Correlated nucleons in configuration space

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Several recent studies have dealt with the effects of short-range correlations on the momentum distribution of nucleons in nuclei. Here we investigate the correlation effects on the density and spectral distribution in coordinate space. A combination of experimental data and spectral functions calculated from realistic nucleonnucleon interactions allows us to resolve a recently uncovered discrepancy with occupation of quasiparticle states derived from $(e, e'p)$ data.

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Much of the understanding of atomic nuclei is based on the assumption that nucleons move, independently from each other, in the average potential created by the interaction with all other nucleons. A more fundamental approach to the understanding of nuclei has to start from the underlying nucleon-nucleon (*N-N*) interaction. Realistic models of the *N*-*N* interaction exhibit a strongly repulsive central interaction at small internucleon distances and a strong tensor component. These features lead to properties of nuclear wave functions that are beyond what is describable in terms of independent particle (IP) motion. In particular, strong shortrange correlations (SRC) are expected to occur.

The effects of the short-range correlations are known for systems where an accurate solution of the Schrödinger equation for a realistic *N*-*N* interaction can be obtained [1]. Very light nuclei (today up to $A \le 10$) and infinite nuclear matter are amongst the systems where this is feasible [2–4]. The corresponding calculations show that in a microscopic description of nuclear systems, the short-range and tensor parts of the *N*-*N* interaction have a very important, not to say dominating, influence, without which not even nuclear binding can be explained.

Due to these short-range correlations the momentum distributions of nucleons acquire a tail extending to very high momenta *k*, and part of the strength, located in IP descriptions at low excitation energy *E*, is moved to very high excitation energies.

In the past, most experimental investigations were confined to rather low momenta and energies, i.e., to the region where the strength is dominated (but not entirely given) by the IP properties. In this region, the consequences of shortrange correlations are indicated primarily by a *depopulation* of states in comparison to the predictions of the IP models (including the long-range correlations which can be described by configuration mixing). According to the calculations mentioned above, a depopulation on the order of 20% is expected.

From the experimental information available up to now, the depopulation of IP strength at low k , E is established [1] (for a caveat see below). Much less is known from a *direct* measurement of the strength of the spectral function $S(k,E)$ at large *k* and *E*. A recent $(e, e'p)$ experiment, performed at high momentum transfer *q* in parallel kinematics by Rohe *et* *al.*, provides a first direct measurement [5,6].

This correlated strength has always been discussed in *k*, *E* space where (part of) it can be separated from the IP strength. In this Rapid Communication, we take an orthogonal look at the correlated strength and discuss it in coordinate space (*r* space). We address this question from both the theory and experiment side.

The trigger for this study lies in difficulties experienced in the past in interpreting data in terms of IP models. For example, fits with IP wave functions of the nuclear charge density often yielded form factors (of IP-dominated transitions) with incorrect *q* dependence. Fits with IP wave functions also have difficulties in reproducing the total densities in the nuclear interior. The origin of these difficulties: total densities have contributions from correlated nucleons that do not appear in observables dominated by quasiparticle properties. The correlated nucleons presumably have a different radial distribution.

The goal of this paper is to derive quasiparticle (QP) and correlated distributions in *r* space using Green's-function theory. We compare the results to the correlated density in *r*-space which we obtain from the difference of the density known from elastic electron scattering—and the QP contributions known from $(e, e'p)$ reactions.

As a side product, this study also sheds light on a recently uncovered problem with QP occupation numbers derived from $(e,e'p)$ experiments with low and high *q*, respectively [7].

The evaluation of the single-particle spectral functions for 12 C, the nucleus we use for our study, has been performed within the framework of the Green's-function method [8] using the techniques described in [9–11]. The nucleon selfenergy $\Sigma_{lj}(p_m, p_n, E)$ is determined in a discrete basis of Bessel functions $\phi_{p_m l j}(r)$ with appropriate boundary conditions at the surface of a spherical box with radius R_{box} . These basis states are identified by the angular momentum quantum numbers *l* and *j* and a radial quantum number (p_m, p_n) referring to the momentum. For a box radius R_{box} of typically 20 fm, it turns out to be sufficient to include around 60 basis states for each partial wave.

