PHYSICAL REVIEW C **81**, 051301(R) (2010)

Narrow states in the three-proton emitter 17Na

N. K. Timofeyuk^{1,*} and P. Descouvemont²

¹*Department of Physics, Faculty of Electronics and Physical Sciences, University of Surrey, Guildford GU2 7XH, United Kingdom* ²*Physique Nucleaire Th ´ eorique et Physique Math ´ ematique, Case Postale 229, Universit ´ e Libre de Bruxelles, B-1050 Brussels, Belgium ´* (Received 5 February 2010; published 19 May 2010)

Based on a microscopic cluster model that reproduces well the ^{17}C spectrum below the neutron threshold, we study the spectrum of its mirror nucleus 17 Na. We find that the 17 Na ground state should be located at 2.4 MeV above the ¹⁶Ne(0_1^+) + *p* threshold, being a broad $1/2^+$ resonance with a width of 1.4 MeV. However, there should exist at least four narrow excited states in ¹⁷Na; three of them, $3/2^+_1$, $7/2^+_1$, and $9/2^+_1$, decay into the ¹⁶Ne(2^+_1) + *p* channel, and the other, $5/2^+_1$, decays mainly into the ¹⁶Ne(0^+_1) + *p* channel. Because the daughter nucleus ¹⁶Ne is a two-proton emitter, the narrow ¹⁷Na states must undergo a sequential three-proton decay via intermediate states in 16 Ne.

DOI: [10.1103/PhysRevC.81.051301](http://dx.doi.org/10.1103/PhysRevC.81.051301) PACS number(s): 21*.*60*.*Gx, 23*.*50*.*+z, 27*.*20*.*+n

Introduction. The development of experimental techniques that allow the fragments of in-flight-decay to be tracked and identified makes it possible to study nuclear structure beyond the proton drip line, where nuclei exist only as resonances in the continuum. With these techniques, one-proton (1*p*) and two-proton $(2p)$ emitters have been identified $[1-5]$. In the medium and heavy mass regions, the lifetime of 1*p* and 2*p* emitters is long, compared to typical nuclear lifetimes, because the strong Coulomb barrier prevents their decay. In the light mass region, however, the Coulomb barrier is lower, the lifetimes are shorter, and, as a consequence, the resonance widths are larger. For example, the ground states of 15F or ¹⁰*,*11N are observed in reaction cross sections as broad *s*-wave resonances. It has been predicted, however, that, even in this region, proton emission from some excited states may be suppressed by structural reasons and that such states can be narrow. For example, three negative-parity states in 15F, 1*/*2−, 5*/*2−, and 3*/*2−, located between 4.5 and 7.6 MeV, are predicted to have a width of only a few kilo–electron volts [\[6,7\]](#page-4-0). Two of these states have recently been identified using in-flight decay tracking techniques [\[1\]](#page-4-0). Their estimated widths are less than $400 \,\text{keV}$. In addition, a state in 15 F with a similarly small width has been seen at 7.6 MeV. It was pointed out in Ref. [\[1\]](#page-4-0) that all these states have a peculiar cluster structure based on excited core states that are in turn 1*p* emitters.

In this paper, we show that narrow states can also exist in the spectrum of another proton-rich nucleus, 17 Na, which is the mirror analog of ${}^{17}C$ famous for its peculiar structure. The neutron binding energy in the ground state ${}^{17}C(3/2^+)$ is only 728 keV, typical of halo nuclei. However, knockout and Coulomb breakup experiments have shown that the weakly bound ${}^{16}C(0^+) + n$ configuration is suppressed in ${}^{17}C(3/2^+)$ and that this state is mainly based on the $^{16}C(2^+) + n$ configuration [\[8,9\]](#page-4-0), where the neutron binding energy is 2.5 MeV. A similar structure should be expected in the mirror nucleus 17 Na(3/2⁺). Therefore, the decay branch $^{17}Na(3/2^+) \rightarrow ^{16}Ne(0^+) + p$ could be suppressed,

and if energetically allowed, the main decay mode would be $17\text{Na}(3/2^+) \rightarrow 16\text{Ne}(2^+) + p$. If its decay energy is below the Coulomb barrier, then its width may be small. Because the decay product 16 Ne is unstable with respect to 2*p* emission, ¹⁷Na should be a three-proton $(3p)$ emitter.

