Two-proton transfer reactions on even Ni and Zn isotopes

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New levels strongly excited by 112-MeV 12 C ions on even Ni and Zn isotopes are J^{π} assigned on kinematical and geometrical arguments, crude shell-model calculations, and distorted-wave Born approximation angular-distribution analysis. These tentative assignments are supported by the Bansal-French model. Because of the contribution of additional collective effects, the two-proton transfer reaction spectra are less selectively fed than those obtained with the analogous two-neutron transfer reactions induced on the same targets in a similar energy range.

I. INTRODUCTION

Most of the high-spin states¹ in f-p shell nuclei have been observed by in-beam spectroscopy using fusionevaporation reactions. These reactions feed only yrast or quasiyrast states that can have many different configurations. On the other hand, transfer reactions are very selective and favor the population of states with a particular configuration.²⁻⁷ The comparison of results from these two kinds of reactions for the same residual nucleus can shed light on the assignment of J^{π} and/or the level configurations.

Obviously the $g_{9/2}$ shell plays an important role in the explanation of high-spin states in the heaviest f-p nuclei, ^{5,6,8} but due to the large size of the involved matrix complete shell-model calculations are scarce. The crude shell model⁹ (CSM), which considers two nucleons moving around an inert target core, predicts the energy of high-spin states formed by two nucleon transfers as the sum of the excitation energies of the corresponding single-particle levels plus (when the two transferred nucleons are identical) an easily calculable pairing energy. Thus the 7⁻ and 8⁺ yrast states in the even isotopes of Zn and Ge have been established as 2n states.⁹ Similarly, the results of (α , ²He) reaction studies^{5,6} on Fe, Ni, Zn, and Ge isotopes agree with the predictions of the CSM.^{9,10}

The 7⁻ and 8⁺ two-proton high-spin states, unobserved up to now, are expected in these nuclei above the similar 2n configuration states.¹⁰ In order to form these states with a maximum cross section, we have performed (¹²C, ¹⁰Be) reactions on Ni and Zn isotopes at an incident energy favoring the transfer of angular momentum that best matches the large angular momentum difference between the grazing waves in the entrance and exit channels. The most important newly observed peaks have been J^{π} assigned according to 2p cluster selection rules, CSM predictions,¹⁰ and angular distribution DWBA analysis. The results are compared with Bansal-French model predictions.¹¹ Among the residual nuclei studied, ⁶⁶Zn is of particular interest since it can be formed with 2n and 2p transfer reactions on 64 Zn and 64 Ni targets, respectively. The relative population of states in both reactions will help to distinguish between 2n and 2p states.

Population of states appears less selective in the $({}^{12}C, {}^{10}Be) 2p$ transfer reactions than in the 2n transfers. This is explained by additional collective effects in the final nuclei that have been taken into account in a semiphenomenological model¹² that considers the coupling of two quasiparticles to a triaxial Davidov rotor in a space including $N=4g_{9/2}$ neutron and proton levels.

II. EXPERIMENTAL PROCEDURE

A beam of 112-MeV ${}^{12}C^{6+}$ was provided by the Strasbourg MP accelerator. Typical beam currents were 50 electrical nA. Targets of self-supporting Ni foils between 80 and 180 μ g/cm² thick, and Zn foils between 85 and 228 μ g/cm² thick on 20- μ g/cm² C backings. These thicknesses have been determined to about 10% with an α -particle gauge. The isotopic enrichments of the ^{58,60,62,64}Ni were 99.89, 99, 97.94, and 98.02% and of the ^{64,66,68}Zn 99.4, 99, and 98.9%, respectively.

Ejectiles were momentum analyzed in a 5- or 10-msr solid angle with a three dipoles and one quadrupole (Q3D) spectrometer $(dp/p=2.10^{-4})$ coupled to the incident beam in an energy-loss mode. A gaseous hybrid counter composed of three proportional counters separated by two ionization chambers of 4- and 15-cm depth was placed along the focal plane at an angle of 45° to the propagation direction of the detected particles. Two 13- μ m Mylar foils separated the isobutane of the counter (100–150 Torr) from the vacuum of the detection chamber, while allowing the particles to reach a 2-cm-thick NE110 scintillator.

The ejectiles, detected in the 11% momentum acceptance of the counter, were identified by their energy loss ΔE in the ionization chambers and residual energy E_R in the scintillator. Correction of the ΔE spectra for entrance angle in the 6° opening of the spectrometer resulted in good mass separation of the detected ejectiles. Position spectra in the first and third proportional counters, separated by 211.5 mm, were obtained by charge division. Conditioned by the ΔE and E_R requirements, the positions allowed reconstruction of the trajectories of the studied ejectiles, the determination of the particle entrance angle, and construction of momentum spectra on the focal plane. The overall resolution was 100-200 keV.

III. ANALYSIS PROCEDURE

To analyze the 2p transfer spectra producing Zn and Ge isotopes we take into account two-nucleon aspects and collective aspects that coexist in these final nuclei.

A. Two-nucleon aspects

1. Semiclassical selection rules

The transfer of a (T=1, S=0) 2p cluster on a 0⁺-target nucleus leads to final states of natural parity and takes place with a large probability when three kinematical conditions, established by Brink, are satisfied within reasonable limits¹³

$$\Delta kR_2 = |k_0R_2 - M_{L_2} - (R_2/R_1)M_{L_1}| \le 2\pi R_2/R \quad , \qquad (1)$$

$$\Delta L = |M_{L_2} - M_{L_1} + k_0 (R_1 - R_2)/2 + Q_{\text{eff}} R / \hbar v \le 2 , \quad (2)$$

$$L_1 + M_{L_1} = \text{even}, \quad L_2 + M_{L_2} = \text{even}.$$
 (3)

 L_1, M_{L_1} and L_2, M_{L_2} are the orbital momenta of the transferred cluster and their z projections in the initial and final nuclei, $k_0 = mv/\hbar$, m is the mass of the transferred cluster, and v is the relative velocity of the two nuclei. R_1 and R_2 are the radii of the projectile and target nuclei,

$$R = R_1 + R_2 ,$$

$$Q_{\text{eff}} = (Q - Z_1^f Z_2^f - Z_1^i Z_2^i) e^2 / R ,$$

where Q is the reaction Q value and $Z_1^f, Z_2^f, Z_1^i, Z_2^i$, the



FIG. 1. Total transition probability of a 2*p*-cluster transfer on ⁵⁸Ni versus the excitation energy and the J^{π} value of the final states in the semiclassical model of Brink [Eq. (4.15) of Ref. 3].

nuclei charges in the initial and final channel. For all reactions considered here k_0R_2 is large (Table I) and final states with large orbital momenta will be favored. Table I gives the values of M_{L_2} and L_2 deduced from Eqs. (1), (2), and (3). In this table the expected orbits favored by these kinematical conditions in the final nuclei are also mentioned.

