

**NUCLEAR CHARGE-DENSITY-DISTRIBUTION PARAMETERS
FROM ELASTIC ELECTRON SCATTERING**

H. DE VRIES, C. W. DE JAGER, and C. DE VRIES

Nationaal Instituut voor Kernfysica en Hoge-Energiefysica, sectie-K (NIKHEF-K)
P.O. Box 4395, 1009 AJ Amsterdam, The Netherlands

A compilation of nuclear charge-density-distribution parameters, obtained from elastic electron scattering, is presented in five separate tables. Data on charge distributions obtained on the basis of a phenomenological model—parameters of nuclei and differences therein between isotopes and between other neighboring nuclei like isotones—are given in Tables I, II, and III. Parameters obtained by a model-independent analysis are given in two additional tables: Table IV gives the coefficients of a Fourier-Bessel series expansion, and Table V gives the positions and amplitudes for the expansion in a sum of Gaussians. References through February 1986 have been covered. © 1987 Academic Press, Inc.

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INTRODUCTION

Since our previous compilation of charge-density-distribution parameters,¹ the analysis of electron scattering experiments has improved greatly. Ambiguities about the root-mean-square (rms) radius of the reference nucleus ¹²C have been resolved,² and the model-independent analysis of data has now generally been adopted. The greater consistency of the present data allows in many cases a combined analysis of experimental data from several electron scattering experiments.

The wealth of new data forced us to change the 1974 policy, where all available data were presented. In the present compilation the number of entries per isotope is limited to at most three. If a selection is necessary, in general the most recent values are given. In those cases where large discrepancies exist between the results of different experiments, we have chosen the values which we judge to be the most reliable. On the other hand we have tried to remain complete in a comprehensive list of all experimental papers.

There were several reasons for producing this compilation. Besides the already mentioned wealth of new and accurate data, large improvements have been made on the theoretical side: shell model calculations in the

lower part of the periodic table³ and density-dependent Hartree-Fock calculations⁴ for medium-heavy and heavy nuclei can now predict charge distributions with unprecedented accuracy. Comparison of the theoretical predictions with the new electron scattering results might guide theoreticians to further advances toward a better understanding of nuclear structure. Furthermore, the previous compilation has frequently been used in systematic studies of charge-density distributions.⁵⁻⁷ For future studies in this area, an updated version of the tables is crucial.

The material is presented in five tables. Table I covers standard charge-density-distribution parameters of nuclei. The next two tables deal with parameter differences between isotopes (Table II) and with differences between other neighboring nuclei (Table III). Results derived from model-independent analyses are presented in the two remaining tables: Table IV gives the coefficients obtained from a Fourier-Bessel analysis (FB), Table V lists the parameters from an analysis with a Sum-of-Gaussians (SOG) method. The extent of these last two tables confirms the expectation we expressed in 1974, namely that model-independent analysis would become more frequently used than analysis with phenomenological models.

In this compilation we no longer include the parameters of magnetization density distributions. The reason for this is twofold: first, Donnelly and Sick published an excellent review on magnetic elastic electron scattering about a year ago;⁸ and second, the presentation of the magnetization density is not so unambiguous as the description of the charge density. The latter is due to the fact that charge scattering is determined mainly by the monopole form factor (for $J = 0^+$ nuclei even exclusively), whereas for most nuclei with a magnetic moment, many different magnetic multipole components (which cannot be separated experimentally) contribute to the magnetic form factor.

This compilation was completed in February 1986 and has taken into consideration all data published up to that date.

Charge-Density-Distribution Parameters

For more extensive information the reader should consult monographs^{9,10} and the introduction to our 1974 compilation.¹ Here we limit ourselves to the following short notes.

The analysis of electron scattering data is restricted by the fact that the form factor can be studied only over a finite range of momentum transfers q_{\min} to q_{\max} . In Plane-Wave Born Approximation, the charge distribution $\rho(r)$ is the Fourier transform of the form factor $F(q)$ and for a spherically symmetric charge distribution is given by

$$\rho(r) = \frac{1}{2\pi^2} \int F(q) \frac{\sin(qr)}{qr} q^2 dq.$$

As a consequence, only the amplitudes of Fourier components of $\rho(r)$ with wavelengths between $2\pi/q_{\max}$ and $2\pi/q_{\min}$ can be extracted from the data. In the past, the limited range of q values made it necessary to describe the data on the basis of a phenomenological model. For experiments which are performed at relatively low q values the data can be described adequately by a two-parameter Fermi distribution. Extension of this model with more parameters showed, however, the limitation of the description: the introduction of a "wine-bottle" parameter w is generally speaking more representative for the behavior of the tail of the charge-density distribution than for that of the inner region.¹¹ This is illustrated by the fact that analyses of the same data set with a three-parameter Fermi and with a three-parameter Gaussian model often yield w values with opposite signs.¹² Also, the inclusion of oscillatory components in $\rho(r)$, which were introduced to fit the data measured at high q values, are at best only significant for light- and medium-heavy nuclei. Even the error bars quoted for the rms radius of $\rho(r)$ do not necessarily represent the full range of rms radii consistent with the experimental data. These limitations can be re-

moved only through the use of a model-independent analysis (see below).