At present, nothing is known about 17 Na. The only theoretical calculation, performed with a deformed Hartree-Fock model, suggests that the last proton orbital may be bound [\[10\]](#page-4-0). However, this is unlikely because 19 Na is unbound [\[11\]](#page-4-0) and 17 Na is obtained from 19 Na by further removal of two deeply bound neutrons. To study 17 Na, we use a two-center microscopic cluster model (MCM) in which excited states of the 16 Ne core are included. First, we show that this model solves the long-standing problem of reproducing the spectrum of its mirror analog 17 C below the neutron decay threshold. Then we use the same model to predict positions and decay widths of the 17 Na states.

¹⁷C *properties in the MCM*. Experimental spectrum of $17C$ and previous theoretical calculations: Experimentally, three levels are known below the neutron threshold in ${}^{17}C$ (see Fig. [1\)](#page-1-0). Their spin parities have been established from the analysis of the *γ*-ray spectrum observed in the $p + {}^{17}C$ inelastic scattering [\[12\]](#page-4-0) and from the study of one-neutron removal reaction from 18 C [\[13\]](#page-4-0). Unbound excited states have been observed in the β decay of ¹⁷B [\[14\]](#page-4-0) and in the three-neutron transfer ${}^{14}C(^{12}C, {}^{9}C)^{17}C$ reaction [\[15\]](#page-4-0). Three unbound excited states have also been recently identified at $E_x = 2.20(3), 3.05(3),$ and 6.13(9) MeV from inelastic proton scattering by detecting neutrons in coincidence with ¹⁶C [\[16\]](#page-4-0) emitted in the ¹⁷C^{*} \rightarrow ¹⁶C + *n* decay. It was suggested in Ref. $[16]$ that these levels have spin parities of $7/2^+$, $9/2^+$, and $5/2^+$ respectively. For the last two levels, this assignment agrees with the one based on the three-nucleon transfer study. However, the spin-parity assignment of 7*/*2⁺ for $E_r = 2.20(3)$ MeV contradicts the observed width of this state equal to $0.53(4)$ MeV. At this energy, the $l = 2$ decay channel ${}^{16}C(2^+) + n$ is closed, while the $l = 4$ decay into the $^{16}C(0^+) + n$ channel should have a width smaller than 1 keV. Also, the state at 2.20(3) MeV identified in the ¹⁷C(p, p)¹⁷C reaction coincides in energy with the state at 2.25(2) MeV seen in the β decay of ¹⁷B. There exist strong arguments for ¹⁷B to

^{*}n.timofeyuk@surrey.ac.uk

FIG. 1. 17C spectra calculated in the MCM with the V2 and MN interactions in comparison to the experimental spectrum (exp) and to the shell model (SM) WBP predictions. Labels correspond to 2*J* . The spin-parity assignment for observed unbound levels corresponds to that suggested by 3*n* transfer [\[15\]](#page-4-0).

have the spin parity of $3/2^-$, as it is obtained from ¹³B(3/2⁻) by adding four neutrons, coupled to spin state 0+, into the *sd* shell. Because no $7/2^+$ states could be populated in the ^{17}B $β$ decay, the spin-parity assignment $7/2^+$ suggested for the $2.20(3)$ MeV state in Ref. [\[16\]](#page-4-0) is most likely erroneous, even though the ¹⁷C(*p*, *p*')¹⁷C^{*} cross sections for this state can be explained by the $L = 2$ excitation to the $7/2^+$ state in the distorted-wave Born approximation calculations [\[16,17\]](#page-4-0). The 2.20(3) MeV state is more likely either $3/2^+_2$ or $5/2^+_2$.