The total transition probability³ depends on the three kinematical conditions and contains the Clebsch-Gordan coefficients of the angular momentum coupling of the cluster in the initial and final nuclei. It is shown (Fig. 1) that for ⁶⁰Zn, for example, the probability of exciting states with spins in the range 4–8 is much greater than for lower spin values when these states are formed by 2p transfer induced by 112-MeV ¹²C ions, on a 0–10 MeV excitation energy range. Therefore $J^+=6^+$, 7^- , and 8^+ states should be well populated.

TABLE I. Favored two-proton stripping orbits at 112-MeV ¹²C incident energy.

	O ₀	$k_0 R_2$	L_1	M	M,	L_2	Favored
Reactions	(MeV)	(ħ)	(ħ)	(<i>ħ</i>)	(<i>ħ</i>)	-2 (ħ)	orbits
58 Ni(12 C, 10 Be) 60 Zn	-18.65	6.44	0	0	4-8	48	d. f.g
60 Ni(12 C, 10 Be) 62 Zn	-15.91	6.51	0	0	4-7	4-7	d, f,g
${}^{62}Ni({}^{12}C, {}^{10}Be){}^{64}Zn$	-13.35	6.58	0	0	4-6	46	d, f
64 Ni(12 C, 10 Be) 66 Zn	-10.81	6.65	0	0	4-6	4-6	d, f
64 Zn(12 C, 10 Be) 66 Ge	-16.98	6.65	0	0	4-7	4-7	d, f,g
$^{66}Zn(^{12}C, ^{10}Be)^{68}Ge$	-14.53	6.72	0	0	4–7	4-7	d, f,g
68 Zn(12 C, 10 Be) 70 Ge	-12.05	6.79	0	0	4-6	4-6	d, f
${}^{58}Ni({}^{12}C, {}^{10}C){}^{60}Ni$	-11.45	6.43	0	0	4-8	4-8	d. f.g
60 Ni(12 C, 10 C) 62 Ni	-13.42	6.50	0	0	4-8	4-8	d. f.g
62 Ni(12 C, 10 C) 64 Ni	-15.35	6.57	0	0	4-9	4–9	d, f,g
64 Ni(12 C, 10 C) 66 Ni	-16.54	6.64	0	0	4-9	4–9	d, f,g
64 Zn(12 C, 10 C) 66 Zn	-12.80	6.64	0	0	4–9	4-9	d, f, g

2. Two-proton transfer spectroscopic factor

The two-proton transfer spectroscopic factor is proportional to the overlap of the wave functions of the two transferred nucleon relative motion and of the 0s cluster. Figure 2 shows this overlap function versus the total spin J for several configurations.¹⁴ It is observed that the wave functions of the states with spins coupled to their maximum, have large overlap with the 0s cluster wave function. For instance, states of configuration $(1g_{9/2})_{8+}^2$ are expected to be fed very selectively.

The fit of the angular distributions by distorted-wave Born approximation (DWBA) calculations allows us to determine the relative spectroscopy factor α for different final states of a nucleus.

3. Shell-model calculations

The great selectivity of two-nucleon stripping reactions results in the preferential excitation of high-spin states of rather simple configuration.² Thus, simple shell-model calculations have been performed to try to explain them, particularly for $J=0^+$, $T_z=0$ target nuclei. When two nucleons are transferred to the orbits j_1 and j_2 outside the core A_0 , the excitation energy in the final nuclear state (J, T) is given¹⁵ by



FIG. 2. Overlap of the wave functions of the 0s cluster and the two transferred proton relative motion versus total spin J.

$$E^{*}(A_{0}+N+N',j_{1},j_{2},J,T) = E_{B}(A_{0},g.s.) - E_{B}(A_{0}+N+N',g.s.) + \epsilon(j_{1}) + \epsilon(j_{2}) + E_{c} + \langle j_{1}j_{2} | V | j_{1}j_{2} \rangle_{JT}$$
(4)

Pr	Previous works				Present work			
E_x (MeV)	J^{π}	Band ^d	E_x (MeV)	dσ/dω ^e (μb/sr)	Configuration	CSM (MeV)	$J^{\pi\mathrm{f}}$	
0.000 ^{a, b}	0	g.s.	0.0					
1.039 ^{a, b}	2+	g.s.	1.04	73				
1.873 ^{a,b}	2+		1.89					
2.450 ^{a, b}	4+	g.s.	2.47					
2.827 ^{a,b}	3 -	NP	2.84	172				
3.077 ^{a, b}	4+		3.06	245	$(\pi f_{5/2})_{A^+}^2$	3.703		
3.746 ^a	5-	NP	3.74	210	$(\pi p_{3/2}\pi g_{9/2})_{5^{-}}$	4.007		
4.182 ^a	(6+)	g.s.			,			
			4.20	292				
4.251 ^{a, c}	(7-)	NP			$(vf_{5/2}vg_{9/2})_{7}^{-}$	4.146	7-	S
4.814 ^a	(7-)		4.80	125				
					$(\pi f_{5/2}\pi g_{9/2})^{7}$	5.122	(7-)	S, T
5.207 ^{a, c}	(8+)	g.s.	5.20	284	$(vg_{9/2})_{8^+}^2$	5.211	8 +	S
5.464 ^a	(9 ⁻)	NP	5.40					
			5.60					
5.74 ^c	(6+)							
			6.00					
			6.85	196	$(\pi g_{9/2})^2_{8^+}$	6.541	(8+)	S, T
			7.55	180	$(\pi g_{9/2}\pi d_{5/2})_{4+}$	(7.398)	(6+)	S, T

TABLE II. Excitation energy, spin, and parity of ⁶⁶Zn levels.

^aReference 19.

^bReference 20.

^cReference 6.

^dReference 21.

^eCross section at $\theta_{lab} = 10^{\circ}$ with an absolute error of 30%.

^fSpin and parity assignments at the right in the table are based on systematic trends S and on crude shell-model (CSM) calculations T.

where the E_B terms are experimental binding energies¹⁶ (taken to be negative), E_c is the Coulomb interaction energy, and $\epsilon(j)$ are the single-particle energies

$$\epsilon(j) = E_B(A_0 + 1, j) - E_B(A_0, g. s.),$$

with

$$E_B(A_0+1,j) = E_B(A_0+1,g.s.) + E^*(A_0+1,j)$$
.

