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³Although we have taken data by separate trigger with acceptance at smaller angles, their analysis is not yet complete.

⁴A more certain, but biased π^0 definition requires two reconstructed photons with π^0 effective mass. However, this definition yields poor efficiency at high P_T . When two reconstructed showers coexist in the same octant, the π^0 energy was taken as that of the highest-energy shower. The inclusion of the lower-energy shower or additional nearby energy in the π^0 definition has no significant effect on our results.

⁵The spectrum has been weighted for geometrical and energy efficiency as determined by calibration runs at the CERN proton synchrotron so that the two π^0 's, when given random angles would have a flat distribution, independent of transverse momentum above $P_T = 1.2$ GeV/c. Trigger corrections due to the ISR intersection angle have been made. The weight of an event varies from 0.3 to 3, with an average weight of 1.

⁶We have also examined the product $\Pi = P_{T_1} \cdot P_{T_2} \cdot P_{TG}$. We find distributions similar to those of the E_T selection.

⁷C. Michael and L. Vanryckegham, University of Liverpool Report No. LTH 31, May 1977 (to be published), and private communication. We compared our spectrum in $\Delta \varphi$ with that predicted by the independentemission model for events with $9 < E_T < 12$ GeV and $P_{T_1} > 1$ GeV; $P_{T_2} > 3$ GeV. The 412 events in our sample have a distribution essentially the same as that of Fig. 2(b). The model predicts a nearly uniform distribution with variations of < 20% over our azimuthal coverage.

 8 The diplot of Fig. 3 is not corrected for efficiency which varies by less than 30% in the populated regions.

⁹Recent measurements of $\Upsilon(9.5) \rightarrow e^+e^-$ in the same apparatus [J. H. Cobb *et al.*, Phys. Lett. <u>72B</u>, 273 (1977)], combined with an assumed 3% branching ratio to electrons, would indicate that there are several hundred decays of $\Upsilon(9.5)$ to hadrons in the data sample from which the present events are drawn.

Atomic Electron Correlation in Nuclear Electron Capture

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The effect of electron-electron Coulomb correlation on orbital-electron capture by the nucleus has been treated by the multiconfigurational Hartree-Fock approach. The theoretical ⁷Be L/K capture ratio is found to be 0.086, and the ³⁷Ar M/L ratio, 0.102. Both ratios are smaller than the independent-particle predictions. Measurements exist for the Ar M/L ratio, and agreement between theory and experiment is excellent.

Benoist-Gueutal's insight¹ that atomic electrons must be included in a complete description of orbitalelectron capture by the nucleus² led to the introduction of atomic exchange and imperfect-overlap factors in the theoretical capture probability.³⁻⁶ All existing work on electron capture has been carried out in the independent-particle approximation; effects due to electron-electron Coulomb correlation have been neglected. Here we report on a first effort to take correlation into account, by using the multiconfigurational Hartree-Fock (MCHF) approach.⁷ We calculate the ⁷Be L/K and ³⁷Ar M/L capture ratios.

The nuclear-electron-capture rate is²

 $\lambda_i = \lambda_i^0 B_i, \quad i = K, L, M, \ldots,$

(1)

where λ_i^0 is the rate obtained when atomic matrix elements are neglected,⁸ and B_i is the exchange-

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overlap correction factor. For example, if the initial and final states are represented by a single Slater determinant, then

$$B_{k} = K \left\{ \langle 2s' | 2s \rangle \langle 3s' | 3s \rangle - \langle 2s' | 1s \rangle \langle 3s' | 3s \rangle [R_{2s}(0)/R_{1s}(0)] - \langle 2s' | 2s \rangle \langle 3s' | 1s \rangle [R_{3s}(0)/R_{1s}(0)] \right\}^{2}, \quad (2)$$

where

$$K = \langle \mathbf{1}s' | \mathbf{1}s \rangle^{2} \langle \mathbf{2}s' | \mathbf{2}s \rangle^{2} \langle \mathbf{2}p' | \mathbf{2}p \rangle^{2q(\mathbf{2}p)} \langle \mathbf{3}s' | \mathbf{3}s \rangle^{2[q(\mathbf{3}s)-1]} \langle \mathbf{3}p' | \mathbf{3}p \rangle^{2q(\mathbf{3}p)}.$$

$$\tag{3}$$

Here, q(nl) is the occupation number of the nlshell, and primes denote the daughter atom. Bahcall²⁻⁶ set K=1, while Vatai^{2,9} retained the factor. Similar expressions exist for B_L and B_M .

