandable that the naive measure does not match the probability given by the spectral density function. Nevertheless, this may be an indication of the involved dynamics. Since the two results from the spectral density functions strongly suggest the elementary interpretation, we might claim the $N^*(1650)$ is possibly compact, but with visible model uncertainties. Note that the results from the spectral density functions are compatible with the conclusion in Ref. [11]: it is necessary to include an s -channel bare state to generate the $N^*(1650)$, but not necessary for $N^*(1535)$.

3. $N(1440) \frac{1}{2}^+$

The $N^*(1440)$ is very interesting in this model; it is definitely dynamically generated in the T^{NP} part [10]. However, the peak structure still shows up significantly in the first spectral density function; see Fig. 4. Meanwhile, the corresponding elementariness is moderate (49.5%). Note that

TABLE IV. The compositenesses and elementarinesses of the selected N^* Gamow states (two-body channels). The percentages in the brackets are the naive measures from Eq. (87).

State	$X_{\pi N}$	$X_{\eta N}$	$X_{K\Lambda}$	$X_{K\Sigma}$	Z
$N(1535) \frac{1}{2}^-$	$0.12 + 0.14i$ (7.5%)	$0.67 + 0.57i$ (35.8%)	$0.03 - 0.01i$ (1.3%)	$0.01 - 0.03i$ (1.2%)	$0.32 - 0.92i$ (39.4%)
$N(1650) \frac{1}{2}^-$	$0.17 + 0.28i$ (22.2%)	$0.60 - 0.15i$ (41.4%)	$0.03 + 0.03i$ (3.0%)	$0.15 - 0.03i$ (10.2%)	$0.12 + 0.02i$ (8.5%)
$N(1440) \frac{1}{2}^+$	$0.69 + 0.37i$ (59.0%)	$0.00 + 0.00i$ (0.2%)	$0.00 - 0.00i$ (0.0%)	$0.00 + 0.00i$ (0.0%)	$0.34 - 0.36i$ (36.9%)
$N(1710) \frac{1}{2}^+$	$0.04 - 0.00i$ (1.4%)	$0.87 + 0.97i$ (44.9%)	$-0.10 + 0.18i$ (6.9%)	$-0.00 + 0.01i$ (0.3%)	$0.02 - 1.17i$ (40.3%)
$N(1720) \frac{3}{2}^+$	$-1.16 - 3.01i$ (32.6%)	$0.04 - 0.04i$ (0.5%)	$0.03 + 0.05i$ (0.6%)	$-0.08 + 0.13i$ (1.5%)	$2.71 + 3.09i$ (41.4%)
$N(1900) \frac{3}{2}^+$	$-0.00 + 0.00i$ (0.1%)	$-0.02 + 0.00i$ (0.2%)	$0.39 + 0.33i$ (6.0%)	$0.90 - 0.17i$ (10.7%)	$-3.04 + 1.22i$ (38.5%)
$N(1520) \frac{3}{2}^-$	$0.19 + 0.39i$ (15.9%)	$-0.00 + 0.00i$ (0.0%)	$0.00 + 0.00i$ (0.0%)	$0.00 - 0.00i$ (0.0%)	$0.87 - 0.66i$ (40.4%)
$N(1675) \frac{5}{2}^-$	$0.02 + 0.17i$ (9.4%)	$0.02 - 0.06i$ (3.3%)	$-0.00 + 0.00i$ (0.0%)	$0.00 + 0.00i$ (0.0%)	$1.01 - 0.49i$ (61.8%)
$N(1680) \frac{5}{2}^+$	$0.22 + 0.50i$ (36.3%)	$-0.00 - 0.00i$ (0.0%)	$0.00 - 0.00i$ (0.0%)	$0.00 - 0.00i$ (0.0%)	$0.68 - 0.48i$ (55.0%)

the first bare state in this channel is just the bare nucleon. As already mentioned, the physical nucleon in this model carries approximately 54% composition of the bare nucleon. Here in the energy region of $N^*(1440)$, i.e., $z \in [M_R - \Gamma_R, M_R + \Gamma_R]$, the set of scattering states carries 34%. The result 49.5% is obtained by further corrections of the Breit-Wigner peak in Eq. (77), and the sum rule Eq. (75) is not violated.