The *spsdpf* shell model with the WBT interaction cannot reproduce the excitation energies and ordering of the two first excited states (see Fig. 1). The modification of this interaction made in Ref. [\[18\]](#page-4-0) does not improve the situation. The bound $17¹⁷C$ spectrum cannot be understood in the two-body potential model with deformation and the 2^+ excitation of the ¹⁶C core either, if standard sets of potentials are used [\[19,20\]](#page-4-0). An *l*-dependent $n + {}^{16}C$ potential and a nonstandard spin-orbit interaction should be used for these purposes, but these lead to difficulties in explaining the spectrum of ${}^{18}C$ [\[20\]](#page-4-0). Also, no success has been made in explaining the ${}^{17}C$ spectrum in the Multi-Channel Algebraic Scattering theory [\[17\]](#page-4-0), in which only the first 2^+ state of the ¹⁶C core was taken into account. This suggests that ${}^{16}C(4^+)$ excitations can be important in ¹⁷C, consistent with a strong population of the ¹⁶C excitation energies around 4 MeV in nucleon knockout from ^{17}C [\[8\]](#page-4-0), where the 2^+_2 , 3^+ , and 4^+ states have been seen.

In Ref. [\[21\]](#page-4-0), an attempt was made to understand the ^{17}C spectrum within a two-cluster MCM. In this model, the valence neutrons in the ¹⁶C core were allowed to occupy only the $0d_{5/2}$ subshell. This gives $0^+, 2^+,$ and 4^+ excitations of the core, all of which were included in the calculations. This model was able to reproduce the positions and separation energies of the $3/2^+$ and $1/2^+$ states by tuning the strength of the spin-orbit force. However, the $5/2^+$ state was still unbound by about 300 keV. This can be a consequence of excluding the 1*s*1*/*² and $0d_{3/2}$ orbitals from the model space used for the ¹⁶C core.

New MCM study of ${}^{17}C$: Here we extend the multichannel MCM study of 17C performed in Ref. [\[21\]](#page-4-0). In a microscopic theory, the Hamiltonian of the system reads

$$
H = \sum_{i=1}^{A} T_i + \sum_{j>i=1}^{A} V_{ij},
$$
 (1)

where T_i is the kinetic energy of nucleon *i*, and V_{ij} a nucleon-nucleon interaction. For large nucleon numbers, the Schrödinger equation associated with this Hamiltonian cannot be solved exactly. We used here the cluster approximation, where the wave function in partial wave J^{π} with channel spin *I* reads

$$
\Psi^{JM\pi} = \sum_{l_l} A \left[Y_l(\Omega_\rho) \otimes \left[\phi_C^{l_c} \otimes \phi_n \right]^l \right]^{JM} g_{l_l}^{J\pi}(\rho). \quad (2)
$$

Here I_C is the spin of ¹⁶C and *l* is the relative angular momentum, taken as 0, 2, and 4. In Eq. (2), $\phi_C^{I_C}$ and ϕ_n are the internal wave functions of the clusters, $g_{I_cII}^{J\pi}(\rho)$ is the radial function depending on the relative coordinate *ρ*, and A is the antisymmetrizer that permutes the last nucleon with the nucleon of the core. A cluster model is well adapted to exotic weakly bound (or unbound) nuclei, as the asymptotic behavior at large *ρ*, crucial for them, is exactly taken into account through the *R*-matrix method. In this method, the total width Γ of unbound states is determined as the imaginary part of a complex eigenvalue (see Ref. [\[22\]](#page-4-0) for details). The partial widths Γ_l in each channel are obtained from the associated eigenvectors and provide the reduced widths γ_l^2 as

$$
\Gamma_l = 2ka\gamma_l^2/|O_l|^2,\tag{3}
$$

where *k* is the wave number, *a* is the channel radius, and $O_l(ka)$ is an outgoing Coulomb function. The reduced widths γ_l^2 do not depend on the energy of the resonance and are proportional to the wave function at $r = a$.

We use two effective nucleon-nucleon (*NN*) forces, Volkov V2 [\[23\]](#page-4-0) and Minnesota (MN) [\[24\]](#page-4-0), complemented by a zerorange spin-orbit force [\[25\]](#page-4-0). Both forces have one adjustable parameter (*m* and *u*, respectively) that gives the strength of the *NN* potentials in odd partial waves.

Unlike in Ref. [\[21\]](#page-4-0), the two ¹⁶C valence neutrons occupy the $0d_{5/2}$, $1s_{1/2}$, and $0d_{3/2}$ orbitals, which gives many excitations in ¹⁶C. We consider only the $2^+_{1,2}$, 3^+_{1} , and 4^+_{1} states, motivated by neutron knockout experiments where they are strongly populated. The single-particle wave functions in 16C were taken from the oscillator shell model with the oscillator radius of 1.6 fm.