The capture ratio for shells i and j, in allowed transitions, is²

$$(\lambda_i/\lambda_j) = (\lambda_i/\lambda_j)^0 (B_i/B_j), \qquad (4)$$

where

$$(\lambda_{i}/\lambda_{j})^{0} = [R_{i}^{2}(0)/R_{j}^{2}(0)](q_{i}^{2}/q_{j}^{2}),$$

$$i, j = K, L_{1}, M_{1}.$$
(5)

The R's are electron radial wave functions, evaluated at the origin, and the q's are neutrino energies. The contributions from L_2 and M_2 electrons are neglected here.

In our MCHF calculation, the ground state is

$$\Psi_{s}(\gamma LS) = \sum_{i} C_{i} \Phi(\gamma_{i} LS), \qquad (6)$$

and the final-state wave function, describing the hole state after capture, is

$$\Psi_{j}'(\gamma LS) = \sum_{i} C_{ji}' \Phi_{i}'(\gamma_{i} LS) \,. \tag{7}$$

The atomic matrix elements become

$$\langle \Psi_{j} | O | \Psi_{g} \rangle = \sum_{i,k} C_{jk}' C_{i} \langle \Phi_{k}' | O | \Phi_{i} \rangle, \qquad (8)$$

where we have $O = \sum_{b} a_{b} R_{b}(0)$, and a_{b} is the destruction operator.⁴ The exchange-overlap correction factor is

$$B_{i} = \sum_{j} |\langle \Psi_{j} | O | \Psi_{g} \rangle / R_{i}(0) |^{2}, \qquad (9)$$

where the summation extends over the states included in the multiconfigurational expansion.

For the ⁷Be L/K-capture-ratio calculation, the ground state is represented by

$$\Psi_{g} = C_{1} \Phi_{1} (1s^{2}2s^{2}) + C_{2} \Phi_{2} (1s^{2}2p^{2}).$$
(10)

The 1s-hole state after K capture is

$$\Psi_{j} = C_{j1}' \Phi_{1}' (1s2s^{2}) + C_{j2}' \Phi_{2}' (1s2p^{2}).$$
(11)

The 2s-hole state after L_1 capture is represented by the single configuration

$$\Psi_i = \Phi'(1s^2 2s). \tag{12}$$

For the ³⁷Ar M/L-capture-ratio calculation, we take the ground-state MCHF wave function to be

$$\Psi_{g} = C_{1}\Phi_{1}(1s^{2}2s^{2}2p^{6}3s^{2}3p^{6}) + C_{2}\Phi_{2}(1s^{2}2s^{2}2p^{6}3p^{6}3d^{2}(^{1}S)) + C_{3}\Phi_{3}(1s^{2}2s^{2}2p^{6}3s^{2}3p^{4}(^{1}S)3d^{2}(^{1}S)) + C_{4}\Phi_{4}(1s^{2}2s^{2}2p^{6}3s^{2}3p^{4}(^{3}P)3d^{2}(^{3}P)) + C_{5}\Phi_{5}(1s^{2}2s^{2}2p^{6}3s^{2}3p^{4}(^{1}D)3d^{2}(^{1}D)).$$
(13)

The 2s-hole state is

$$\Psi_{j} = C_{j1} \Phi_{1}' (1s^{2}2s2p^{6}3s^{2}3p^{6}) + C_{j2} \Phi_{2}^{-1} (1s^{2}2s2p^{6}3s^{2}3p^{4}(^{1}S)) + C_{j3} \Phi_{3}' (1s^{2}2s2p^{6}3s^{2}3p^{4}(^{3}P)^{4}P 3d^{2}(^{3}P)) \\ + C_{j4} \Phi_{4}' (1s^{2}2s2p^{6}3s^{2}3p^{4}(^{3}P)^{2}P 3d^{2}(^{3}P)) + C_{j5} \Phi_{5}' (1s^{2}2s2p^{6}3s^{2}3p^{4}(^{1}D) 3d^{2}(^{1}D)).$$
(14)