The last term in Eq. (4), which is the two-body matrix element (TBME), describing the interaction between active particles outside the core, has been calculated¹⁷ or experimentally estimated.⁶ It is generally weak for the configurations considered in this work. In our calculations done in the crude shell-model approximation^{9,10} the TBME and the Coulomb energy terms have been neglected. The positions of the aligned configuration high-spin states corresponding to the filling of the available empty shells of the target nuclei ($f_{5/2}, g_{9/2}$, and $2d_{5/2}$ shells) reported in Tables II–VIII (CSM column), have been calculated from recent mass value¹⁶ and single-particle energies reported in Table IX.

B. Collective aspects

Previous experimental work has shown that prominent features of Zn and Ge spectra are three close-lying $J^{\pi}=8^+$ that are qualitatively explained by an 8^+ of the ground-state band and the alignment of a proton and a neutron $g_{9/2}$ pair, several negative-parity bands, and a γ band. A recent semiphenomenological model¹² gives a quantitative explanation of these characteristics in ^{64,66}Zn and ^{66,68,70}Ge.

In this model a pair of quasiparticles of angular momentum J is coupled to an asymmetric triaxial rotor core of angular momentum R to form a state of total angular momentum I. For the positive-parity states both particles are in the $g_{9/2}$ single-particle orbital, and for the negative-parity states one nucleon is in $g_{9/2}$, while the other is in a $2p_{3/2}$, $2p_{1/2}$, or $1f_{5/2}$ orbital. Amplitude mixing in the wave functions is explicitly given for ⁶⁸Ge and ⁶⁶Ge. The lowest 8⁺ state characterized by R (=I) only, corresponds to the ground-state rotational band (97.2% of the wave function for ⁶⁸Ge). The second 8⁺ state contains an aligned neutron pair $v_1^2 = (vg_{9/2})_{J^{\pi}=8^+}^2$ and rotations from the core with a kinematical moment

	Previous works							
E_x (MeV)	J^{π}	Band ^c	E_x (MeV)	dσ/dω ^d (μb/sr)	Configuration	CSM (MeV)	$J^{\pi\mathrm{e}}$	
0.000 ^{a, b}	0+	g.s.	0.00					
0.991 ^{a, b}	2+	g.s.	0.98	76				
1.799 ^{a,b}	2+	U	1.80					
2.307 ^{a, b}	4+	g.s.	2.40					
2.998 ^{a,b}	3-	NP						
			3.06	434				
3.078 ^{a, b}	4+				$(\pi f_{5/2})_{4^+}^2$	3.514		
3.925 ^{a,b}	5-	NP			4			
3.993	6+	g.s.						
4.156	5	U U	4.10	530	$(\pi p_{3/2}\pi g_{9/2})_{s}$	4.096		
4.237	6+							
4.635 ^{a,b}	7	NP	4.65	185	$(vf_{5/2}g_{9/2})_{7}$	4.645		
4.981 ^{a,b}	7-							
5.151ª	(6,7)							
			5.30	434	$(\pi f_{5/2}\pi g_{9/2})_{7}$	5.058	7-	S , T
5.680 ^{a, c}	$8^{(-)}, (9)^{-}$	NP	5.70	186	1		(8 +)	S
6.124 ^a	(9 ⁻)				$(\nu g_{9/2})_{a^+}^2$	6.156		
			6.30		8			
6.766 ^a			6.70	315	$(\pi g_{9/2})^2_{++}$	6.602	(8 +)	Т
			7 40		0 / 2 8			
			7.90	521	$(\pi g_{9/2} \pi d_{5/2})_{+}$	7.572	(6+)	T^{f}
			1120		······································		() /	-

TABLE III. Excitation energy, spin, and parity of ⁶⁴Zn levels.

^aReference 22.

^dCross section at $\theta_{lab} = 10^{\circ}$ with an absolute error of 30%.

^eSpin and parity assignments at the right in the table are based on systematic trends S and on crude shell-model (CSM) calculations (T).

^fThe 7.90-MeV peak is chosen as (6⁺) rather than the 7.40-MeV one on yield considerations.

^bReference 20.

^cReference 23.

Pr	evious work	S.			Present work	43		
E_x (MeV)	J^{π}	Band ^e	E_x (MeV)	<i>dσ/dω</i> ^f (μb/sr)	Configuration	CSM (MeV)	$J^{\pi g}$	
0.00 ^{a, b}	0+	g.s.	0.0					
0.95 ^{a,b}	2+	g.s.	0.96	19				
1.81 ^{a,b}	2+		1.80					
2.19 ^{a, b}	4+	g.s.	2.20					
3.20 ^{b, c, e}	(3 ⁻)	NP					3-	S
			3.31	116				
3.216 ^d	(4 +)				$(\pi f_{5/2})_{a^+}$	3.610	4+	S, T
3.71 ^e	6+	g.s.			•			
4.05 ^{b, e}	5-	NP	4.17	141	$(\pi p_{3/2}\pi g_{9/2})_{s}$	4.391		
4.347 ^e	6+				J			
4.54 ^b	6+		4.50	47				
			4.70		$(vf_{5/2}vg_{9/2})^{7}$	4.795	7-	S
4.90 ^{a, e}	(7-)				0 372 0 772			
5.122 ^e	5,7	NP	5.19	171	$(\pi f_{5/2}\pi g_{9/2})_{7^{-1}}$	5.361	7-	Т
			6.30		$(\nu g_{9/2})_{g^+}^2$	(6.673)	(8+)	S
			7.54	106	$(\pi g_{9/2})_{8^+}^{2}$	7.112	(8+)	Т
			8.30	30	$(\pi g_{92/}\pi d_{5/2})_{4+}$	(7.80)	(6+)	Т

TABLE IV. Excitation energy, spin, and parity of ⁶²Zn levels.

^aReference 24.

^bReference 25.

^cReference 26.

^dReference 27.

^eReference 28.

^fCross section at $\theta_{lab} = 10^{\circ}$ with an absolute error of 30%.

^gSpin and parity assignments at the right in the table are based on systematic trends S and on crude shell-model (CSM) calculations T.