The theoretical and experimental spectra of ${}^{16}C$ are shown in Fig. [2.](#page-2-0) The V2 potential gives a correct excitation energy for ${}^{16}C(2^+)}$, which is very important for predicting the widths of the 17 Na decay modes into two different channels. With MN, the excitation energy of the $2₁⁺$ state is only 1.39 MeV, which is about 300 keV lower than the experimental value of 1.766 MeV.

Let us now discuss the MCM 17 C spectrum shown in Fig. 1 and compared to experiment. The parameter *m* (or u) and the spin-orbit amplitude have been chosen to reproduce both the ${}^{16}C + n$ threshold in ¹⁷C and the excitation energy of ¹⁷C(1/2⁺). These values, $m = 0.64922$

FIG. 2. Theoretical and experimental ¹⁶C and ¹⁶Ne spectra.

and $S_0 = 31.28 \text{ MeV}$. fm⁵ for V2 ($u = 0.76706$ and $S_0 =$ 38*.*95 MeV*.*fm5 for MN), are in the range of typical values used previously in the MCM for other cluster systems. With these parameters, the energies of the three low-lying states are in very good agreement with experiment. In particular, the level ordering is correctly reproduced by the MCM. For comparison, the shell model spectrum calculated with the WBP interaction [\[26\]](#page-4-0) using the NuShell code [\[27\]](#page-4-0) is also shown in Fig. [1.](#page-1-0)

To understand the structure of ${}^{17}C$ bound states in terms of core excitations, we have calculated the spectroscopic factors for the $\langle {}^{17}C \rangle$ ¹⁶C \rangle overlap in MCM. We list them in Table I in comparison to shell model predictions. We confirm that the ground state ${}^{17}C(3/2^+)$ is based on the ${}^{16}C(2_1^+) + n$ configuration, which has a spectroscopic factor similar to that obtained in the shell model. We also confirm the shell model results for the strong 2^+_2 and 4^+ core excitations in ¹⁷C(3/2⁺), but we do not get significant strength for ${}^{16}C(3^+_1) + n$.

The first $1/2_1^+$ excited state is mainly based on the $^{16}C(0^+) + n$ and $^{16}C(2_2^+) + n$ configurations. Unlike in the shell model, the contributions from the 2^+_1 and 3^+_1 states are small. As for the second excited state, ${}^{17}C(5/2_1^+)$, its structure

in terms of core excitations is similar to the shell model one, except for the *s*-wave ${}^{16}C(3^+) + n$ and *d*-wave ${}^{16}C(2_2^+) + n$ configurations, where the spectroscopic factors are close to zero.

The MCM predicts that the lowest unbound $3/2^+_2$, $5/2^+_2$, $7/2_1^+$, and $9/2_1^+$ levels form a group in a narrow region around 2 MeV above the ${}^{16}C + n$ threshold (see Fig. [1\)](#page-1-0). In the shell model these levels are also around 2 MeV but with a larger spread. For V2, the $7/2^+_1$ and $9/2^+_1$ resonances have a width of the order of 10^{-12} and 10^{-6} MeV, respectively, as their decay to the only open $l = 4$ channel ${}^{16}C(0_1^+) + n$, which is suppressed by the strong centrifugal barrier. The $5/2^+_2$ state is below the ${}^{16}C(2_1^+) + n$ threshold and has a width of 15 keV, while the $3/2^+_2$ state is 80 keV above the ¹⁶C(2⁺) + *n* threshold and has the width of 265 keV owing to the presence of the *s*-wave component in this channel. For MN, all these levels have widths of less than 20 keV.