The 3s-hole state after M_1 capture is

$$\Psi_{j} = C_{j1}\Phi_{1}'(1s^{2}2s^{2}2p^{6}3s^{3}p^{6}) + C_{j2}\Phi_{2}'(1s^{2}2s^{2}2p^{6}3s^{2}3p^{4}(^{1}D)3d) + C_{j3}\Phi_{3}'(1s^{2}2s^{2}2p^{6}3s^{3}p^{4}(^{1}S)3d^{2}(^{1}S)) + C_{j4}\Phi_{4}'(1s^{2}2s^{2}2p^{6}3s^{3}p^{4}(^{3}P)^{4}P^{3}d^{2}(^{3}P)) + C_{j5}\Phi_{5}'(1s^{2}2s^{2}2p^{6}3s^{3}p^{4}(^{3}P)^{2}P^{3}d^{2}(^{3}P)).$$
(15)

	<i>K</i> capture			L_1 capture	
	$ 1s\rangle$	$ 2 s\rangle$	2p angle	$ 1s\rangle$	$ 2s\rangle$
$\langle 1s' $	0.972 09	- 0.190 99		0.962 47	- 0.155 91
$\langle 2s' $	0.171 93	0.96785		0.08271	0.88283
<2 p'			0,992 60		

TABLE I. MCHF $\langle nl' | nl \rangle$ overlap integrals for ₄Be electron capture.

	$ 1s\rangle$	$ 2s\rangle$	$ 2p\rangle$	$ 3s\rangle$	3 <i>p</i> >	$ 3d\rangle$
			L_1 capture	е		
$\langle 1s' $	0,99873	-0.02977		-0.00630		
$\langle 2 s' $	0.02705	0.992 50		-0.10496		
$\langle 2p' $			0.998 58		-0.02279	
$\langle 3s' $	0.00798	0.10177		0.99274		
$\langle 3p' $			0.02142		0,99927	
$\langle 3d' $						0.99954
			M_1 capture	е		
$\langle 1s' $	0.99875	-0.02921		-0.00628		
$\langle 2s' $	0.02623	0.99228		- 0.09736		
$\langle 2p' $			0.99445		-0.08177	
$\langle 3s' $	0.00702	0.09020		0.98913		
⟨3 <i>p</i> ′			0.07552		0.99047	
⟨3 <i>d</i> ′						0.93200

TABLE II. MCHF $\langle nl' | nl \rangle$ overlap integrals for ₁₈Ar electron capture.

TABLE III.	Electron radial wave-function ratios
$R_{ns}^{2}(0)/R_{n's}^{2}(0)$), exchange-overlap correction factors
B_{i} , and capture	re ratios λ_i / λ_i .

Element	Quant	tity	Result
⁷ ₄ Be	$R_{2s}^{2}(0)/R_{1s}^{2}(0)$,	HF MCHF ^{.c}	0.0332 0.0300
	B_K ,	HF (V) ^a HF (B) ^b MCHF ^c	0.816 0.900 0.792
	B_L ,	HF (V) ^a HF (B) ^b MCHF ^c	2.222 3.045 2.259
	λ_L/λ_K ,	HF (V) ^a HF (B) ^b MCHF ^c	0.090 0.112 0.086
$^{37}_{18}{ m Ar}$	$R_{3s}^{2}(0)/R_{2s}^{2}(0)$,	HF MCHF ^c	0.0977 0.0669
	B_L ,	HF (V) ^a HF (B) ^b MCHF ^c	1.121 1.171 1.098
	Β _Μ ,	HF (V) ^a HF (B) ^b MCHF ^c	1.322 1.549 1.674
	λ_M/λ_L ,	HF (V) ^a HF (B) ^b MCHF ^c Experiment ^d	$\begin{array}{c} 0.115\\ 0.129\\ 0.102\\ 0.104 \substack{+0.007\\-0.003}\end{array}$

^aHartree-Fock, Vatai's approach (Refs. 2 and 9). ^bHartree-Fock, Bahcall's approach (Refs. 2 and 3-6).

 $^{\rm c}{\rm Present}$ multiconfigurational HF calculation. $^{\rm d}{\rm Ref.}$ 11.