Present work Previous works E_x E_x $d\sigma/d\omega^{e}$ CSM $J^{\pi\,\mathrm{f}}$ (MeV) J^{π} (MeV) **Band**^d $(\mu b/sr)$ Configuration (MeV) 0.000^{a} $\mathbf{0}^+$ 1.004^a 2+ 1.01 3 g.s. **4**⁺ 2.193^a 2.19 3 g.s. 3.504^a 3-NP 3.59 10 $(\pi f_{5/2})^2_{4^+}$ 3.71 3.531 3.82^{b,d} (0-4) 3.84 4.200^{d} (6⁺) S g.s. 4.36^{a, b, d} (5⁻) (2-6) 4.40 30 $(\pi p_{3/2}\pi g_{9/2})_{5^{-1}}$ 4.746 Т 5.35^{a,b} $(vf_{5/2}vg_{9/2})_{7^{-1}}$ (7⁻) S, T**4**⁺ 5.30 27 (5.45) $(\pi f_{5/2}\pi g_{9/2})_{7^{-1}}$ 5.660 6.63^{a,c,d} (0-4) $(\nu g_{9/2})_{8^+}^2$ 6.95 16 (7.24)(8+) S 7.38^{a,c} 7.35 7.66^{a, b} $(\pi g_{9/2})_{8^{+}}^{2}$ $(\pi g_{9/2} \pi d_{5/2})_{6^{+}}$ $(\pi d_{5/2})_{4^{+}}^{2}$ 7.98 8 7.788 (8^+) Т 8.30 6 8.326 (6⁺) Τ 8.73^{a,c,} 8.863 (4⁺) Т 8.75

TABLE V. Excitation energy, spin, and parity of ⁶⁰Zn levels.

^aReference 29.

^bReference 30.

^dReference 32.

^eCross section at $\theta_{lab} = 10^{\circ}$ with an absolute error of 30%.

Spin and parity assignments at the right in the table are based on systematical trends S and on crude shell-model (CSM) calculations T.

^cReference 31.

	Previous works				Present work			
E_x (MeV)	J^{π}	Band ^c	E_x (MeV)	dσ/dω ^d (μb/sr)	Configuration	CSM (MeV)	Ј ^{πе}	
0.000 ^a	0+	g.s.	0.0					
1.039 ^a	2+	g.s.	1.03	72				
2.153 ^a	4+	g.s.	2.13					
2.561 ^a	3 -	NP	2.55	173				
3.059 ^a	4+		3.05	317	$(\pi f_{5/2})_{4+}^2$	3.071		
3.297ª	6+	g.s.	2.50	101	4			
2 41 78		ND	3.50	101				
3.41/	3	NP			,	2 002		
3.368"	(2 -6)				$(\pi p_{3/2}\pi g_{9/2})_{5^{-1}}$	3.893		
3.955ª	7-	NP	4.02	220	$(vf_{5/2}vg_{9/2})_{7^{-}}$	3.744		
4.202 ^a	8+	g.s.			$(\nu g_{9/2})_{8^+}^2$	(4.142)		
			4.33	591				
4.30 ^b	(7-)				$(\pi f_{5/2}\pi g_{9/2})_{7^{-1}}$	4.467	7-	T, S
4.430 ^a	8+				,			
4.984 ^a			4.92					
			5.42					
			5.73		$(\pi g_{9/2})_{8^+}^2$	5.863	(8+)	Т
			6.60		$(\pi g_{9/2}\pi d_{5/2})_{6^+}$	(6.377)	(6+)	Т
			7.04		$(\pi d_{5/2})_{4+}^2$	(6.891)	(4 +)	Т

TABLE VI. Excitation energy, spin, and parity of ⁷⁶Ge levels.

^aReference 33.

^bReference 34.

^cReference 35.

^dCross section at $\theta_{lab} = 10^{\circ}$ with an absolute error of 30%.

^eSpin and parity assignments at the right in the table are based on systematic trends S and on crude shell-model (CSM) calculations T.

Pre	vious work	s		Present work						
E_x (MeV)	J^{π}	Band ^d	E_x (MeV)	<i>dσ/dω</i> ^e (μb/sr)	Configuration	CSM (MeV)	$J^{\pi\mathrm{f}}$			
0.000 ^{a, b}	0+	g.s.	0.0							
1.016 ^{a, b, c}	2 +	g.s.	1.03	22						
1.778 ^{a,b,c}	2+		1.80							
2.268 ^{a, b, c}	4+	g.s.	2.28							
2.649 ^{a, b}	3 -	NP	2.66	148						
3.041 ^a	(4 +)		3.05	116	$(\pi f_{5/2})^2_{A^+}$	2.838	4+	<i>S</i> , <i>T</i>		
3.650 ^c	5 -	NP	3.67	130	$(\pi p_{3/2}\pi g_{9/2})_{5}^{-}$	4.194				
3.696 ^{a, c}	6+	g.s.			5					
4.054 ^c	7-	NP	4.08	292	$(vf_{5/2}vg_{9/2})_{7}$	(4.056)				
			4.61	156	$(\pi f_{5/2}\pi g_{9/2})_{7^{-}}$	4.553	7-	Т		
4.837 ^{a, c}	(8 +)	g.s.	4.81	210	$(vg_{9/2})_{s^+}^2$	(4.790)	8+	S		
5.151 ^a			5.18		ö					
			5.56							
			6.30		$(\pi g_{9/2})_{8^{+}}^{2}$	6.268	(8+)	Т		
			6.96		$(\pi g_{9/2}\pi d_{5/2})_{6^+}$	(6.940)	(6+)	Т		

TABLE VII. Excitation energy, spin, and parity of ⁶⁸Ge levels.

^aReference 36.

^bReference 37.

^cReference 38.

^dReference 39.

^eCross section at $\theta_{lab} = 10^{\circ}$ with an absolute error of 30%.

^fSpin and parity assignments at the right in the table are based on systematic trends S and on crude shell-model (CSM) calculations T.