The first four unbound levels in the experimental spectrum are indeed concentrated around 2 MeV above the threshold and they should probably correspond to the theoretical levels $3/2_2^+$, $5/2_2^+$, $7/2_1^+$, and $9/2_1^+$. Because the states at 2.06 and 3.10 MeV belonging to this group have not been seen in the β decay of ¹⁷B, they probably have spin parities of $7/2_1^+$ and $9/2_1^+$. The spin assignment of $9/2_1^+$ for the state at $3.05(3)$ MeV agrees well with both the (p, p') and the 3*n* transfer experiments. The state at 2.06 MeV could be $7/2_1^+$ according to the 3*n* transfer study. Therefore, the two remaining states, at 2.25 and 2.64 MeV, should be the $3/2^+_2$ and $5/2^+_2$ states. It is very important to reproduce the positions of the 17° C states for the 17° Na study. For these purposes, we refitted the parameter *m* of V2 individually for each of these states to reproduce their positions precisely. In these new calculations, the $7/2_1^+$ has a width of 30 meV, as it is still below the ¹⁶C(2⁺) + *n* threshold, while the 9/2⁺ state, which is now above this threshold, has a width of 18 keV, which is smaller than the width of 0.10(5) MeV observed in the 3*n* transfer. The 2.25-MeV state has a width of either 8 or 40 keV, depending on whether it has spin parity $5/2^+_2$ or $3/2^+_2$. In both cases it disagrees with the width of 0.53(4) MeV deduced from inelastic proton scattering. The 2.64-MeV state has a width of 368 or 595 keV for the $3/2^+_2$ or $5/2^+_2$ prescription, respectively. No experimental width for this state is available.

TABLE I. Spectroscopic factors $S_l = S_{l,j=l-1/2} + S_{l,j=l+1/2}$ for the $\langle {}^{17}C | {}^{16}C \rangle$ overlap calculated in the MCM with the V2 and MN potentials in comparison with the shell model (SM) values.

$I_C({}^{16}C)$		${}^{17}C(3/2^+)$			${}^{17}C(1/2^+)$			${}^{17}C(5/2^+)$		
		V2	MN	SM	V ₂	MN	SM	V ₂	MN	SM
0^{+}_{1}	$\boldsymbol{0}$				0.828	0.841	0.644			
	2	0.010	0.010	0.035				0.558	0.537	0.701
2^{+}_{1}	$\overline{0}$	0.328	0.383	0.163				0.037	0.066	0.096
	2	1.260	1.243	1.445	0.034	0.016	0.415	0.520	0.593	0.226
2^{+}_{2}	$\overline{0}$	0.030	0.033	0.225				0.050	0.030	0.014
	2	0.127	0.152	0.090	0.366	0.408	0.372	θ	0.002	0.631
4^{+}_{1}	2	0.372	0.389	0.381				0.969	0.965	0.916
3^{+}_{1}	$\boldsymbol{0}$							Ω	θ	0.301
	2	0.026	0.002	0.285	0.091	0.254	1.027	0.060	0.128	0.003

N. K. TIMOFEYUK AND P. DESCOUVEMONT PHYSICAL REVIEW C **81**, 051301(R) (2010)

TABLE II. Spins, energies, and widths in various ${}^{16}Ne + p$ channels of 17Na states with the V2 interaction. All energies and widths are given as mega–electron volts.

J^{π}	E	E_{x}	$\Gamma(0^+_1)$	$\Gamma(2^+_1)$
$1/2_1^+$	2.40	θ	1.36	
$3/2^+$	2.57	0.17	0.001	0.024
$5/2^+$	2.97	0.57	0.123	0.021
$7/2^+$	4.35	1.95	8×10^{-8}	0.025
$(5/2^+_2)^a$	4.38	1.98	0.032	1.561
$(3/2^+_2)^a$	5.27	2.87	0.146	1.237
$(3/2^+_2)^b$	4.63	2.23	0.108	0.828
$(5/2^+_2)^b$	5.32	2.92	0.059	2.509
$9/2^+$	5.41	3.01	6×10^{-6}	0.210

^a Assuming that the ¹⁷C states at 2.25 and 2.64 MeV are $5/2^+_2$ and $3/2^+_2$, respectively.

 $3/2_2^+$, respectively.
^bAssuming that the ¹⁷C states at 2.25 and 2.64 MeV are $3/2_2^+$ and $5/2^+_2$, respectively.

All levels above 2 MeV, calculated for spin less than 9*/*2+, have a width much larger than 1 MeV for both V2 and MN. No level with a width typical of those observed in the ¹⁴C(¹²C,⁹C)¹⁷C reaction, 0.10(5) $\leq \Gamma \leq 0.66(20)$ MeV, is predicted in this region. In particular, we find for the 5*/*2⁺ 4 level a width of 2.5 MeV, which is mainly caused by the *s*-wave decay into the ${}^{16}C(2_1^+) + n$ channel. This spin parity has been assigned to the level 6.2 MeV both in 3*n* transfer and in proton inelastic scattering. However, the observed widths of this state, $0.35(15)$ MeV from 3*n* transfer and $0.26^{+0.4}_{-0.26}$ MeV from (p, p') , are much smaller than the theoretical predictions.