One of the major advantages of electron scattering as a nuclear probe is the fact that the interaction is purely electromagnetic and hence is well known. This implies that for a given charge distribution, electron scattering cross sections can be calculated accurately by phase-shift analysis. Two theoretical problems remain to be solved in the field of electron scattering: radiative and dispersion corrections. Radiative corrections are in good agreement with experiments involving relative measurements, but deviations have been observed in absolute measurements. This seems to indicate that higher-order diagrams have to be taken into account. Dispersive effects or virtual nuclear excitations might also play a role. No accurate calculations are available as yet, but there is experimental evidence that the minimum of the elastic form factor might be appreciably affected. More will be said about this subject in the paragraph where electron scattering results are compared with muonic x-ray data.

Model-Independent Analysis

The higher accuracy of the experimental data and the larger q range covered have led to the use of more refined models to describe the finer details of the charge-density distribution. However, the interpretation of the results has not always been carried out unambiguously. Several attempts have been made to describe the charge distribution by sets of orthonormal functions. A viable model-independent analysis will have to incorporate some model dependence to account for the fact that data are available only over a finite q range. The limitation should be based on physical arguments. At present the majority of experimental results are analyzed by two different model-independent approaches: Fourier-Bessel analysis or sum of Gaussians.

Fourier-Bessel Analysis

The Fourier-Bessel series expansion was introduced by Dreher et al.¹³ For practical reasons $\rho(r)$ is assumed to be zero beyond a certain cutoff radius R . The first N ($=Rq_{\max}/\pi$) coefficients of this series expansion are determined directly from the experimental data. The behavior of the form factor $F(q)$, at q values beyond the maximum value of q for which data are available, is assumed to be limited by a q^{-4} and an $\exp(-aq^2)$ decrease. These assumptions originate from expectations for the distribution of the nucleons inside the nucleus and for the finite extension of the nucleons, respectively. They yield an upper limit for the contributions of the higher Fourier components of the series expansion. The results depend to a certain degree on the value of the cutoff radius R . An advantage of this method is that the uncertainties

in the charge distribution originating from the experimental errors and from lack of knowledge about the large- q behavior can be determined separately.

Unfortunately, several definitions and normalizations have been used in the literature. In this compilation, we use

$$\rho(r) = \begin{cases} \sum_v a_{v,0} j_0(v\pi r/R) & \text{for } r \leq R \\ 0 & \text{for } r \geq R, \end{cases}$$

where $j_0(qr)$ denotes the Bessel function of order zero.

For the normalization we have adopted the convention that the integral over the charge distribution equals the nuclear charge Ze . This normalization has the advantage that the difference between different nuclei can be deduced directly from the difference between the Fourier-Bessel coefficients

$$\Delta\rho(r) = \sum_v \delta a_{v,\Delta A} j_0(v\pi r/R) \quad \text{for } r \leq R$$

with

$$\delta a_{v,\Delta A} = a_{v,A1} - a_{v,A2}$$

provided that both sets of coefficients have been determined for the same value of the cutoff radius R .

In Table IV we present the Fourier-Bessel coefficients with the above-mentioned definition and the value of the cutoff radius that has been used. Since the analysis frequently involves several data sets, all of these are also indicated. A final remark should be made about the errors in the coefficients, which are not presented in the tables. Since the errors are strongly correlated, the uncertainties in the charge distribution can be determined only from the full correlation matrix. But since this matrix is never published, it would not make sense to present the errors in the Fourier-Bessel coefficients.

Sum of Gaussians

This parametrization was first introduced by Sick.¹¹ The width γ of the Gaussians is chosen equal to the smallest width of the peaks in the nuclear radial wave functions calculated by the Hartree-Fock method. Only positive values of the amplitudes of the Gaussians are allowed so that no structures narrower than γ can be created through interference. An advantage of the use of Gaussians is that values of $\rho(r)$ at different values of r are decoupled to a large extent because of the rapid decrease of the Gaussian tail. The results of the analysis are independent of the number of Gaussians, provided this number is sufficiently large to allow a good fit to the data. In this approach, all authors use the same definition,

$$\rho(r) = \sum_i A_i \{ \exp(-[(r - R_i)/\gamma]^2) + \exp(-[(r + R_i)/\gamma]^2) \},$$

where the coefficients A_i are given by

$$A_i = ZeQ_i/[2\pi^{3/2}\gamma^3(1 + 2R_i^2/\gamma^2)].$$

In this definition the values of Q_i indicate the fraction of the charge contained in the i th Gaussian, normalized such that

$$\sum_i Q_i = 1.$$

Table V gives a list of the positions R_i and amplitudes Q_i fitted to the data. The rms radius of the Gaussians and the rms radius of the charge distribution deduced are also given. The data sets used in the analysis are mentioned.

Electron Scattering and Muonic x-Rays

The information on nuclear charge-density distributions can be improved significantly by simultaneous analyses of electron scattering data and muonic x-ray data. Whereas electron scattering maps the Fourier transform of the charge-density distribution, muonic transition energies are sensitive to a special moment of this distribution, the so-called Barrett moment,¹⁴

$$\langle r^k e^{-\alpha r} \rangle = (4\pi/Ze) \int \rho(r) r^k e^{-\alpha r} r^2 dr.$$

The parameters k and α are discussed in Ref. 15.

The inclusion of these Barrett moments in the electron scattering analysis reduces the uncertainty in the lowest-order Fourier-Bessel coefficients. Effectively, the muonic information is equivalent to an electron scattering experiment at low momentum transfer. Inclusion of the precisely known value of the muonic Barrett moment greatly improves the overall normalization error as well, resulting in a substantial reduction of the uncertainties in the combined analysis.

One would expect that the inclusion of muonic data in an electron scattering analysis would only reduce the errors. However, a comparison of results from only electron scattering and only muonic x-ray data shows that the muonic results yield rms radii up to 0.02 fm larger than those deduced from electron scattering.^{16,17} Therefore we have indicated it in the tables when muonic results have been included in the analysis.