Р	revious works				Present work			
E_x (MeV)	J^{π}	Band ^c	E_x (MeV)	dσ/dω ^d (μb/sr)	Configuration	CSM (MeV)	Ј ^{πе}	
0.000 ^{a, b}	0+	g.s.	0.0					
0.957 ^{a,b}	2+	g.s.	0.93	1				
2.174 ^{a,b}	4+	g.s.	2.17					
2.799 ^{a, c}	(3-)	NP	2.80	37			3-	S
3.025 ^{a, c}	(3-,5)							
3.080 ^a			3.06	11	$(\pi f_{5/2})_{*+}$	2.699	4 ⁺	<i>S</i> , <i>T</i>
3.656 ^{a,b,c}	(6 ⁺)	g.s.	3.60		4			
3.685 ^{a,b,c}	5-	NP						
3.830 ^{a,b,c,}	(3 ⁻ ,5 ⁻)		3.83	94	$(\pi p_{3/2} \pi g_{9/2})_{r}$	4.357	5-	<i>S</i> , <i>T</i>
3.841 ^a					1 0/2 0//2 5			
4.207 ^{a,b,c}	7-	NP	4.20	83	$(vf_{5/2}vg_{9/2})_{-}$	4.557		
4.544 ^{a,b,c}			4.59	53	$(\pi f_{5/2} \pi g_{9/2})_{}$	4.548	(7-)	<i>S</i> . <i>T</i>
			4 92	45	0 0/2 0//2 7			,
5.495 ^{a,b,c}	$(6-9)^{-}$	NP	1.52	15				
	(0)))		5.50	37	$(vg_{9/2})^2_{a+}$	5.662	8+	S
5.534ª 6.505ª	(8+)				0772 8			
			6.63	71	$(\pi g_{9/2})_{a^+}^2$	6.397	(8+)	Т
			7.27	76	$(\pi g_{0/2} \pi d_{5/2})$	(7.179)	(6 ⁺)	Т

TABLE VIII. Excitation energy, spin, and parity of ⁶⁶Ge levels.

^aReference 19.

^bReference 40.

^cReference 41.

^dCross section at $\theta_{lab} = 10^{\circ}$ with an absolute error of 30%.

^eSpin and parity assignments at the right in the table are based on systematic trends S and on crude shell-model (CSM) calculations T.

TABLE IX. Single-particle energies (MeV) used in the crude shell-model calculations.

	$2p_{3/2}$	$1f_{5/2}$	$2p_{1/2}$	lg _{9/2}	$2d_{5/2}$
⁵⁹ Cu ^a	0	0.914	0.491	3.043	3.580
⁶¹ Cu ^b	0	0.970	0.475	2.721	(3.406)
⁶³ Cu ^c	0	0.962	0.670	2.506	3.476
⁶⁵ Cu ^d	0	1.115	0.771	2.534	(3.391)
65Gad	0	0.191	(0.062)	2.040	(2.822)
⁶⁷ Ga ^e	0	0.359	0.167	2.074	(2.746)
⁶⁹ Ga ^f	0	0.574	0.319	1.970	(2.484)
⁵⁹ Ni ^a	0	0.339	0.465	3.055	4.462
⁶¹ Ni ^b	0	0.067	0.283	2.122	2.697
⁶³ Ni ^c	0.155	0.087	0	(1.292)	2.297
⁶⁵ Ni ^d	0.692	0	0.063	1.013	1.920
⁵⁹ Zn ^{a,g}	0	(0.90)	(0.54)	(2.68)	
⁶¹ Zn ^b	0	0.124	(0.088)	(2.002)	
⁶³ Zn ^c	0	0.193	0.248	1.704	
⁶⁵ Zn ^d	0.115	0	(0.054)	1.065	1.370
⁶⁵ Ge ^h	0	0.111		1.216	
⁶⁷ Ge ^e	(0.123)	(0.018)	(0)	(0.752)	
⁶⁹ Ge ^f	0.233	0	0.087	(0.398)	
^a Reference 42.			^e Reference 46.		

^bReference 43.

^cReference 44.

^dReference 45.

^fReference 47. ^gReference 50.

^hReference 51.

200

160

COUNTS 120

R=0 and R=2 (64 and 20%, respectively of the wave functions for 68 Ge). The third band (v_2^2) corresponds to two $g_{9/2}$ neutrons coupled to 6⁺ with the core. Petrovici and Faessler¹² predict an 8⁺ state formed by two aligned protons $(\pi g_{9/2})_{J^{\pi}=8^+}^2$ at higher energies, or for ⁶⁸Ge about 1 MeV higher than the third 8⁺ state.

While the 3⁻ level is almost an equal admixture of $(\pi p_{3/2}\pi g_{9/2})_{3^-}$ R=0 and $(\nu p_{3/2}\nu g_{9/2})_{3^-}$ R=0 configurations, two favored negative-parity bands NP1 and NP2 with total angular momentum I = J + R are at higher spins characterized by $(vg_{9/2}vf_{5/2})_{J=7}$ and $(\pi g_{9/2}\pi f_{5/2})_{J=7}$. NP3 is an unfavored negative-parity I = R + J - 1band characterized with and $(vg_{9/2}vf_{5/2})_{J=7}$ at higher spins.

IV. RESULTS

The measured spectra are presented in Figs. 3-7. In Tables II-VII the experimental excitation energies, given with a precision of 50 keV, are compared with previously known levels for ^{66,64,62,60}Zn and ^{70,68,66}Ge nuclei, and absolute differential cross sections are reported. These latter have been determined to about 30% and, depending on the transition and on the target for the well-fed transitions, range from 20-500 μ b/sr for 2p transfer on Ni isotopes (Fig. 5), 100-600 μ b/sr for 2p transfer on Zn isotopes (Fig. 6), and 60–300 μ b/sr for 2n transfer on Ni isotopes (Fig. 7). The 2n values have to be compared with $(\alpha, {}^{2}\text{He})$ cross sections, namely about 50 μ b/sr for transitions to the preferentially populated final states.⁶ Angular distributions in 1° slices have been obtained over the 6° aperture of the spectrometer and fitted with exact finite range (EFR) DWBA calculations (code PTOLEMY) and parameters taken from Ref. 18. Figure 8 shows the fit to the partial angular distribution of a 3^{-} level in 66 Zn and a 7^{-} level in ⁶⁸Ge. Though the shapes of the angular distributions are not typical enough to serve as a J^{π} signature, it is possible to distinguish from the slope between large and low spin values. Values of the deduced

relative spectroscopic factor α are effectively found close to 1 for high-spin states expected as 2N levels, but a more refined analysis taking account of multistep processes would be more adequate.

The 66 Zn spectrum obtained by 2p transfer (Fig. 3) differs from that of the 2n transfer spectrum (Fig. 4) in that it shows a larger number of peaks superimposed on a more important background.