The MCM spectrum of 17Na *and its decay properties*. Based on MCM with the V2 potential that reproduces the positions of the 17 C bound states, we have calculated the energies and widths of the $1/2_1^+$, $3/2_1^+$, and $5/2_1^+$ states in ¹⁷Na. They are reported in Table II. These states lie above both the ¹⁶Ne(0_1^+) + p and $^{16}Ne(2^+_1) + p$ decay thresholds, so we show the partial widths $\Gamma(0_1^+)$ and $\Gamma(2_1^+)$ for decay to these individual channels as well. The $1/2_1^+$ state becomes the ground state. It mainly decays to the *s*-wave ${}^{16}Ne(0_1^+) + p$ channel and has the large width of 1.36 MeV. The $3/2₁⁺$ state is located just above $1/2₁⁺$ but it has a different decay mode, ${}^{16}Ne(2_1^+) + p$, with the small width of only 24 keV. The $5/2₁⁺$ state is also narrow, with $\Gamma(0_1^+) = 125$ and $\Gamma(2_1^+) = 20 \,\text{keV}$. Because the MCM overpredicts the location of the $5/2₁⁺$ state by 46 keV, we can expect that the energy and the width of the $5/2₁⁺$ in ¹⁷Na are also slightly overestimated.

For the analogs of the first four unbound states in ${}^{17}C$ we used the modified V2 potentials discussed in the previous section. The decay scheme of the lowest part of the 17 Na spectrum is shown in Fig. 3. We find that the ${}^{17}Na(7/2_1^+)$ level should be very narrow. It decays into the *d*-wave ${}^{16}Ne(2^+_1) + p$ and *s*-wave ${}^{16}Ne(4_1^+) + p$ channels with partial widths of $\Gamma(2_1^+) = 25 \,\text{keV}$ and $\Gamma(4_1^+) = 98 \,\text{keV}$, respectively. It should be noted that the theoretical value of the latter threshold is underestimated by 630 keV, and therefore, the energy in this

FIG. 3. ¹⁷Na decay scheme with the V2 interaction.

channel is too high. Decreasing this energy by tuning the Majorana parameter *m*, we obtain a partial width $\Gamma(4_1^+) =$ 4 keV . Thus, the ¹⁷Na(7/2⁺) state should be as narrow as ¹⁷Na(3/2⁺₁). A similar situation occurs for the ¹⁷Na(9/2⁺₁) state. The partial width for the decay into ${}^{16}Ne(2_1^+) + p$ is predicted to be 211 keV. A similar width is expected for the decay into the ¹⁶Ne(4_1^+) + *p* channel. Tuning the energy of this channel to reproduce the position of the ${}^{17}C(9/2_1^+)$ state with respect to the ¹⁶Ne(4^+_1) + *p* threshold, similarly to what has been done in the case of $7/2₁⁺$, we get a partial width $\Gamma(4_1^+) = 78 \,\text{keV}.$

For the analogs of the $5/2^+_2$ and $3/2^+_2$ states in ¹⁷C, we made two predictions on the assumption that (a) the 2.25- and 2.64-MeV states in ¹⁷C are $5/2^+_2$ and $3/2^+_2$, and (b) the 2.25and 2.64-MeV states in ¹⁷C are $3/2^+_2$ and $5/2^+_2$. In both cases these states are broad and decay into the *s*-wave ¹⁶Ne(2_1^+) + *p* channel with a width between 0.8 and 2.5 MeV. As for all other excited states, the MCM predicts widths of 2 MeV and higher. Unless our model strongly overestimates the contribution from the ¹⁶Ne(0_1^+) and ¹⁶Ne(2_1^+) states (which is possible, as it gives overestimated widths for excited states above 3 MeV in the mirror nucleus ${}^{17}C$), no narrow states can be expected in this area.