There are several possible explanations for this discrepancy. Whereas on the muonic side the calculation of the nuclear polarization correction is supposedly well under control, a remanent discrepancy might be present due to a short-range muon-nucleon interaction described by a scalar or a vector boson.¹⁶ Another explanation might be that the electron scattering results are not corrected for dispersive effects. Rough estimates¹⁸ predict the dispersive corrections to be small. Recent experiments,¹⁹ however, seem to indicate that the dispersive effects might be ap-

preciable. More conclusive experiments are necessary to settle this problem. If these effects are studied in more detail and the cause of the discrepancies mentioned can be resolved, one may expect still further improvement in our knowledge about charge distributions of nuclei, which is already impressively accurate.

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POLICIES

<i>Literature coverage</i>	All available experimental papers were covered, including preprints, theses, and internal reports. If the same data have been described in several papers, we have given only the most extensive and easiest-to-access reference. No theoretical papers have been included, unless they contain a reanalysis of experimental data.
<i>Tabulated results</i>	The maximum number of entries per isotope is limited to three. If more results were available we have in general listed the most recent values. In cases where conflicting results were present we have listed those which we considered to be most reliable. References containing additional information are given in a separate list at the end of the corresponding table.
<i>Models</i>	In those cases where the same data have been analyzed with different models, only results obtained with one particular model have been tabulated, following wherever possible the preference given in the original publication. Unless otherwise stated, the results are for the monopole charge distribution (C_0) only.
<i>Errors</i>	The errors quoted have been taken from the original papers. Generally, they represent the total error: the sum of the statistical (one standard deviation) and the systematic errors. No effort has been made to standardize the various types of error analyses used. The errors are given in parentheses following a tabulated value. For example, $0.3359(36) = 0.3359 \pm 0.0036$.
<i>References</i>	A comprehensive list of all experimental papers is given at the end of the five Tables. Reference keys are in the style Si79, where Si refers to the name of the first author and 79 to the year. In cases where this key is ambiguous a letter is added, as in Ca82a.
<i>Neighboring nuclei</i>	By observing the cross-section ratios, differences in charge-distribution parameters between neighboring nuclei can be determined to a higher accuracy than the parameters themselves. Therefore, the results of such analyses have been listed separately in Tables II and III. Not included there are the results of experiments where neighboring nuclei have been measured simultaneously, but not analyzed in terms of cross-section ratios. On the other hand, charge-distribution parameters which have been obtained solely through cross-section ratios of neighboring nuclei are omitted from Table I.

EXPLANATION OF TABLES

TABLE I. Charge-Density-Distribution Parameters**TABLE II. Differences in Charge-Density-Distribution Parameters between Isotopes****TABLE III. Differences in Charge-Density-Distribution Parameters between Neighboring Nuclei (Not Isotopes)**

Nucleus The absence of a mass number indicates that the tabulated values are for a natural isotopic admixture. The asterisk after a nucleus indicates that additional information is available in the references given at the end of the tabulation.

model The normalization of the charge distribution is such that

$$4\pi \int \rho(r)r^2 dr = Ze.$$

If no entry is given in this column, the model used is described in the "remarks."

EXPLANATION OF TABLES continued

HO Harmonic-oscillator model:

$$\begin{aligned}\rho(r) &= \rho_0(1 + \alpha(r/a)^2)\exp(-(r/a)^2) \\ \alpha &= \alpha_0 a_0^2 / (a^2 + \frac{3}{2} \alpha_0(a^2 - a_0^2)) \\ a_0^2 &= (a^2 - a_p^2)A/(A - 1) \\ \alpha_0 &= (Z - 2)/3; \quad a_p^2 = \frac{2}{3} \langle r^2 \rangle_{\text{proton}}\end{aligned}$$

MHO Modified harmonic-oscillator model, with the same expression for $\rho(r)$ as in HO but with α as an additional free parameter

MI Model-independent evaluation of the form factor by the expression

$$F(q^2) = 1 - \frac{1}{6} q^2 \langle r^2 \rangle$$

FB Model-independent analysis by means of a Fourier-Bessel expansion for the charge distribution

SOG Model-independent analysis by means of an expansion for the charge distribution as a sum of Gaussians

2pF Two-parameter Fermi model

$$\rho(r) = \rho_0 / (1 + \exp((r - c)/z))$$

3pF Three-parameter Fermi model

$$\rho(r) = \rho_0(1 + wr^2/c^2)/(1 + \exp((r - c)/z))$$

3pG Three-parameter Gaussian model

$$\rho(r) = \rho_0(1 + wr^2/c^2)/(1 + \exp((r^2 - c^2)/z^2))$$

UG Uniform Gaussian model

$$\rho(r) = \rho_0 \int \exp(-(r-x)^2/g^2)x^2 dx$$

$\langle r^2 \rangle^{1/2}$ Root-mean-square radius of the charge distribution

$$\langle r^2 \rangle = (4\pi/Ze) \int \rho(r)r^4 dr$$

c or a In this column the values are given for the parameter c if the 2pF, 3pF, or 3pG model has been used and for the parameter a if the HO or MHO model has been used.

z or α In this column the values are given for the parameter z if the 2pF, 3pF, or 3pG model has been used and for the parameter α if the HO or MHO model has been used.

w The parameter w of the 3pF and 3pG models

q-range The momentum transfer range covered by the data used in the analysis

ref. Source of tabulated data, keyed to the list of references following the tables

remarks The symbols \dagger and $\$$ denote that additional information can be found in Tables IV (Fourier-Bessel) and V (Sum-of-Gaussians), respectively. The entries indicated by a letter or number are explained at the end of each table.