The locally constructed spectral density function gives a smaller elementariness (31.5%). In Table IV, the complex compositeness of the πN channel is rather large, and the

naive measure of the elementariness is also small (36.9%). However, we should admit that which channel has the biggest composition is model dependent. It was indicated in Ref. [43] that the $N^*(1440)$ has a large component of σN and also in the JüBo model the σN channel plays an important role in the P_{11} partial wave [52]. All in all, the $N^*(1440)$ tends to be a molecular state here. Actually, in the studies of the quark model, the $N^*(1440)$ deviates from the simple picture of a qqq state. For example in Ref. [13] the $N^*(1440)$ is depicted by a core of three valence quarks complemented by a meson cloud, while in Ref. [14] the coupling of the $N^*(1440)$ to the πN

TABLE V. The compositenesses and elementarinesses of the selected N^* Gamow states (three-body channels). The percentages in the parentheses are the naive measures from Eq. (87). The meaning of the channel indices are $\rho N(1) \rightarrow |J-L| = \frac{1}{2}, S = \frac{1}{2}$; $\rho N(2) \rightarrow |J-L| = \frac{1}{2}, S = \frac{3}{2}$; $\rho N(3) \rightarrow |J-L| = \frac{3}{2}, S = \frac{3}{2}$; $\pi \Delta(1) \rightarrow |J-L| = \frac{1}{2}$; $\pi \Delta(2) \rightarrow |J-L| = \frac{3}{2}$.

State	$X_{\rho N(1)}$	$X_{\rho N(2)}$	$X_{\rho N(3)}$	$X_{\pi \Delta(1)}$	$X_{\pi \Delta(2)}$
$N(1535) \frac{1}{2}^-$	$-0.17 + 0.27i$ (13.0%)		$-0.00 + 0.00i$ (0.2%)		$0.02 - 0.04i$ (1.6%)
$N(1650) \frac{1}{2}^-$	$-0.02 - 0.11i$ (7.4%)		$-0.01 - 0.06i$ (4.3%)		$-0.04 + 0.02i$ (3.0%)
$N(1440) \frac{1}{2}^+$	$-0.00 + 0.01i$ (0.6%)	$-0.01 + 0.00i$ (0.6%)		$-0.02 - 0.03i$ (2.7%)	
$N(1710) \frac{1}{2}^+$	$0.00 - 0.00i$ (0.0%)	$-0.01 + 0.01i$ (0.3%)		$0.17 + 0.01i$ (5.9%)	
$N(1720) \frac{3}{2}^+$	$0.19 - 0.07i$ (2.1%)	$0.58 - 0.07i$ (5.9%)	$0.10 - 0.07i$ (1.2%)	$-1.40 + 0.01i$ (14.1%)	$-0.01 - 0.01i$ (0.1%)
$N(1900) \frac{3}{2}^+$	$0.01 - 0.01i$ (0.1%)	$3.04 - 1.60i$ (40.3%)	$-0.00 - 0.00i$ (0.0%)	$-0.27 + 0.22i$ (4.1%)	$-0.00 - 0.00i$ (0.0%)
$N(1520) \frac{3}{2}^-$	$0.00 + 0.00i$ (0.1%)	$-0.01 + 0.01i$ (0.5%)	$-0.58 + 0.26i$ (23.5%)	$0.01 - 0.00i$ (0.3%)	$0.52 - 0.00i$ (19.3%)
$N(1675) \frac{5}{2}^-$	$-0.00 - 0.00i$ (0.1%)	$-0.14 + 0.26i$ (16.2%)	$0.00 - 0.01i$ (0.4%)	$0.10 + 0.12i$ (8.8%)	$-0.00 + 0.00i$ (0.0%)
$N(1680) \frac{5}{2}^+$	$0.00 - 0.01i$ (0.4%)	$0.00 - 0.00i$ (0.2%)	$-0.01 - 0.00i$ (0.7%)	$0.00 + 0.00i$ (0.0%)	$0.11 - 0.01i$ (7.4%)