Conclusions. Based on the MCM with the V2 potential that reproduces the 17 C bound spectrum, we predict that the ¹⁷Na ground state should be a broad $1/2^+$, $l = 0$ resonance. However, there should be at least four narrow states, $3/2₁⁺$, $5/2_1^+$, $7/2_1^+$, and $9/2_1^+$, in the ¹⁷Na spectrum. The decay product of these states, ¹⁶Ne, is unstable with respect to the two-proton emission. Therefore, 17 Na is in fact a three-proton emitter with a decay path ¹⁷Na \rightarrow ¹⁶Ne^{*} + *p* \rightarrow ¹⁴O + 2*p* + p . Consequently, 17 Na states can be identified by detecting $14O + p + p + p$ events in coincidence.

The narrow 17 Na states can be populated using the $14O(^{12}C, ^9Li)^{17}Na$ reaction in inverse kinematics with the $14O$ radioactive beam. It is a mirror analog of the 3*n* transfer ¹⁴C(${}^{12}C, {}^{9}C$)¹⁷C used in Ref. [\[15\]](#page-4-0) to study the unbound spectrum of ¹⁷C. This reaction strongly populates the ¹⁷C(5/2⁺₁) and ${}^{17}C(9/2_1^+)$ states at 0.31 and 3.10 MeV, the proton-rich analogs of which should be narrow. It also populates the

NARROW STATES IN THE THREE-PROTON EMITTER 17Na PHYSICAL REVIEW C **81**, 051301(R) (2010)

state at 2.06 MeV, which we think should be assigned to the $7/2₁⁺$ state. Its mirror analog should also be narrow. No states have been observed in the ${}^{14}C(^{12}C, {}^{9}C)^{17}C$ reaction at 0.21, 2.25, or 2.64 MeV. Therefore, we can expect that their mirror analogs (which are broad) will not be populated in the $^{14}O(^{12}C, ^{9}Li)^{17}Na$ reaction either, so that three distinctive narrow peaks would be seen in the 17 Na spectrum populated by the 3*p* transfer. The widths of these peaks can be studied both by analysis of the missing mass spectra and by detection in coincidence of their decay products. The 17 Na spectrum can be also studied in the charge exchange reaction $(^{17}Ne,^{17}Na)$. The Borromean nucleus 17 Ne has two protons orbiting the 15 O core. The charge exchange in this core would create the proton unstable nucleus ${}^{15}F = {}^{14}O + p$. With two additional protons around it, one gets the $3p$ emitter 17 Na. The knowledge of its energy levels and widths will test the accuracy of available theories beyond the proton drip line.

Finally, we have made predictions only for positive-parity states. No negative-parity states have yet been seen in ${}^{17}C$. In other $A = 17$ nuclei, states of the opposite to the groundstate parity are located near 3 MeV in ^{17}O and ^{17}F and near 2 MeV in 17 N. From the shell model point of view, they are intruders and their description requires taking into account many major shells. In the $17\text{Na} = 16\text{Ne} + p$ system, intruder states can be expected at proton energies $E_p > 4$ MeV, assuming that they are 2 MeV above the 17 Na ground state. Although such energies are above the Coulomb barrier, these states can be narrow because of small spectroscopic factors typical of intruder states. To predict reliably the positions and the spectroscopic factors of such states in the MCM is a challenging task for the future.

N.K.T. thanks I. Mukha for valuable discussions and acknowledges the UK STFC ST*/*F012012*/*1 grant.