The selectivity of the 2n reaction is well explained by the predominant excitation of the configurations $(vf_{5/2}vg_{9/2})_{7^{-}}, (vg_{9/2})_{8^{+}}^{2}$, and $(vg_{9/2}v2d_{5/2})_{6^{+}}$ as shown by Jahn⁶ with the $(\alpha, {}^{2}He)$ reaction. The yet not reported peak at 6.25 MeV could correspond to the $(v2d_{5/2})^2_{4^+}$ configuration calculated with the CSM to lie at 5.819 MeV. The same predominance of the $(\pi f_{5/2}\pi_{9/2})_{7^{-}}$, $(\pi g_{9/2})_{8^{+}}^{2}$, and $(\pi g_{9/2} \pi 2d_{5/2})_{6^{+}}$ stretched natural-parity configurations, predicted at 5.119, 6.535, and 7.374 MeV, respectively, was expected for the 2p spectrum.¹⁰ While the peaks at 5.20, 6.85, and 7.55 MeV are good candidates for these states, there are several other strong peaks, some corresponding to known states.¹⁹ Indeed, in the 2p transfer case the $2p_{3/2}$, $2p_{1/2}$, and $f_{5/2}$ shells are all available. The filling of these shells leads to configurations of final spin 3^- , 4^+ , 5^- , and could explain the fact that the peaks from the 3^- , 4^+ , and 5^- levels at 2.84, 3.06, and 3.74 MeV, respectively, are strongly excited in the ⁶⁴Ni(¹²C, ¹⁰Be)⁶⁶Zn reaction, but not, or weakly, in ${}^{64}Zn({}^{12}C, {}^{10}C){}^{66}Zn$. In particular, the 5⁻ state (seen in γ -ray spectroscopy) probably has large $(\pi p_{3/2} \pi g_{9/2})_{5^{-1}}$ and/or $(\pi p_{1/2} \pi g_{9/2})_{5^{-}}$ components. Two peaks corresponding to energies close to those of 2n states, i.e., the 4.20-MeV 7⁻ yrast state and the 5.25-MeV state, previously ascribed $(vf_{5/2}vg_{9/2})_{7^{-}}$ and $(vg_{9/2})_{8^{+}}^{2}$, are surprisingly also well excited in the 2p transfer. This argues for 2p components in these states also.

The same considerations hold for the 60,62,64 Zn (Fig. 5) and the ^{66,68,70}Ge spectra (Fig. 6) which appear less selective as the number of neutrons increases. This is contrary to what is observed in the 2n transfer spectra on ^{60, 62, 64, 66}Ni (Fig. 7).

The greater neutron number (2-10) outside the closed $1f_{7/2}$ shell probably contributes to the loss of selectivity of the 2p transfer reactions on the even Ni and Zn iso-



Be)

(Q = -10.806 MeV). This is to be compared with the ⁶⁶Zn spectrum of Fig. 4 obtained in the same conditions by a 2ntransfer reaction (Q = -12.083 MeV).





FIG. 5. Two-proton transfers on 58,60,62,64 Ni. The respective Q values are -18.647, -15.908, -13.350, and -10.806 MeV.

topes, while the reduced proton number (0-2) outside the closed $1f_{7/2}$ shell is of no consequence for the same targets in the 2*n* transfer spectra (Fig. 7) in which the $(v2p_{1/2}vg_{9/2})_{5^{-}}$, $(vf_{5/2}vg_{9/2})_{7^{-}}$, $(vf_{5/2}v2d_{5/2})_{5^{-}}$, $(vg_{9/2})_{8^{+}}^{2}$, and $(vg_{9/2}v2d_{5/2})_{6^{+}}$ configurations dominate.⁶

The background, which increases with the Q value of the reaction is more important in the 2p transfer spectra [Q(2p)=-27.19 MeV compared with Q(2n)=-31.84 MeV] and adds to the difficulty of their interpretation.

In Tables II–VIII are reported the J^{π} proposed on the basis of previous data, angular distribution analysis, systematic spectral trends S among even isotopes, and shellmodel calculations of two-particle configurations (CSM). States related by E2 transitions in the referenced γ -ray



FIG. 6. Two-proton transfers on 64,66,68 Zn. The respective Q values are -16.989, -14.538, and -12.052 MeV.



FIG. 7. Two-neutron transfers on ^{58,60,62,64}Ni. The respective Q values are -11.453, -13.424, -15.347, and -16.768 MeV. The J^{π} values are those deduced from the $(\alpha,^{2}\text{He})$ reactions (Ref. 6).

scheme are indicated in the tables by g.s. and NP for ground-state positive, and negative-parity bands, respectively.

V. DISCUSSION

Two-proton aspects are dominant in the measured $({}^{12}C, {}^{10}Be)$ reaction spectra and the configurations $(\pi g_{9/2})_{8^+}^2$, $(\pi g_{9/2}\pi f_{5/2})_{7^-}$, $(\pi g_{9/2}\pi d_{5/2})_{6^+}$, and $(\pi g_{9/2}\pi p_{3/2})_{5^-}$ allow us to explain the observation of most of the strongest peaks. Table X shows the agreement between experimental and CSM calculated excitation energies. In Table XI are reported the TBME values deduced from the present work.

In Fig. 9 are drawn for each of the main configurations the experimental 2p-binding energies versus the mass



FIG. 8. Angular distributions over the 6° aperture of the spectrometer The continuous line is a DWBA fit with the parameters of Ref. 18.

TABLE X. Crude shell model and experimental energies (MeV) of assumed 2p states in Zn and Ge

isotope	sotopes.													
	$(\pi f_{5/2}\pi g_{9/2})_{7^{-1}}$		$(\pi g_{9/2})_{8^+}^2$		$(\pi g_{9/2}\pi$	$(\pi g_{9/2}\pi d_{5/2})_{6^+}$		$(d_{5/2})^2_{4^+}$						
⁶⁰ Zn	5.660	5.30	7.788	7.98	8.326	8.30	8.863	8.75						
⁶² Zn	5.361	5.19	7.112	7.54	(7.80)	7.54	(8.482)							
⁶⁴ Zn	5.058	5.30	6.602	6.70	7.572	7.90	8.542							
⁶⁶ Zn	5.122	5.20	6.541	6.85	(7.398)	7.55	(8.255)							
⁶⁶ Ge	4.548	4.59	6.397	6.63	7.179	7.27	(7.961)							
⁶⁸ Ge	4.553	4.61	6.268	6.30	6.940	6.96	(7.612)							
⁷⁰ Ge	4.467	4.33	5 863	5 7 3	(6.377)	6 60	(6.891)							





FIG. 9. Two-proton binding energies deduced from the present work for Zn and Ge even isotopes and reported one-proton binding energies for $^{59-65}$ Cu and $^{65-69}$ Ga odd isotopes. The coefficients of the least-square linear fits are reported in Table XII.

TABLE XI. Two-body matrix elements for Zn and Ge isotopes (MeV).