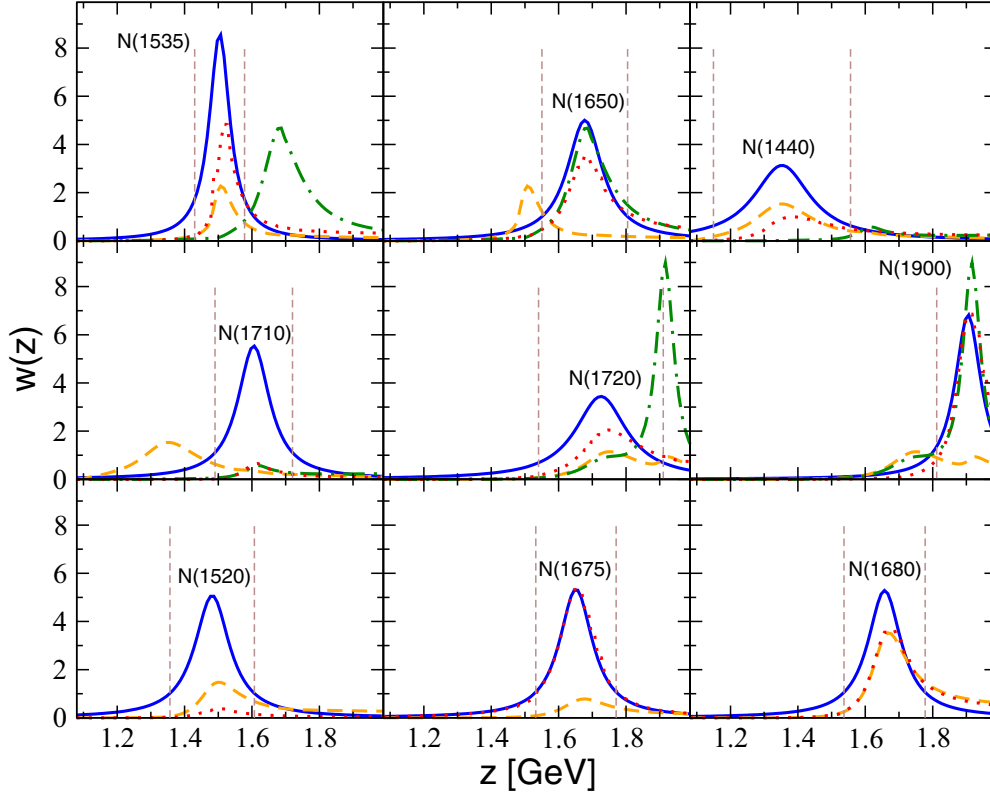


FIG. 4. The spectral density functions for all the selected N^* states. Blue solid line: the Breit-Wigner denominator in Eq. (57). Orange dashed (green dash-dotted) line: the first (second) spectral density function from the model. Red dotted line: the locally constructed function in Eq. (82) [for $N^*(1675)$ Eq. (83) is applied]. The vertical lines label the integral region of Eq. (77).

channel differs a lot from the prediction of the simple quark model. Note also that the Roper resonance is often considered as breathing mode of the nucleon; see e.g., Refs. [66,67] and references therein.⁵ We point out that the interplay of a bare Roper state with pion loops was already considered in Ref. [68] in the framework of chiral dynamics with explicit resonance fields.

4. $N(1710)_{\frac{1}{2}}^{+}$

The quantum numbers of the $N^*(1710)$ are the same as for the $N^*(1440)$, and their elementarinesses are both not large from the spectral density functions. As seen from Fig. 4, the first function from the model does not show a significant structure near its pole mass, while the second has a very weak structure. The local construction is very similar to the latter. The total elementariness of this state does not exceed 21%. The complex compositeness of the ηN channel is quite large. As given in Table IV, the naive measure of the ηN compositeness is 44.9%, whereas the elementariness is around 40.3%. So all the three results suggest the $N^*(1710)$ to be more of composite nature. This corroborates the findings of Ref. [45] that the $N^*(1710)$ is dynamically generated in the JüBo2022

analysis while the second bare s channel of the P_{11} wave pole moved far into the complex plane ($z_0 = 1513 - i405$ MeV).

Note that the pole of $N^*(1710)$ in the JüBo_omegaN solution [60] (fit A) is only 20 MeV below the ωN threshold, and the coupling is large. One may expect a quantitative change of the result for this state with a large component in the ωN channel. However, a systematic study can only be carried out after ωN photoproduction is fully analyzed.