- [1] I. Mukha *et al.*, Phys. Rev. C **79**[, 061301\(R\) \(2009\)](http://dx.doi.org/10.1103/PhysRevC.79.061301)
- [2] M. Pfützner et al., [Eur. Phys. J. A](http://dx.doi.org/10.1140/epja/i2002-10033-9) 14, 279 (2002); J. Giovinazzo *et al.*, Phys. Rev. Lett. **89**[, 102501 \(2002\).](http://dx.doi.org/10.1103/PhysRevLett.89.102501)
- [3] B. Blank *et al.*, Phys. Rev. Lett. **94**[, 232501 \(2005\).](http://dx.doi.org/10.1103/PhysRevLett.94.232501)
- [4] I. Mukha *et al.*, Phys. Rev. Lett. **99**[, 182501 \(2007\).](http://dx.doi.org/10.1103/PhysRevLett.99.182501)
- [5] I. Mukha *et al.*, [Nature \(London\)](http://dx.doi.org/10.1038/nature04453) **439**, 298 (2006).
- [6] L. Canton, G. Pisent, J. P. Svenne, K. Amos, and S. Karataglidis, Phys. Rev. Lett. **96**[, 072502 \(2006\).](http://dx.doi.org/10.1103/PhysRevLett.96.072502)
- [7] H. T. Fortune and R. Sherr, Phys. Rev. Lett. **99**[, 089201 \(2007\).](http://dx.doi.org/10.1103/PhysRevLett.99.089201)
- [8] V. Maddalena *et al.*, Phys. Rev. C **63**[, 024613\(R\) \(2001\).](http://dx.doi.org/10.1103/PhysRevC.63.024613)
- [9] U. Datta Pramanik *et al.*, [Eur. Phys. J. A](http://dx.doi.org/10.1140/epjad/i2005-06-173-2) **25**, 339 (2005).
- [10] H. Kitagawa, N. Tajima, and H. Sagawa, [Z. Phys. A](http://dx.doi.org/10.1007/s002180050345) **358**, 381 [\(1997\).](http://dx.doi.org/10.1007/s002180050345)
- [11] D. R. Tilley *et al.*, [Nucl. Phys.](http://dx.doi.org/10.1016/0375-9474(95)00338-1) **595**, 1 (1995).
- [12] Z. Elekes *et al.*, [Phys. Lett. B](http://dx.doi.org/10.1016/j.physletb.2005.04.007) **614**, 174 (2005).
- [13] Y. Kondo *et al.*, Phys. Rev. C **79**[, 014602 \(2009\).](http://dx.doi.org/10.1103/PhysRevC.79.014602)
- [14] G. Raimann *et al.*, Phys. Rev. C **53**[, 453 \(1996\).](http://dx.doi.org/10.1103/PhysRevC.53.453)
- [15] H. G. Bohlen *et al.*, [Eur. Phys. J. A](http://dx.doi.org/10.1140/epja/i2006-10275-5) **31**, 279 (2007).
- [16] Y. Satou *et al.*, [Phys. Lett. B](http://dx.doi.org/10.1016/j.physletb.2008.01.022) **660**, 320 (2008).
- [17] S. Karataglidis *et al.*, [Nucl. Phys. A](http://dx.doi.org/10.1016/j.nuclphysa.2008.09.007) **813**, 235 (2008).
- [18] M. Stanoiu *et al.*, Phys. Rev. C **78**[, 034315 \(2008\).](http://dx.doi.org/10.1103/PhysRevC.78.034315)
- [19] D. Ridikas *et al.*, [Nucl. Phys. A](http://dx.doi.org/10.1016/S0375-9474(98)00657-5) **628**, 363 (1998).
- [20] A. Yakhelef, N. K. Timofeyuk, J. Al-Khalili, and I. J. Thompson, [Few-Body Syst.](http://dx.doi.org/10.1007/s00601-010-0086-8) **47**, 213 (2010).
- [21] P. Descouvemont, [Nucl. Phys. A](http://dx.doi.org/10.1016/S0375-9474(00)00184-6) **675**, 559 (2000).
- [22] P. Descouvemont and M. Vincke, Phys. Rev. A **42**[, 3835 \(1990\).](http://dx.doi.org/10.1103/PhysRevA.42.3835)
- [23] A. B. Volkov, Nucl. Phys. **74**[, 33 \(1965\).](http://dx.doi.org/10.1016/0029-5582(65)90244-0)
- [24] D. R. Thompson, M. LeMere, and Y. C. Tang, [Nucl. Phys. A](http://dx.doi.org/10.1016/0375-9474(77)90007-0) **286**[, 53 \(1977\).](http://dx.doi.org/10.1016/0375-9474(77)90007-0)
- [25] D. Baye and N. Pecher, Bull. Soc. Acad. Roy. Belg. **67**, 835 (1981).
- [26] B. A. Brown, [Prog. Part. Nucl. Phys.](http://dx.doi.org/10.1016/S0146-6410(01)00159-4) **47**, 517 (2001).
- [27] B. A. Brown and W. D. M. Rae, NUSHELL@MSU, MSU-NSCL Report, 2007 (unpublished).