$\Delta \langle r^2 \rangle^{1/2}$ $\langle r^2 \rangle^{1/2}(A_2) - \langle r^2 \rangle^{1/2}(A_1)$ with $A_2 > A_1$.

Δc $c(A_2) - c(A_1)$ with $A_2 > A_1$.

Δz $z(A_2) - z(A_1)$ with $A_2 > A_1$.

Δw $w(A_2) - w(A_1)$ with $A_2 > A_1$.

EXPLANATION OF TABLES continued

TABLE IV. Fourier-Bessel Coefficients

rms	Value of the root-mean-square radius $\langle r^2 \rangle^{1/2}$ of the charge distribution
a1 ... a17	List of the Fourier-Bessel coefficients a_v , with $v = 1$ to 17. The coefficients are defined by
	$\rho(r) = \sum_v a_v j_0(v\pi r/R) \quad \text{for } r \leq R,$
	$\rho(r) = 0 \quad \text{for } r > R.$
	The normalization is chosen such that $4\pi \int \rho(r)r^2 dr = Ze$. 0.25182e -1 means 0.25182×10^{-1} .
ref.	Reference for the data analysis
q-range	The momentum-transfer range covered by the data used in the analysis
data-sets	References for the data sets used in the analysis. The symbol μ indicates that muonic x-ray data have been used as a constraint in the analysis.
R	Value of the cutoff radius, beyond which the charge density is assumed to be identical to zero

TABLE V. Sum-of-Gaussians Parameters

rms	Value of the root-mean-square radius $\langle r^2 \rangle^{1/2}$ of the charge distribution
R_i, Q_i	Position and amplitude of the Gaussians. The coefficients are defined by
	$\rho(r) = \sum_i A_i \{ \exp(-[(r - R_i)/\gamma]^2) + \exp(-[(r + R_i)/\gamma]^2) \}$
	with
	$A_i = ZeQ_i/[2\pi^{3/2}\gamma^3(1 + 2R_i^2/\gamma^2)].$
	The values of Q_i indicate the fraction of the charge contained in the i th Gaussian, normalized such that $\sum_i Q_i = 1$.
ref.	Reference for the data analysis
q-range	The momentum-transfer range covered by the data used in the analysis
data-sets	References for the data sets used in the analysis. The symbol μ indicates that muonic x-ray data have been used as a constraint in the analysis.
RP	The rms radius of the Gaussians: $RP = \gamma \sqrt{\frac{3}{2}}$

TABLE I. Charge-Density-Distribution Parameters
See page 500 for Explanation of Tables

Nucleus	model	$\langle r^2 \rangle^{1/2}$ [fm]	c or a [fm]	z or α [fm]	w	q-range [fm $^{-1}$]	ref.	remarks
n*		0.3359(36) 0.3455(26)			0 0		Kr73 Ko76	1 1
${}^1\text{H}^*$	MI	0.85(2)			0.33 - 1.42		Th72	a,2
	MI	0.84(1)			0.36 - 11.50		Ho76	3
	MI	0.862(12)			0.36 - 1.18		Si80	4
${}^2\text{H}^*$	MI	2.095(6)			0.22 - 0.71		Be73a	a,h,5
	MI	2.116(6)			0.21 - 0.77		Si81	h,6
${}^3\text{H}^*$	FB	1.68(3)			0.51 - 2.83		Be84	\dagger ,b,7
	SOG	1.76(4)			0.55 - 4.79		Ju85	\$,8
${}^3\text{He}^*$	SOG	1.844(45)			0.59 - 4.47		MC77	\$
	FB	1.877(19)			0.18 - 10.1		Re84b	\dagger ,9
	MI	1.976(15)			0.45 - 1.92		Ot85	a,h,10
${}^4\text{He}^*$	MI	1.696(14)			0.17 - 0.50		Gu82	h
	SOG	1.676(8)			0.14 - 7.70		Si82	\$,g,h,11
	MI	1.671(14)			0.51 - 2.00		Ot85	a,h,12
${}^6\text{Li}^*$	MHO	2.54(5)			0.69 - 2.52		Su67	a,h,13
	MHO	2.56(5)			0.56 - 3.66		Li71a	h,14
	MI	2.57(10)			0.09 - 0.90		Bu72	i,15
${}^7\text{Li}^*$	HO	2.39(3)	1.77(2)	0.327	0.69 - 2.62		Su67	a,h,16
	MI	2.41(10)			0.09 - 0.90		Bu72	i,15
${}^9\text{Be}^*$	HO	2.519(12)	1.791(9)	0.611	0.26 - 0.70		Ja72	
	HO	2.50(9)	1.77(6)	0.631	0.15 - 0.53		Fe73a	i,17
${}^{10}\text{B}^*$	HO	2.45(12)	1.71(8)	0.837	0.69 - 2.81		St66b	b
${}^{11}\text{B}^*$	HO	2.42(12)	1.69(8)	0.811	0.69 - 2.81		St66b	b
		2.37			0.61 - 1.76		Ri71	b,18
${}^{12}\text{C}^*$	FB	2.472(15)			0.10 - 4.01		Ca80a	\dagger ,19
	SOG	2.471(6)			0.13 - 3.70		Si82	\$,g,h,20
	FB	2.464(12)			0.25 - 2.75		Re82	\dagger
${}^{13}\text{C}$	MHO	2.440(25)	1.635(9)	1.403(16)	0.3 - 3.4		He70a	h,21
${}^{14}\text{C}$	MHO	2.56(5)	1.73(4)	1.38(12)	1.04 - 2.16		KI73	h
${}^{14}\text{N}^*$	HO	2.58	1.76	1.234	0.86 - 1.62		Da70	a,22
	HO	2.540(20)	1.729(14)	1.291	0.29 - 0.48		Sc75	h
	3pF	2.524(23)	2.570(8)	0.5052(2)	-0.180(7)		La82	i,23
${}^{15}\text{N}^*$	MHO	2.65	1.81	1.250	0.86 - 1.62		Da70	a,h,22
	3pF	2.70(3)	2.334(35)	0.498(5)	0.139(30)		Ge72	24
	FB	2.611(9)			0.22 - 3.17		Vr86	\dagger ,h,25
${}^{16}\text{O}^*$	3pF	2.730(25)	2.608(36)	0.513(5)	1.05 - 3.97		Si70b	\$,h,26
	HO	2.718(21)	1.833(14)	1.544	0.29 - 0.48		Sc75	h
	FB	2.737(8)			0.38 - 2.85		La82	\dagger ,i