5. $N(1720)_{\frac{3}{2}}^{+}$

In the P_{13} wave of the πN channel, two bare s -channel states are included and we observe two resonance poles, the $N^*(1720)$ and the $N^*(1900)$. The peak of $N^*(1900)$ is very strong, which affects the analyses of the $N^*(1720)$ since it is not very narrow; cf. Table III. Such interference severely increases the ambiguity. As seen from Table III, the elementariness from the first spectral density function is around 34.1%, while the second gives 68.5%, which mainly comes from the shoulder of the $N^*(1900)$ peak; see Fig. 4. Therefore an unambiguous interpretation of the numbers above is difficult. The locally constructed function gives a value of more than 60%, preferring the elementary nature. On the other hand, the off-shell residues of the πN and $\pi \Delta$ channels are extremely large, so the naive measure of the elementariness is less than 50%. In a word, we cannot draw any firm conclusion on this state.

⁵Specifically, the calculation of the decay width in Ref. [66] does not support the interpretation of the $N^*(1440)$ state as the radial excitation of the nucleon.

6. $N(1900) \frac{3}{2}^+$

The $N^*(1900)$ can hardly be seen in the purely hadronic observations, but is believed to be very important for KY photoproduction [69–73]. In the JüBo model, it is not significant if photoproduction is not considered. For instance, in the JüBo_omegaN solution [60], the pole lies in a rather distant region in the complex plane (the pole width $\Gamma_R > 600$ MeV) and the couplings are very weak. However, in the JüBo2022 solution [45] it becomes a significant narrow resonance.

We meet a situation similar to the $N^*(1650)$ for this state. The second spectral density function from the model in Fig. 4 simply suggests 100% elementariness for this state and also the local construction indicates nearly 100% elementariness. In contrast, the complex compositeness in Table V show a very large ρN composition of this state. Note that ρN is an effective three-body channel, for which the constraints from data are relatively weak.

7. $N(1520) \frac{3}{2}^-$

The D_{13} wave of πN is rather clear; there is only one s -channel bare state and only one resonance pole. The spectral density functions in Fig. 4, whether directly from the model or locally constructed, indicate a small elementariness of this state. Furthermore, in Table V, the compositions of $N^*(1520)$ in the three-body channels are relatively large. The model uncertainty for this state seems to be small. All this suggests the $N^*(1520)$ to be composite.

8. $N(1675) \frac{5}{2}^-$

As shown in Fig. 4, the spectral density function directly given by the model provides a very small elementariness for $N^*(1675)$. However, the local construction from Eq. (81) fails for this state. The residue of the $\rho N(2)$ channel, i.e., $S = 3/2$ and $|J - L| = 1/2$, is quite large, and the imaginary part of the $\rho N(2)$ loop function is positive in Eq. (81). We have checked the unitarity condition: every L_k'' must have a negative imaginary part when the energy is $z - i0^+$, with z real and higher than the corresponding threshold. Even though $\text{Im} L_{\rho N(2)}''(M_R - i0^+) < 0$, at the pole $\text{Im} L_{\rho N(2)}''(M_R - i\Gamma_R/2) > 0$. The $N^*(1675)$ is not broad, but the sign is still changed by the finite width. If one uses Eq. (83), the elementariness would be nearly 100%, which is not so trustworthy because the couplings are rather small, and the pole width almost totally comes from the constant width in the mass parameter M_0 . On the bright side, the constant width does not cause negative values near the pole mass; see Fig. 4. In addition, the violation of the sum rule is not large for this construction, $\int_{m_\pi+m_N}^\infty w^{\text{lc}}(z)dz \simeq 0.97$, deviating from the standard value of 1 only by 3%, and this spectral density function from plan B does not show any negative value in the energy region we study.

Furthermore, the complex compositenesses in Tables IV and V prefer a larger elementariness, which is opposite to the spectral density function from the model. So we cannot draw a certain conclusion on the $N^*(1675)$.