The self-energy contains a Hartree-Fock contribution Σ_{lj}^{HF} calculated in terms of a nuclear matter *G* matrix plus com-

FIG. 1. Comparison of QP and correlated densities from theory.

plex correction terms, $\Delta\Sigma_{lj}$, which account for the inclusion of two-particle one-hole and two-hole one-particle contributions. These correction terms are calculated directly for the finite nucleus ${}^{12}C$, describing the intermediate particle states by plane waves orthogonalized with respect to the occupied hole states. This is a good approximation to describe the effects of SRC; however, it tends to underestimate the spectral strength due to long-range correlations at missing energies slightly above the two-hole one-particle threshold.

The single-particle Green's function is determined from this self-energy $\Sigma_{lj}^{\text{HF}} + \Delta \Sigma_{lj}$ by solving the Dyson equation in the box basis described above. From the imaginary part of this Green's function one can calculate the spectral function in this basis [11] or determine it in configuration space by the transformation

$$
S_{lj}(r,r';E) = \sum_{m,n} \phi_{p_mlj}(r) S_{lj}(p_m,p_n;E) \phi_{p_nlj}^*(r'),
$$

using the Bessel functions $\phi_{p_m l j}(r)$ described above. The spectral function can be split into the QP contribution S_{ij}^{QP} , which only occurs in the $s_{1/2}$ and $p_{3/2}$ partial waves, and in the continuum contribution, S_{ij}^{cont} , which originates from the imaginary components in the self-energy. This leads to the single-particle density

$$
\rho(r) = \sum_{lj} S_{lj}^{QP}(r,r) + \sum_{lj} \int_{\varepsilon_{2hlp}}^{\infty} dE S_{lj}^{cont}(r,r;E)
$$

$$
= \rho_{QP}(r) + \rho_{corr}(r),
$$

where the integration over missing energies *E* starts at the threshold of two-hole one-particle configurations. We have assigned the label "*corr*" to the part of the single-particle density, which originates from the continuum part of the spectral function to indicate that this correlated density is absent in the IP model.

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Results for these contributions to the point density of protons in 12 C are displayed in Fig. 1. The calculation of these densities has been performed using the CD-Bonn potential for the *N*-*N* interaction [12]. A fraction of the proton density, which accounts for around five protons, is described by the QP part and the rest is covered by the correlated density.

In order to allow for a better comparison of the radial shape of the density contributions, Fig. 1 also contains the correlated density $\rho_{corr}(r)$ multiplied by a factor of 3. The comparison shows quite clearly that the correlated density is located much more in the center of the nucleus than the QP contribution.

The correlated single-particle density is distributed over partial waves *lj* including those which are unoccupied in the IP model. A large fraction of the correlated strength, however, is contained in partial waves with *l*=0 (around 31%) and *l*=1 (around 37%). One also should note that, contrary to what one naively could expect, the strength in the higher *l* states does not contribute at large *r*; the corresponding large values of *E* pull the radial wave functions to lower *r*.

For 12 C an extensive set of elastic electron scattering data is available [13–16], covering the range of momentum transfer between 0.13 and 3.7 fm⁻¹. The carbon rms radius is precisely known from a μ -x-ray experiment [17].

These data have been used to determine a modelindependent charge density using the SOG approach of [18]. The procedure employed has been described in [19]. This yields the charge density as a function of *r*, together with an error bar that covers both the statistical and systematic uncertainties of the data, as well as the model error.

In order to obtain the point density, we have unfolded the effect of the finite size of the proton and neutron charge density. This has been done by parametrizing the point density, folding it with $\rho_p(r) + \rho_n(r)$, and fitting the resulting density to the charge density as determined above. The contribution of the electromagnetic spin-orbit term to the charge density turned out to be negligible. The resulting folded density agrees within \sim 1% with the one given in [20].

The resulting point density is shown in Figs. 2 and 3. The error bars in general are too small to be seen.

For carbon, quite an extensive set of $(e, e[']p)$ data is available [21–28]; a compilation is discussed in [7]. Part of this data has been taken at low *q* with the goal to determine the 1*p* and 1*s* quasiparticle momentum distributions and occupation numbers. Some data have been taken at large *q* mainly in connection with the determination of nuclear transparencies for high-energy protons.