TABLE I. Charge-Density-Distribution Parameters
See page 500 for Explanation of Tables

Nucleus	model	$\langle r^2 \rangle^{1/2}$ [fm]	c or a [fm]	z or α [fm]	w	q-range [fm $^{-1}$]	ref.	remarks
^{17}O	HO	2.662(26)	1.798(18)	1.498		0.58 - 0.99	Si70a	27
^{18}O	HO	2.727(20)	1.841(14)	1.513		0.58 - 0.99	Si70a	27
$^{19}\text{F}^*$	2pF	2.900(15)	2.59(4)	0.564(14)		0.55 - 1.01	Ha73b	28
	2pF	2.90(2)	2.58(4)	0.567(13)		0.46 - 1.79	Oy75	
$^{20}\text{Ne}^*$	2pF	3.040(25)	2.805(15)	0.571(5)		0.22 - 1.04	Mo71	27
	2pF	3.004(25)	2.740(46)	0.572(17)		0.21 - 1.12	Kn81	i
	3pF	2.992(8)	2.791(9)	0.698(5)	-0.168(8)	0.49 - 1.80	Be85	d
^{22}Ne	2pF	2.969(21)	2.782(12)	0.549(4)		0.2 - 1.1	Mo71	27
^{23}Na	UG	2.94(6)				0.4 - 2.02	Sa69a	b,29
$^{24}\text{Mg}^*$	3pF	3.075(15)	3.108(33)	0.607(9)	-0.163(30)	0.58 - 1.99	Av74	i
	3pF	2.985(30)	3.192(34)	0.604(6)	-0.249(20)	0.74 - 3.46	Li74	\$,h,30
	2pF	3.08(5)	2.98(5)	0.551(32)		0.20 - 1.15	Le76	i
^{25}Mg	2pF	3.11(5)	2.76(5)	0.608(32)		0.20 - 1.15	Le76	i
	3pF	3.003(11)	3.22(5)	0.58(4)	-0.236(34)	0.19 - 2.07	Eu77a	i,31
^{26}Mg	2pF	3.06(5)	3.05(5)	0.523(32)		0.20 - 1.15	Le76	i
$^{27}\text{Al}^*$	2pF	3.06(9)	3.07(9)	0.519(26)		0.51 - 1.59	Lo67	
	2pF	3.05(5)	2.84(5)	0.569		0.23 - 0.59	Fe73a	b,e,i,32
	FB	3.035(2)				0.47 - 2.70	Ro86	
$^{28}\text{Si}^*$	3pG	3.106(30)	1.95(9)	2.076(10)	0.286(12)	0.74 - 3.71	Li74	\$,h
	2pF	3.15(4)	3.14(6)	0.537(32)		0.16 - 1.1	Br77	i
	3pF	3.086(18)	3.340(9)	0.580(3)	-0.233(9)	0.25 - 1.49	Mi82	†,i,33
^{29}Si	2pF	3.13(5)	3.17(8)	0.52(4)		0.16 - 1.1	Br77	i
	3pF	3.079(21)	3.338(12)	0.547(2)	-0.203(12)	0.25 - 1.49	Mi82	†,i,33
^{30}Si	3pF	3.176(22)	3.252(21)	0.553(2)	-0.078(22)	0.25 - 1.49	Mi82	†,i,33
$^{31}\text{P}^*$	3pF	3.19(3)	3.369(25)	0.582(6)	-0.173(24)	0.73 - 2.85	Si72	h,34
	FB	3.187(11)				0.30 - 2.85	Me76	i,35
	FB	3.187				0.25 - 2.64	Mi82	†,i,36
$^{32}\text{S}^*$	3pG	3.239(30)	2.54(9)	2.191(10)	0.160(12)	0.74 - 3.71	Li74	\$,f,h,30
	FB	3.240(11)				0.30 - 3.71	Me76	f,1,37
	FB	3.248(11)				0.47 - 2.56	Ry83a	†
$^{34}\text{S}^*$	FB	3.281(13)				0.47 - 2.56	Ry83a	†
$^{36}\text{S}^*$	FB	3.278(11)				0.47 - 2.56	Ry83a	†
^{35}Cl	3pF	3.388(17)	3.476(32)	0.599(5)	-0.10(2)	0.60 - 1.70	Br80	38
^{37}Cl	3pF	3.384(17)	3.554(27)	0.588(5)	-0.13(2)	0.60 - 1.70	Br80	39