	Previous								
Configuration	values	⁶⁰ Zn	⁶² Zn	⁶⁴ Zn	⁶⁶ Zn	⁶⁶ Ge	⁶⁸ Ge	⁷⁰ Ge	Mean values
$(f_{5/2})^2_{4^+}$	+0.46117	+0.18	-0.30	-0.45	-0.64	+0.36	+0.21	-0.02	-0.09 ± 0.35
$(p_{3/2}g_{9/2})_{5^{-1}}$		-0.35	-0.22	+0.00	-0.27	-0.53	-0.52	-0.39	$-0.33{\pm}0.18$
$(f_{5/2}g_{9/2})_{7^{-1}}$	$-0.10{\pm}0.35^{6}$	-0.36	-0.17	+0.24	-0.32	+0.04	+0.06	-0.14	-0.09 ± 0.22
$(g_{9/2})_{8^{+}1}^{2}$	$+0.22\pm0.51^{6}$	+0.19	+0.43	+0.10	+0.31	+0.23	+0.03	-0.13	+0.17±0.17
$(g_{9/2}d_{5/2})_{6^+,1}$	$+0.18{\pm}0.63^{6}$	-0.03	+0.50	+0.33	+0.15	+0.09	+0.02	+0.22	$+0.18{\pm}0.17$

^aThese two-proton configuration values include the Coulomb interaction energy.

number A of the final nucleus

$$B_{2p}(A_0+2,J) = E_B(A_0+2,g.s.) - E_B(A_0,g.s.) + E^*(A_0+2,J) .$$

These binding energies vary linearly with A, as observed⁶ for the 2n states fed by the $(\alpha, {}^{2}\text{He})$ reaction on the 1f-2p shell nuclei and as predicted by the Bansal-French model.¹¹

In this model, the 2p-binding energy of a level in a nucleus of mass A and atomic number Z is written

$$B_{2p}(A,JT) = B_{2p}(A_0 + 2,JT_0) - 2a(A_0 + 2 - A) + b(T-1) + c(A_0 + 2 - A + 2T_Z)$$
(5)

with the isospin $T = T_Z = (A/2) - Z$ for the considered nuclei. Then,

$$B_{2p}(A,JT) = \left[2a + \frac{b}{2} \right] A + B_{2p}(A_0 + 2, JT_0) -2a(A_0 + 2) - b(Z + 1) + c(A_0 + 2 - 2Z).$$

For isotopes of an element Eq. (5) depends on A only:

 $B_{2p} = kA + l \; .$

This linear dependence results from the linear dependence of the 1p-binding energies versus A. In effect

$$B_{1p}(A,JT) = B_{1p}(A_0 + 1,JT_0) - a(A_0 + 1 - A) + \frac{b}{2}(T - \frac{1}{2}) + \frac{c}{2}(A_0 + 1 - A + 2T_z)$$

which can be written

$$B_{1p}(A,JT) = \left[a + \frac{b}{4}\right]A + B_{1p}(A_0 + 1, JT_0) - a(A_0 + 1)$$
$$-\frac{b}{2}(Z + \frac{1}{2}) + \frac{c}{2}(A_0 + 1 - 2Z) .$$

For isotopes of an element

$$B_{1p}(A,JT) = k'A + l'$$

with

$$k' = a + \frac{b}{4} = \frac{k}{2} ,$$

$$l' = B_{1p}(A_0, JT_0) - a(A_0 + 1)$$

$$- \frac{b}{2}(Z + \frac{1}{2}) + \frac{c}{2}(A_0 + 1 - 2Z) .$$

The coefficients k, l, k', and l' are determined from a least-square fit of the experimental binding-energy values (Fig. 9) and reported in Table XII. The values of k, l are clearly correlated to k', l' and we have $k = k'_1 + k'_2$, $l = l'_1 + l'_2$ to less than 10%. The linearity observed in the 2p case reflects the linearity in the 1p case.¹⁰

In each of the Tables II–VIII appears one positiveparity g.s. and one negative-parity (NP) band which could be associated with the ground-state rotational (GSRB) and one of the negative-parity bands of Petrovici and Faessler.¹²

The ground-state band previously known up to 6^+ in 62,64,66 Zn and 66,68,70 Ge is excited weakly in the corre-

TABLE XII. Coefficients k, l, k' and l' deduced from the linear least-square fits of the experimental two-proton and one-proton binding-energy values.

,

		$(f_{5/2})_{4^+}^2$	$(p_{3/2}, g_{9/2})_{5^{-}}$	$(f_{5/2},g_{9/2})_{7}^{-}$	$(g_{9/2})_{8^+}^2$	$(g_{9/2}d_{5/2})_{6^+}$
Zn	k	1.413	1.406	1.313	1.515	1.515
	1	- 79.895	-80.207	-75.493	-90.239	-91.151
Ge	k	1.235	1.315	1.208	1.458	1.400
	1	-74.425	-80.479	-80.132	-92.722	- 89.535
		2p _{3/2}	1f _{5/2}	2 <i>p</i> _{1/2}	1g _{9/2}	$2d_{5/2}$
Cu	k'	0.671	0.641	0.619	0.758	0.696
	ľ	-36.180	-35.326	-33.575	-44.282	-41.184
Ga	k'	0.665	0.570	0.601	0.683	0.750
	l'	-39.341	-33.301	-35.219	-42.542	-47.687

sponding spectra. In 60 Zn the γ -ray scheme is not well known.²⁹ A recent (3 He, $n\gamma$) study³² performed at a 12-MeV incident energy leads to low J^{π} value levels, weakly or not fed in our spectrum. The probable doublet peak we observe at 4.40 MeV could contain the 6⁺ ground-state band member, if it follows the trend in the location of the 6⁺ states in the other nuclei.

Of the three 8^+ branches in forks reported in some of the Zn and Ge isotopes there subsists only one 8^+ state in our spectra, very close to the calculated position of the $(\nu g_{9/2})_{8^+}^2$ configuration. In ⁶⁸Ge where the triple forking was first observed, 48 the nature of the 8^+ states at 4.837, 5.050, and 5.367 MeV differs according to several interpretations.^{12, 38, 49} Recent g-factor measurements³⁹ order them as v_1^2 , v_2^2 , and GSRB states. This does not agree with the order calculated by Petrovici and Faessler¹² and lends support to the importance of the neutrons of the target outside the closed $f_{7/2}$ shell, as mentioned in Sec. IV. This leads us to propose that the peaks at 5.50 MeV in 66 Ge and 5.20 MeV in 66 Zn are 8^+ levels of the same $(v_{9/2})^2$ configuration. The same assumption for peaks at 5.70 MeV in ⁶⁴Zn and 6.30 MeV in ⁶²Zn is more questionable because the correspondence with the calculated energy is not as good and the collectivity effect is expected to decrease with the neutron number.