The MCHF wave functions, including the amplitudes C, were computed with the Froese Fischer program.⁷ The electrostatic interaction matrix elements were calculated with Hibbert's program.¹⁰ The one-electron overlap integrals are listed in Tables I and II. The electron radial-wave-function ratios at the origin and the overlap-exchange correction factors B_i as well as the electron-capture ratios are listed in Table III. For comparison, theoretical single-configuration HF capture ratios² and the experimental result¹¹ for ³⁷Ar are also listed; there is no measurement of the ⁷Be L/K ratio.

Electron correlation is seen to have a substantial effect on nuclear capture ratios when outer electrons are involved. Compared with singleconfiguration HF results according to Vatai's approach,² the MCHF L/K capture ratio of ⁷Be is reduced by 4.4%; the ³⁷Ar M/L ratio is reduced by 11% and brought into excellent agreement with experiment.^{2, 11}

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Population of Resonant ¹²C+¹²C States via the Reaction ¹²C(¹⁶O, α)²⁴Mg

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Excitation functions for the reaction ${}^{12}C({}^{16}O, \alpha){}^{24}Mg$ have been measured from $E_{1ab}({}^{16}O) = 62$ to 110 MeV using a counter telescope at $\theta_{1ab} = 7.5^{\circ}$. Selective population of relatively few states at very high excitation energies in ${}^{24}Mg$ [$E_x({}^{24}Mg) > 20$ MeV] is seen. A possible correspondence is found between these states and the narrow resonances reported in ${}^{12}C + {}^{12}C$ reactions. In addition, a possible correspondence between averaged yields in ${}^{12}C({}^{16}O, \alpha){}^{24}Mg$ and gross structure seen in several ${}^{12}C + {}^{12}C$ reaction channels is cited.

The reaction ${}^{12}C({}^{16}O, \alpha)^{24}Mg$ has been studied extensively because of its striking final-state selectivity. It is interesting to consider whether this behavior reflects special structures in the initial, compound, and final systems. In this work we have significantly extended the ${}^{12}C({}^{16}O, \alpha)^{24}Mg$ excitation functions. An apparent correlation has been found between individual pronounced transitions in that reaction and resonant states which have been seen in ${}^{12}C + {}^{12}C$ reactions. Furthermore, the envelope of the ${}^{12}C({}^{16}O, \alpha)^{24}Mg$ transitions shows a weak correlation with gross structure variations in the ${}^{12}C$ strength function as indicated by several ${}^{12}C + {}^{12}C$ reaction-channel excitation functions.

High-resolution spectra from the reaction ${}^{12}C({}^{16}O, \alpha)^{24}Mg$ were measured from $E_{1ab}({}^{16}O)$ = 62.0 to 100.0 MeV in 1-MeV steps. Two measurements at $E_{1ab}({}^{16}O)$ = 105.0 and 110.0 MeV were also taken. The experiment was performed at the Brookhaven National Laboratory Tandem Van de Graaff facility and employed a surface-barrier counter telescope placed at θ_{1ab} =7.5°±0.25°. Targets were nominally 45- μ g/cm² natural carbon. The experimental resolution was typically 90 keV.

The primary objective of the study was to determine if the ${}^{12}C({}^{16}O, \alpha)^{24}Mg$ spectra at higher energies show any gross- or fine-structure correlations to previously known ${}^{24}Mg$ states which have been observed via ${}^{12}C + {}^{12}C$ resonance reactions. The α spectra contain numerous previously unseen transitions in the range of $E_x({}^{24}Mg) = 20$ to 35 MeV—a region in which most ${}^{12}C + {}^{12}C$ resonances have been recorded. The α -transition yields exhibit compound-nuclear fluctuations; to average out this effect and to enhance the persistently strong transitions, the spectra were averaged over the incident ¹⁶O energy. This was performed by linearizing the 39 individual α spectra to a common ²⁴Mg excitation-energy scale. Kinematic corrections were performed so the energysummed spectra would reflect center-of-mass cross sections. A smooth evaporative background was subtracted by hand from each spectrum to further enhance strong discrete transitions in the summed spectra. Figure 1 shows three typical linearized spectra and background curves. It is understood that the magnitude of the underlying



FIG. 1. Typical ${}^{12}C({}^{16}O, \alpha)^{24}Mg$ spectra at $E_{1ab} = 63$ (curve *a*), 77 (curve *b*), and 91 MeV (curve *c*). They have been linearized in $E_x({}^{24}Mg)$ and the smooth curves are hand drawn to represent the background that is to be subtracted.