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Nucleus	model	$\langle r^2 \rangle^{1/2}$ [fm]	c or a [fm]	z or α [fm]	w	q-range [fm $^{-1}$]	ref.	remarks
^{36}Ar	2pF	3.327(15)	3.54(4)	0.507(15)		0.54 - 0.96	Fi76	d,i
$^{40}\text{Ar}^*$	3pF	3.48(4)	3.73(5)	0.62(1)	-0.19(4)	0.8 - 1.81	We74	i
	2pF	3.393(15)	3.53(4)	0.542(15)		0.54 - 0.96	Fi76	d,i
	FB	3.423(14)				0.29 - 1.81	Ot82	$\dagger,h,40$
$^{39}\text{K}^*$	UG	3.40(7)				0.4 - 1.92	Sa69a	b,f,41
	3pF	3.408(27)	3.743(25)	0.585(6)	-0.201(22)	0.64 - 3.43	Si73a	$\$,f,h,42$
$^{40}\text{Ca}^*$	3pF	3.482(25)	3.766(23)	0.586(5)	-0.161(23)	0.53 - 3.24	Si73a	h,43
	SOG	3.479(3)				0.53 - 3.69	Si79	$\$,g,44$
	FB	3.450(10)				0.35 - 3.69	Em83b	$\dagger,45$
$^{48}\text{Ca}^*$	3pF	3.470	3.7369	0.5245	-0.030	0.49 - 3.37	Be67d	$\$,h,46$
	FB	3.451(9)				0.35 - 3.55	Em83b	\dagger
$^{48}\text{Ti}^*$	2pF	3.713(15)	3.843(15)	0.588(5)		1.03 - 1.62	Sh78c	i
	FB	3.597(1)				0.61 - 2.20	Se85	\dagger,i
$^{50}\text{Ti}^*$	FB	3.573(2)				0.61 - 2.20	Se85	\dagger,i
$^{51}\text{V}^*$	2pF	3.58(4)	3.94(3)	0.505(14)		0.56 - 1.79	Pe73	i
	2pF	3.615(31)	3.91(4)	0.532(29)		0.23 - 0.78	Go74	i
$^{50}\text{Cr}^*$	2pF	3.638(13)	3.979(30)	0.520(13)		0.15 - 0.79	La76	i
	2pF	3.707(15)	3.941(15)	0.566(5)		0.97 - 1.62	Sh78b	i
	FB	3.662(4)				0.15 - 2.59	Li83	\dagger,g
$^{52}\text{Cr}^*$	2pF	3.613(17)	4.01(4)	0.497(19)		0.15 - 0.79	La76	i
	2pF	3.684(15)	3.984(15)	0.542(5)		0.97 - 1.62	Sh78b	i
	FB	3.643(3)				0.15 - 2.59	Li83	\dagger,g
^{53}Cr	2pF	3.726(15)	4.000(15)	0.557(5)		0.97 - 1.62	Sh78b	i
^{54}Cr	2pF	3.673(14)	4.021(31)	0.524(13)		0.15 - 0.79	La76	i
	2pF	3.776(15)	4.010(15)	0.578(5)		0.97 - 1.62	Sh78b	i
	FB	3.689(4)				0.15 - 2.59	Li83	\dagger,g
^{55}Mn	2pF	3.68(11)	3.89(12)	0.567		0.23 - 0.50	Th69	e,47
$^{54}\text{Fe}^*$	3pG	3.680(13)	3.518(17)	2.270(12)	0.403(15)	0.51 - 2.22	Wo76	i
	2pF	3.675(17)	4.075(38)	0.506(18)		0.15 - 0.79	La76	i
	2pF	3.732(15)	4.074(14)	0.536(6)		0.97 - 1.62	Sh78b	i
$^{56}\text{Fe}^*$	3pG	3.729(13)	3.505(17)	2.325(12)	0.380(15)	0.51 - 2.22	Wo76	i
	2pF	3.721(20)	4.106(45)	0.519(21)		0.15 - 0.79	La76	i
	2pF	3.801(15)	4.111(13)	0.558(6)		0.97 - 1.62	Sh78b	i
^{58}Fe	2pF	3.783(19)	4.027(19)	0.576(7)		1.02 - 1.77	Li71b	c
	3pG	3.767(13)	3.585(18)	2.354(12)	0.328(15)	0.51 - 2.22	Wo76	i
$^{59}\text{Co}^*$	2pF	3.80(5)	4.08(5)	0.569		0.23 - 0.56	Fe73a	e,48
	3pG	3.775(37)	3.656(18)	2.339(12)	0.0331(16)	0.51 - 2.22	Sc77	\dagger,g,i
	2pF	3.864(15)	4.158(20)	0.575(5)		0.96 - 1.61	Sh78a	