9. $N(1680) \frac{5}{2}^+$

Unlike the $N^*(1675)$, the uncertainties of the $N^*(1680)$ in this study are not large. The spectral density functions given by the model and the local construction agree with each other well in Fig. 4, both of which lead to an elementariness larger than 67%. The complex compositenesses in Tables IV and V may indicate non-negligible compositions of πN and $\pi \Delta$ channels, but the naive measure of the elementariness is still more than 55%. It is expected that the $N^*(1680)$ tends to be elementary.

Note again that the pole position of the $N^*(1680)$ is closer to the ωN threshold in the JüBo_omegaN solution [60], and the coupling is not small. There is a possibility that the $N^*(1680)$ becomes more composite when ωN is considered.

C. Analyses of the Δ states

The elementariness for the selected Δ states is summarized in Table VI. The complex quantities of the Gamow states are listed in Table VII. The corresponding spectral density functions are plotted in Fig. 5. The results are explained and discussed one by one in what follows.

1. $\Delta(1620) \frac{1}{2}^-$

There are significant uncertainties for this state. The spectral density function given by the model in Fig. 5 is very small, which indicates strongly that this state only very weakly related to the s -channel bare state. The local construction gives an exactly moderate elementariness 50.0%. However, the complex compositenesses in Table VII are small, and the naive measure of the elementariness is even 69.4%. Since the three scenarios lead to three different results, it is hard to draw a conclusion for this state. Note again that, although in JüBo2022 with the inclusion of the mixed-isospin $\gamma p \rightarrow K \Sigma$ reactions more experimental information on the Δ states was taken into account, the database for the Δ states is still smaller than that for the N^* states. Thus it is understandable that some outputs are uncertain.

2. $\Delta(1232) \frac{3}{2}^+$

As the lowest Δ state, there has been a lot of discussions on the $\Delta(1232)$ in the literature. In the quark model the $\Delta(1232)$ can be described by a qqq core, though with the Goldstone mesons as also the degrees of freedom [7]. Reference [14] also claims the deficiency of the simple quark model when studying the $\Delta(1232)$. As for the studies of the compositeness, there are different proposals, for example Ref. [74] finds a sizable πN component in the resonance, while Ref. [43] obtains a larger elementariness. Note that the interplay of the compact (bare) state with the pion loops around the nucleon was already discussed in Ref. [68] within baryon chiral perturbation theory with explicit resonance fields, and it was shown that both components are required to achieve a proper description of the P_{33} partial wave of pion-nucleon scattering.

In the JüBo model the situation is even more complicated, and unfortunately we cannot draw firm conclusions. The Δ particle appears as the ground state in the initial or final $\pi \Delta$ channel and also as s -channel intermediate states in various processes. In addition, there are some u -channel potentials

TABLE VI. The elementariness of the selected N^* states. The label “(F)” means that the local construction of Eq. (81) has failed and Eq. (83) is used instead.

State	Pole position (MeV)	$\mathcal{Z}_1$	$\mathcal{Z}_2$	$\mathcal{Z}_{tot}$	$\mathcal{Z}^{lc}$	$\tilde{\mathcal{Z}}$
$\Delta(1620) \frac{1}{2}^-$	$1607 - 42i$	18.9%		18.9%	50.0%	69.4%
$\Delta(1232) \frac{3}{2}^+$	$1215 - 46i$	53.4%	0.9%	53.8%	82.6% (F)	30.5%
$\Delta(1600) \frac{3}{2}^+$	$1590 - 68i$	39.7%	13.3%	47.8%	77.5%	69.7%
$\Delta(1700) \frac{3}{2}^-$	$1637 - 148i$	59.7%		59.7%	44.9%	47.8%

with Δ being exchanged. Those three Δ 's should be physically the same, but due to practical reasons there are technical simplifications. The initial/final state Δ only couples to the πN channel, and the coupling constant and bare mass are fixed; see Refs. [10,48]. The Δ pole in the amplitude is generated by the s -channel bare state, which couples to all the $I = 3/2$ channels, and the bare mass and coupling constants are fit parameters. Meanwhile, the u -channel exchanged Δ is just regarded as a stable particle with the mass 1232 MeV. Because of numerical limitations, it is at the moment impossible to overcome these inconsistencies.