The low-*q* data, taken with good energy and momentum resolution mainly at NIKHEF and Saclay, have been analyzed in the framework of a distorted wave Born approximation (DWBA) using optical potentials known from protoncarbon scattering. The QP radial wave functions have been parametrized using Woods-Saxon (WS) potentials. Lapikas *et al.* [7] have made a coherent analysis of the entire data set. The occupation of the QP orbits, obtained by summing the experimental spectroscopic factors, turns out to be rather low in comparison to what is known for other nuclei [1]; the summed 1*p* plus 1*s* strength amounts to 3.4 protons only (56% occupation).

The high-*q* data, determined in part with moderate energy

FIG. 2. Comparison of QP and correlated densities.

and momentum resolution, were taken mainly at SLAC and JLAB. The data were summed over a large region of initial momentum *k* and removal energy *E*, and fitted using WS radial wave functions and theoretical transparencies. When using the most reliable transparencies, confirmed by other experiments, Lapikas *et al.* found a much higher occupancy of the QP orbits, 5.0–5.6 protons (\sim 87% occupation).

This discrepancy—which is very embarrassing to the practitioners of $(e,e'p)$ as it sheds serious doubts on the quantitative interpretation of $(e, e[']p)$ data—obviously needs

FIG. 3. Comparison of QP and correlated densities. The crosses indicate the tail of the density obtained when using the low-*q* QP occupation.

to be better understood, and is discussed in more detail below. This difference also has led to speculations about *q*-dependent QP occupations, for which we see no physical basis.

A partial reason for the difference between the low-*q* and high-*q* results is immediately clear: The low-*q* data cover the region of missing momenta of typically $\langle 180 \text{ MeV}/c \rangle$ and missing energy $E \le 50$ MeV, the high-q data extend to 300 MeV/*c*. The high-*q* data also cover a larger range in missing energy; they are integrated up to typically 80 MeV. In this larger *k*, *E* range, there is not only QP strength, but also a fraction of the correlated strength is integrated over. Before making a valid comparison, this correlated strength needs to be removed.

In order to correct for this effect, we start from the high $q(e,e'p)$ data taken in a recent JLAB experiment [5] in quasielastic kinematics, which minimize final state interactions (FSI) and meson exchange current contributions (MEC). This experiment yields, in agreement with the previous JLAB and SLAC experiments, 5.2 protons in the integration region $E < 80$ MeV, $k < 300$ MeV/*c*. This number we correct for the continuum contribution using the calculated spectral function discussed above. With this correction the discrepancy between the low-*q* and high-*q* results is significantly reduced; the QP occupations now are 3.4 vs 4.5 protons for the low-*q*/high-*q* data, respectively. Given the uncertainties of these numbers—believed to be perhaps 10%—there still is a worrisome incompatibility.

In order to proceed, we have to choose. We have decided to use the occupation coming from the high-*q* measurements, as we judge the interpretation of these data to be safer. The low-*q* data suffer from uncertainties in the treatment of the final state interaction. Due to the low energies of the outgoing proton (70 MeV for the NIKHEF data [25]) coupledchannel effects not treated in the usual DWBA analysis should be relevant. van der Steenhoven *et al.* [25,29] have shown that inclusion of these effects would increase, for the rather soft nucleus ¹²C, the QP occupation by \sim 20%. For the kinematics of the low-*q* experiments, the calculations of Boffi *et al.* [30,31] also predict significant MEC effects that would lead to a further increase of the QP occupation.

The value for the QP occupation deduced from the high-*q* measurement is also compatible with the correlated strength measured directly in the recent experiment by Rohe [5]. This measurement agrees with theoretical predictions for the correlated strength of \sim 20%. The summed QP strength from the low- q (e,e'p) data (which thus includes the fragmentation due to long-range correlations), on the other hand, would correspond to greater than 40% correlated strength, i.e., be unrealistically high. Furthermore, as we will see below, our choice of the QP occupation is confirmed by a consistency check in our analysis.

Before proceeding to the calculation of the QP density, one more effect of correlations must be removed from the results of Lapikas *et al.* While the *p*- strength is located in discrete states where no ambiguity occurs, the *s* strength is located in the continuum between 20 and \sim 50 MeV removal energy. In this region, also the correlated strength contributes, and affects the *shape* of the fitted WS momentum distribution. The correlated strength has a momentum distribution that falls much more slowly with increasing *k* than the 1*s* QP strength; when fitting the sum with a WS parametrization, the resulting WS momentum distribution extends somewhat too far in momentum, i.e., it would have too small a radial extension in *r* space.