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Nucleus	model	$\langle r^2 \rangle^{1/2}$ [fm]	c or a [fm]	z or α [fm]	w	q-range [fm $^{-1}$]	ref.	remarks
$^{58}\text{Ni}^*$	3pF	3.764(10)	4.3092	0.5169	-0.1308	0.58 - 2.64	Fi70	g,h,49
	SOG	3.772(2)				0.58 - 3.80	Ca80b	\$,g,50
	FB	3.769(13)				0.25 - 2.65	Be83	†
$^{60}\text{Ni}^*$	3pF	3.796(10)	4.4891	0.5369	-0.2668	0.52 - 2.28	Fi70	g,h,49
	3pG	3.793(13)		3.691(16)	2.337(12)	0.352(15)	Wo76	†,i
	FB	3.797(13)				0.25 - 2.65	Be83	†
^{61}Ni	3pF	3.806(10)	4.4024	0.5401	-0.1983	0.52 - 2.28	Fi70	g,h,49
$^{62}\text{Ni}^*$	3pF	3.822(10)	4.4425	0.5386	-0.2090	0.52 - 2.28	Fi70	g,h,49
	3pG	3.830(13)		3.742(17)	2.360(12)	0.338(16)	Wo76	i
	2pF	3.827(13)		4.262(26)	0.521(12)	0.15 - 0.79	Ke77	i
$^{64}\text{Ni}^*$	2pF	3.907(26)	4.212(28)	0.578(7)	-0.2284	0.68 - 1.61	Kh70b	c
	3pF	3.845(10)		4.5211		0.52 - 2.64	Fi70	g,h,49
	3pG	3.850(13)		3.842(16)		0.51 - 2.22	Wo76	i
$^{63}\text{Cu}^*$	3pG	3.876(36)	3.707(19)	2.412(12)	0.339(16)	0.51 - 2.22	Sc77	†,g,i
	2pF	3.933(15)		4.163(27)	0.606(11)	0.15 - 0.79	Ke77	i
	2pF	3.947(13)		4.218(14)	0.596(5)	0.96 - 1.61	Sh78a	
$^{65}\text{Cu}^*$	3pG	3.892(36)	3.807(19)	2.405(12)	0.315(16)	0.51 - 2.22	Sc77	†,g,i
	2pF	3.986(19)		4.158(35)	0.632(8)	0.15 - 0.79	Ke77	i
	2pF	3.954(13)		4.252(15)	0.589(5)	0.96 - 1.61	Sh78a	
$^{64}\text{Zn}^*$	2pF	3.965(17)	4.285(9)	0.584(9)	0.342(15)	0.30 - 1.09	Ne72	i,51
	3pG	3.923(13)		3.664(18)		0.51 - 2.22	Wo76	i
	2pF	3.923(13)		4.297(25)		0.15 - 0.79	Ke77	i
$^{66}\text{Zn}^*$	2pF	3.991(27)	4.286(29)	0.595(11)	0.299(15)	0.96 - 1.63	Li73	c,i
	3pG	3.931(13)		3.757(18)		0.51 - 2.22	Wo76	i
	2pF	3.952(15)		4.340(28)		0.15 - 0.79	Ke77	i
$^{68}\text{Zn}^*$	2pF	3.979(31)	4.353(32)	0.567(14)	0.293(16)	0.96 - 1.63	Li73	c,i
	3pG	3.963(13)		3.870(18)		0.51 - 2.22	Wo76	i
	2pF	3.958(17)		4.393(32)		0.15 - 0.79	Ke77	i
^{70}Zn	2pF	4.044(18)	4.409(10)	0.583(9)	0.342(18)	0.30 - 1.09	Ne72	i,51
	3pG	3.986(13)		3.878(20)		0.51 - 2.22	Wo76	i
	2pF	3.993(20)		4.426(37)		0.15 - 0.79	Ke77	i
^{70}Ge	2pF	4.07(2)	4.44(2)	0.585(7)		0.65 - 1.14	Kl75	i
	FB	4.043(2)				0.35 - 2.90	Ma84	†,g
^{72}Ge	2pF	4.05(3)	4.45(2)	0.573(7)		0.65 - 1.14	Kl75	i
	FB	4.060(2)				0.35 - 2.90	Ma84	†,g
^{74}Ge	FB	4.075(2)				0.35 - 2.90	Ma84	†,g
^{76}Ge	FB	4.081(2)				0.35 - 2.90	Ma84	†,g