Anyway, we may start the analyses with the initial/final state Δ . Its bare mass in this model is $M_0 = 1415$ MeV, with the coupling to the πN channel $g^2/(4\pi) = 0.36$. The resulting pole position in the propagator of Eq. (60) is $1211 - 37i$ MeV. This is very similar to the single channel toy model, and the spectral density function can be obtained directly from the imaginary part of Eq. (60); see Fig. 6. The resulting elementariness is 69.6%, suggesting the elementary interpretation. One can also calculate the complex elementariness of the Gamow state by directly applying Eq. (40): $Z = 0.64 - 0.32i$ with the compositeness (only of πN channel) $X = 1 - Z = 0.36 + 0.32i$. Then the naive measure of the elementariness is 59.5%, qualitatively in agreement with the spectral density function.

As for the pole of $\Delta(1232)$ in the amplitude, there are two s -channel bare states in the P_{33} wave, and the spectral density functions show a typical mixture. The two structures, the first of which is the $\Delta(1232)$, both show up in the first function. The elementariness directly from the model is 53.8%. The local construction of Eq. (81) fails. Similarly to the $N^*(1675)$ state discussed in the last section, here the $\pi \Delta$ residue is large

and the imaginary part of the $\pi \Delta$ loop function is positive. Then the whole imaginary part is positive, making the g_1^2 in Eq. (81) negative. If we use Eq. (83), the constructed spectral density function gives an elementariness about 82.6%. This “plan B” does not apparently violate the sum rule either: $\int_{m_\pi+m_N}^\infty w^{lc}(z)dz \simeq 0.93$. Besides, this spectral density function from plan B does not show any negative value in the energy region we study. The spectral density function may suggest the $\Delta(1232)$ to be more elementary than composite, but on the contrary the Gamow state has a large composition of the $\pi \Delta$ channel, leading to a rather small measure of the elementariness; see Table VII. We cannot claim this is certainly true because of the inconsistencies discussed above.

3. $\Delta(1600) \frac{3}{2}^+$

Though the $\Delta(1600)$ and the $\Delta(1232)$ are in the same wave, there is much less uncertainty for the $\Delta(1600)$. In Fig. 5, it is seen that the peak of $\Delta(1600)$ mainly stays on the first curve. The total elementariness from the two is rather moderate (47.8%), while the local construction succeeds and results in 77.5%. On the other hand, as shown in Table VII, the biggest partial compositeness is of the $\pi \Delta$ channel, while the naive measure of the elementariness is 69.7%. Combining the three results, the $\Delta(1600)$ is likely to be an elementary state. This is in accordance with the observation in Ref. [45] that the $\Delta(1600)$ is now largely induced by a bare s -channel pole, whereas it was dynamically generated in previous JüBo studies [53–55].

4. $\Delta(1700) \frac{3}{2}^-$

The result for the $\Delta(1700)$ in this study is interesting. The three scenarios agree with each other, but all give rather

TABLE VII. The compositeness and elementariness of the selected Δ Gamow states. The percentages in the parentheses are the naive measures from Eq. (87). The meaning of the channel indices are $\rho N(1) \rightarrow |J-L| = \frac{1}{2}, S = \frac{1}{2}; \rho N(2) \rightarrow |J-L| = \frac{1}{2}, S = \frac{3}{2}; \rho N(3) \rightarrow |J-L| = \frac{3}{2}, S = \frac{3}{2}; \pi \Delta(1) \rightarrow |J-L| = \frac{1}{2}; \pi \Delta(2) \rightarrow |J-L| = \frac{3}{2}$.