Calculation of the energy of the $(vg_{9/2})_{8^+}^2$ configuration in ⁶⁰Zn requires the position of the $g_{9/2}$ single-particle state in ⁵⁹Zn that is yet unknown.⁴² A level seen at 2.68 MeV in the ⁵⁸Ni $(p, \pi^-)^{59}$ Zn reaction⁵⁰ is a possible candidate and would lead to a $(vg_{9/2})_{8^+}^2$ excitation energy of 7.24 MeV, which is close to the 6.95-MeV peak in ⁶⁰Zn that systematics indicates as an 8⁺ state.

Three strong peaks observed in the ⁶⁶Zn and ^{70,68}Ge spectra belong to a known negative-parity band (see Tables II, VI, and VII). Respectively, the 3⁻ members are at 2.84, 2.66, and 2.55 MeV, while the 5⁻ members are at 3.74, 3.67, and 3.50 MeV (not resolved in our spectra from the 6^+ GSRB member in the two last cases) and the 7⁻ members are at 4.20 MeV (also mixed with the 6⁺ GSRB member), 4.08 and 4.02 MeV. In the isotones ⁶⁴Zn and 66 Ge (see Tables III and VIII) the 3⁻, 5⁻, 7⁻ members of the same band are reported at 2.998, 3.925, and 4.635 MeV and at (2.799), 3.685, and 4.207 MeV, respectively. All of these states correspond to intense peaks in our spectra except the 5⁻ members that are weakly excited. The agreement between the location of the 7⁻ members, and the calculated energy of the $(vf_{5/2}vg_{9/2})_{7^{-1}}$ configuration as well as the apparent dependence of the relative yields on the neutron number of the target nuclei lead us to propose that we excite the negative-parity band NP1 of Petrovici and Faessler¹² that is composed of mixed neutron and proton configurations for lower spins, and levels of mainly $(vf_{5/2}vg_{9/2})_{J=7}$ I=J+R nature for higher spins.

A second negative-parity band NP2 built on a 5⁻ state is predicted¹² a little higher in energy with a main configuration $(\pi f_{5/2}\pi g_{9/2})_{J=7}$ I=J+R. In fact, the calculated energies of the high-spin stretched 2*p* states, which are expected to be strongly excited, correspond to those of the remaining strong spectral peaks. However,

few of these states were previously J^{π} assigned or even known (see Tables II–VIII). We propose 4^+ as J^{π} for the 3.71- or 3.82-MeV state in ⁶⁰Zn and the 3.06-MeV state in ⁶⁶Ge, because these levels are close to the calculated energy of the $(\pi f_{5/2})_{4^+}^2$ configuration and also in agreement with the systematic trend. The $(\pi p_{3/2}\pi g_{9/2})_{5^{-1}}$ configuration compatible with all of the previous assignments (see Tables II-VIII) is often close in energy to the 5⁻ member of the NP1 band, and in some cases, it is difficult to distinguish between them. In ⁶⁴Zn and ^{66,70}Ge, where the NP1 member corresponds to the lowest reported 5⁻ state, we attribute on yield arguments the $(\pi p_{3/2} \pi g_{9/2})_{5^{-}}$ configuration to the 4.156-, 3.830- $(3^{-}, 5^{-})$, and 3.568-MeV $(2^{-}-6^{-})$ states, respectively. In ⁶⁸Ge two 5⁻⁻ states are reported at 3.582 and 3.650 MeV, but are not distinguishable in our spectrum. Because of the systematic trends and the relative γ -ray decay rate³⁹ of the 4.054 MeV, 7^- NP1 member to the 3.582 MeV, $5^$ state (52%) and the 3.650 MeV, 5^- state (31%), we argue that the 5^- NP1 member is at 3.582 MeV and that the 3.650 MeV, 5⁻ state, is mainly of a $(\pi p_{3/2} \pi g_{9/2})_{5^-}$ configuration. The same configuration could explain the important peaks at 4.40 MeV previously J assigned (2-6) in 60 Zn, at 4.17 MeV in 62 Zn and 3.74 MeV in 66 Zn.

For all the considered nuclei there exist peaks in the vicinity of the predicted values for the $(\pi f_{5/2}\pi g_{9/2})_{7^{-}}$, $(\pi g_{9/2})_{8^{+}}^{2}$, and $(\pi d_{5/2}\pi g_{9/2})_{6^{+}}$ configurations (see Tables II–VIII). Reported (7⁻) states at 4.814 MeV in ⁶⁶Zn and 4.30 MeV in ⁷⁰Ge reinforce our proposition for the 4.80-MeV peak (Fig. 5) and the 4.33-MeV peak (Fig. 6).

VI. CONCLUSION

The present study of the $({}^{12}C, {}^{10}Be)$ 2*p* transfer reactions with 112-MeV ${}^{12}C$ on Ni and Zn isotopes reveals new 2*p* levels in the ${}^{60,62,64,66}Zn$ and ${}^{66,68,70}Ge$ final nuclei and leads to high-spin and parity assignments that are proposed on systematical trends, and crude shell-model calculations, and are supported by the Bansal-French model.¹¹

This work complements the $(\alpha, {}^{2}\text{He})$ 2n transfer reaction data.⁶ The $(\pi g_{9/2})_{8^+}^2$ and $(\pi g_{9/2} \pi d_{5/2})_{6^+}$ states located a few MeV higher than the corresponding 2n states are not populated by fusion-evaporation reactions, but some proposed $(\pi f_{5/2}\pi g_{9/2})_{7^{-}}$ states could be identified with 7⁻ states seen in γ spectroscopy. The lesser selectivity of the 2p transfer as compared to the 2n transfer on the same targets can be related to the greater number of neutrons than protons outside the closed $f_{7/2}$ shell in the target. This fact adds some collectivity to the spectra that are well explained by the asymmetric rotor model with an admixture of two quasiparticles.¹² It explains why we can have serious doubts about whether DWBA calculations could be trusted for the two-proton transfer nuclei discussed here, but a more refined theoretical analysis taking into account multistep processes lies beyond the scope of this work.

Increasing the incident energy should enhance the selectivity as has been found for 2p transfers on ⁵⁴Fe and

lighter targets with 40-MeV/nucleon ¹²C projectiles,⁷ and the evolution of the spectra with energy could be interesting to locate the two-proton stretched states.

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