We have used the theoretical 1*s* QP and correlated strength in the region used in [7] for the determination of the 1*s* momentum distribution to calculate a correction to the WS parametrization fitted to the sum. We find that the extent of the wave function in *k* space needs to be reduced by 11%. This modified WS shape has been confirmed [32] by an independent analysis employing the recent JLAB data of [5] and the correlated spectral function of [3].

With these QP radial wave functions, and occupations renormalized to the one derived above from the high-*q* experiments, we can compute the QP density in *r* space. As the radial wave functions fitted by Lapikas *et al.* refer to relative coordinate between the proton and the c.m. of the $(A-1)$ system, we need to rescale the radial coordinate by a factor $11/12$, with the corresponding adjustment in height to conserve the normalization.

The resulting QP density is shown in Figs. 2 and 3. The difference of the 12 C point density and this QP density yields the correlated density also shown in Figs. 2 and 3. The "error bar" on the correlated density is not straightforward to calculate due to the various adjustments that had to made in the analysis; $\pm 20\%$ for $r < 3$ fm is probably a realistic estimate.

The first observation one can make from Figs. 2 and 3 concerns the fact that, as in the theoretical analysis of Fig. 1, the correlated density deduced from experimental data is significantly more concentrated towards the nuclear interior than the QP density. One also observes that the correlated density in the nuclear interior gives a very significant contribution, of order 30%, of the central density, i.e., larger than one could have expected from the number of 20% or so of correlated nucleons. This explains why attempts to explain the total densities in terms of QP orbitals cannot be very successful.

Figure 3 shows another important feature: The experimental QP density at large *r* agrees perfectly with the point density measured via elastic electron scattering. Such an agreement should occur, as at large *r*—outside the range of the nuclear potential—the density is entirely given by the tail of the least-bound QP orbit, the $1p_{3/2}$ state in ¹²C. More deeply bound states, or correlated nucleons with large removal energy, cannot contribute.

The good agreement between QP density and point density at large *r* also confirms the correctness of our choice of QP occupations. Had we used the occupation derived from the low-*q* experiments and, as above, the shape of $R(r)$ determined in [7] from the fit to the world $(e, e[']p)$ data, we would have obtained the tail indicated in Fig. 3 by the crosses. These are obviously significantly too low. This comparison thus provides an *a posteriori* justification of our procedure.

In Figs. 2 and 3 we also show the correlated density obtained by theory. Considering the above-mentioned uncertainty of the experimental result, we consider the agreement between theory and experiment as a good one. The size of the correlated contribution in the nuclear interior is very similar, the rapid falloff of the correlated density at large *r* also agrees within our estimated uncertainty. The correlated density deduced from experimental data seems to contain more spectral strength in partial waves with $l > 0$ than the theoretical one.

The large contribution of the correlated density in the nuclear interior shows that the neglect of this correlated contribution in the standard IP calculations (e.g., all the shellmodel descriptions) is not justified.

Starting from the $(e, e^t p)$ data, we have constructed the QP density for ${}^{12}C$ in coordinate space. The difference to the total density, obtained from elastic electron scattering, provides the density distribution of the correlated nucleons. We find that it is significantly more concentrated towards the nuclear interior. We also find good agreement with the theoretical calculation of the correlated density distribution.

The large contribution of the density related to shortrange *N-N* correlations, \sim 30% in the nuclear interior, together with the fact that the *shape* of the correlated density differs strongly from the QP density, explains the poor performance of QP wave functions in explaining many observables. Due to the shape difference, the shortcoming of the neglect of the correlated contribution also cannot satisfactorily be "compensated" by using effective quantities like effective charges, etc.

As a side product, our analysis provides a solution to the puzzle raised in [7], the pronounced disagreement between QP occupations derived from the low-*q* and high-*q* $(e, e'p)$ experiments. We find that only the high-*q* occupation (suitably corrected for the correlated contribution not considered in [7]), is compatible with the independent information from elastic electron scattering.

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