State	$X_{\pi N}$	$X_{\rho N(1)}$	$X_{\rho N(2)}$	$X_{\rho N(3)}$	$X_{\pi \Delta(1)}$	$X_{\pi \Delta(2)}$	$X_{K\Sigma}$	Z
$\Delta(1620) \frac{1}{2}^-$	$0.10 - 0.03i$ (8.6%)	$-0.02 - 0.09i$ (8.0%)		$-0.00 - 0.00i$ (0.3%)		$0.13 - 0.09i$ (13.0%)	$0.01 + 0.00i$ (0.7%)	$0.79 + 0.20i$ (69.4%)
$\Delta(1232) \frac{3}{2}^+$	$0.63 + 1.16i$ (28.5%)	$0.02 - 0.01i$ (0.4%)	$-0.00 + 0.01i$ (0.2%)	$0.06 - 0.03i$ (1.4%)	$1.54 - 0.58i$ (35.4%)	$-0.01 - 0.01i$ (0.3%)	$-0.00 + 0.15i$ (3.3%)	$-1.24 - 0.70i$ (30.5%)
$\Delta(1600) \frac{3}{2}^+$	$-0.04 + 0.07i$ (3.5%)	$0.00 - 0.01i$ (0.3%)	$0.01 - 0.03i$ (1.3%)	$-0.00 - 0.00i$ (0.1%)	$-0.42 + 0.21i$ (21.0%)	$-0.00 - 0.00i$ (0.0%)	$-0.08 + 0.05i$ (4.1%)	$1.53 - 0.29i$ (69.7%)
$\Delta(1700) \frac{3}{2}^-$	$-0.03 + 0.05i$ (2.3%)	$-0.00 + 0.01i$ (0.4%)	$-0.03 + 0.02i$ (1.3%)	$-0.03 + 0.00i$ (1.3%)	$-0.01 - 0.03i$ (1.5%)	$0.45 - 0.95i$ (45.4%)	$0.00 + 0.00i$ (0.0%)	$0.64 + 0.91i$ (47.8%)

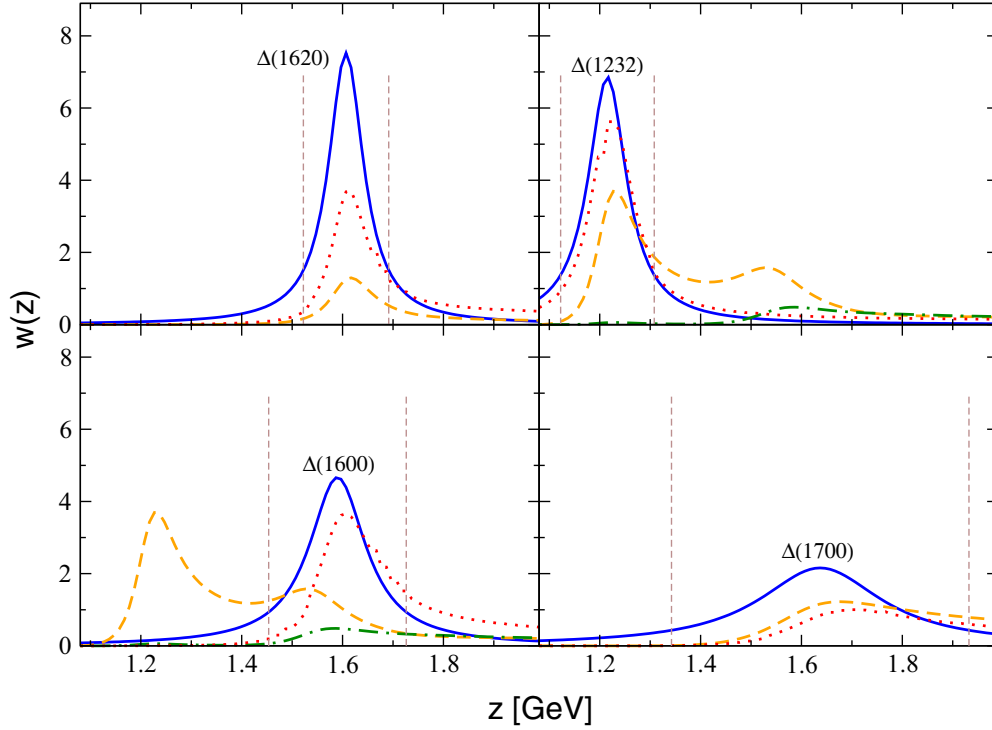


FIG. 5. The spectral density functions for all the selected Δ states. Blue solid line: the Breit-Wigner denominator in Eq. (57). Orange dashed (green dash-dotted) line: the first (second) spectral density function from the model. Red dotted line: the locally constructed function in Eq. (82) [for $\Delta(1232)$ Eq. (83) is applied]. The vertical lines label the integral region of Eq. (77).

moderate values. As seen from Fig. 5, the spectral density function from the model and the local construction behave similarly. The model gives the value of the elementariness as 59.7%, while the local construction results in is 44.9%. In Table VII, the naive measure of the elementariness is also moderate (47.8%), and the largest composition seems to be $\pi\Delta$. One might claim that physically this state is

really half elementary and half composite, but notice that the width of the $\Delta(1700)$ resonance (almost 300 MeV) is the biggest among all the selected states, causing visible ambiguities.

V. CONCLUSION AND OUTLOOK

This paper studies the composition of the N^* and Δ resonances, based on the coupled-channels dynamics of the Jülich-Bonn model. Results from three scenarios are compared to draw the conclusions: the spectral density functions directly from the model, the local construction from the pole parameters, and the complex compositenesses. The first two scenarios stem from the physical scattering states, which give probabilities as output, while the last is defined by the Gamow states and is different from the first two in the basic philosophy. The three results roughly depict the model dependence of this study. We have selected 13 states, 8 of which have reached relatively certain results: $N(1535)_{\frac{1}{2}}^{-}$, $N(1440)_{\frac{1}{2}}^{+}$, $N(1710)_{\frac{1}{2}}^{+}$, and $N(1520)_{\frac{3}{2}}^{-}$ have chances of being composite, whereas $N(1650)_{\frac{1}{2}}^{-}$, $N(1900)_{\frac{3}{2}}^{+}$, $N(1680)_{\frac{5}{2}}^{+}$, and $\Delta(1600)_{\frac{3}{2}}^{+}$ have tendencies to be elementary. For those 8 states, at least two of the three scenarios result in qualitative agreements.

However, model uncertainties do exist. Some of the them are caused by the three-body effective channels; especially the uncertainty of the $\Delta(1232)$ state is highly related to the $\pi\Delta$ channel. Including rigorous three-body unitarity and crossing symmetry is a challenge left for future studies. Moreover,

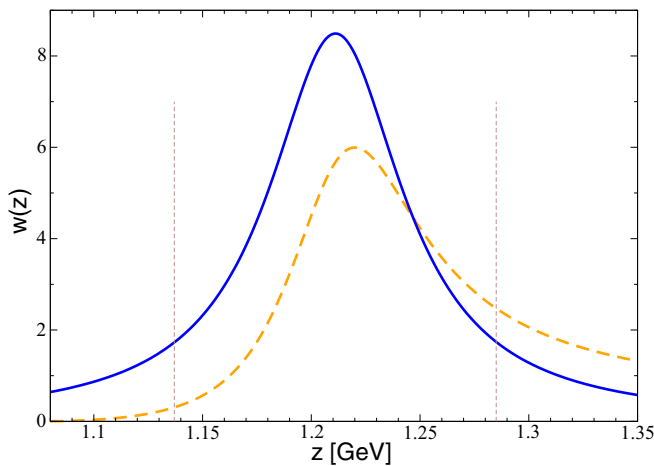


FIG. 6. The spectral density functions the initial/final state $\Delta(1232)$. Blue solid line: the Breit-Wigner denominator in Eq. (57). Orange dashed line: the spectral density function from the model. The vertical lines label the integral region of Eq. (77).

theoretically, the connections between the spectral density functions and the Gamow states still need to be understood in a deeper way. Note that for $N(1650)_{\frac{1}{2}}^{-}$, $N(1900)_{\frac{3}{2}}^{+}$, and $\Delta(1232)_{\frac{3}{2}}^{+}$ the complex compositenesses of the Gamow states lead to opposite conclusions on the spectral density functions. Moreover, relating the results which are based on a pure hadronic model to the descriptions in quark models appears to be difficult. However, it is clear that a compact state in such an approach must have a mapping onto a quark model state, which we have assumed throughout. All in all, this topic itself is difficult in nature. One cannot expect very clear interpretations and definite conclusions for every state, since only the states like the deuteron can be highly related to physical observables. We believe this paper presents the presently achievable status of such type of investigation. We also hope this paper can inspire the community toward further constructive discussions on this topic.

In the future, the ωN channel can be included in the study of the compositions, after the ωN photoproduction is included in this model. Moreover, including more high-quality data also from other photoproduction channels will lead to refined pole values. The same methods can be applied on the other sectors, e.g., the Λ^* resonances or the P_c

exotic states, which are also being studied in the Jülich-Bonn model.

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