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Nucleus	model	$\langle r^2 \rangle^{1/2}$ [fm]	c or a [fm]	z or α [fm]	w	q-range [fm $^{-1}$]	ref.	remarks
$^{88}\text{Sr}^*$	3pG	4.26(1)	4.254(10)	2.548(6)	0.47(3)	0.71 - 2.65	Al73	h
	2pF	4.17(2)	4.83(1)	0.496(11)		0.41 - 1.01	Fi74	i
	FB	4.188(5)				0.50 - 2.50	St76	$\dagger, 52$
$^{89}\text{Y}^*$	2pF	4.24	4.76(5)	0.571(29)		0.49 - 1.81	Sh67a	c
	3pG	4.24(2)	4.45(3)	2.526(23)	0.25	0.41 - 1.15	Si73c	d,i,53
	2pF	4.27(2)	4.86(1)	0.542(11)		0.41 - 1.01	Fi74	i
$^{90}\text{Zr}^*$	3pG	4.274(22)	4.434(20)	2.528(3)	0.350(25)	0.53 - 2.80	Fa71	h
	3pG	4.28(2)	4.46(5)	2.569(32)	0.25	0.40 - 1.15	Si73c	d,i,53
	FB	4.258(8)				0.50 - 2.50	Ro76	\dagger, i
^{91}Zr	3pG	4.309(22)	4.325(20)	2.581(3)	0.433(25)	0.53 - 2.43	Fa71	h
^{92}Zr	3pG	4.300(22)	4.455(20)	2.550(3)	0.334(25)	0.53 - 2.43	Fa71	h
	FB	4.294(11)				0.50 - 2.50	Ro76	\dagger, i
^{94}Zr	3pG	4.332(22)	4.494(20)	2.585(3)	0.296(25)	0.53 - 2.43	Fa71	h
	FB	4.315(10)				0.50 - 2.50	Ro76	\dagger, i
^{96}Zr	3pG	4.396(22)	4.503(20)	2.602(3)	0.341(25)	0.89 - 2.80	Fa71	h
^{93}Nb	2pF	4.31	4.87(5)	0.573(29)		0.49 - 1.81	Sh67a	c
	2pF	4.331(10)	4.953(6)	0.541(5)		0.15 - 0.79	Ja76	54
^{92}Mo	3pG	4.28(7)	4.61(10)	2.52(7)	0.19(11)	0.64 - 1.86	Ph72	i
	3pG	4.34(2)	4.56(4)	2.606(23)	0.21	0.40 - 1.15	Si73c	d,i,53
	FB	4.294(16)				0.56 - 1.96	Dr75	\dagger, i
^{94}Mo	FB	4.334(16)				0.56 - 1.96	Dr75	\dagger, i
^{96}Mo	FB	4.364(16)				0.56 - 1.96	Dr75	\dagger, i
^{98}Mo	FB	4.388(16)				0.56 - 1.96	Dr75	\dagger, i
^{100}Mo	FB	4.430(16)				0.56 - 1.96	Dr75	\dagger, i
^{104}Pd	FB	4.437(10)				0.29 - 2.38	La86	\dagger, i
^{106}Pd	FB	4.467(11)				0.29 - 2.38	La86	\dagger, i
^{108}Pd	FB	4.524(10)				0.29 - 2.38	La86	\dagger, i
^{110}Pd	2pF	4.639(19)	5.301(23)	0.581(9)		0.36 - 1.00	Li76	i
	FB	4.541(10)				0.29 - 2.40	La86	\dagger, i
^{110}Cd	2pF	4.578(7)	5.33(2)	0.535(7)		0.25 - 1.07	Gi75	i,55
^{112}Cd	2pF	4.608(7)	5.38(2)	0.532(9)		0.25 - 1.07	Gi75	i,55
^{114}Cd	2pF	4.629(8)	5.40(2)	0.537(9)		0.25 - 1.07	Gi75	i,55
	2pF	4.632(17)	5.314(23)	0.571(9)		0.36 - 1.00	Li76	i
^{116}Cd	2pF	4.639(8)	5.42(2)	0.532(9)		0.25 - 1.07	Gi75	i,55

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Nucleus	model	$\langle r^2 \rangle^{1/2}$ [fm]	c or a [fm]	z or α [fm]	w	q-range [fm $^{-1}$]	ref.	remarks
In*	2pF	4.646(12)	5.357(7)	0.563(4)		0.15 - 0.79	Ja76	56
^{112}Sn	2pF 3pG	4.655(23) 4.586(5)	5.375(26) 4.962(7)	0.560(10) 2.638(3)	0.285(12)	0.49 - 1.40 0.64 - 2.37	Kh70b Fi72	c h,57
^{114}Sn	3pG	4.602(5)	4.971(7)	2.636(3)	0.320(12)	0.64 - 2.37	Fi72	h,57
$^{116}\text{Sn}^*$	3pG 2pF SOG	4.619(5) 4.626(15) 4.627	5.062(7) 5.358(22)	2.625(3) 0.550(9)	0.272(12)	0.64 - 2.65 0.36 - 1.00 0.36 - 3.60	Fi72 Li76 Ca82a	h,57 i \$,g,i,58
^{117}Sn	3pG	4.625(5)	5.058(7)	2.625(3)	0.295(12)	0.64 - 2.37	Fi72	h,57
$^{118}\text{Sn}^*$	2pF 2pF 3pG	4.679(16) 4.676(17) 4.634(5)	5.412(18) 5.442(21) 5.072(7)	0.560(5) 0.543(7) 2.623(3)	0.304(12)	0.49 - 1.40 0.84 - 1.75 0.64 - 2.37	Kh70b Li72b Fi72	c c h,57
^{119}Sn	3pG	4.639(5)	5.100(7)	2.618(7)	0.290(12)	0.64 - 2.37	Fi72	h,57
$^{120}\text{Sn}^*$	2pF 3pG	4.640 4.646(5)	5.315(25) 5.110(7)	0.576(11) 2.619(3)	0.292(12)	0.46 - 1.08 0.64 - 2.37	Ba67b Fi72	h,57
^{122}Sn	3pG	4.658(5)	5.088(7)	2.611(3)	0.378(12)	0.64 - 2.37	Fi72	h,57
$^{124}\text{Sn}^*$	2pF 3pG SOG	4.695(17) 4.670(5) 4.677	5.490(21) 5.150(7)	0.534(7) 2.615(3)	0.311(12)	0.84 - 1.75 0.64 - 2.65 0.36 - 3.60	Li72b Fi72 Ca82a	c h,57 \$,g,i,58
Sb*	2pF	4.63(9)	5.32(11)	0.57(6)		0.56 - 1.31	Ha56	f
^{138}Ba	3pG	4.836	5.3376	2.6776	0.3749	0.56 - 2.84	He70b	h,59
^{139}La	2pF	4.85	5.71(6)	0.535(27)		0.74 - 1.87	Sh67b	c,f
$^{142}\text{Nd}^*$	3pF 2pF 2pF	4.920 4.863(34) 4.993(35)	5.6135 5.774(26) 5.839(33)	0.5868(24) 0.513(16) 0.569(18)	0.096(14)	0.55 - 2.97 0.23 - 0.59 0.22 - 0.73	He71a Ca73 Ma74	h,60 i,61 i
^{144}Nd	2pF	4.926	5.6256	0.6178(30)		0.55 - 2.97	He71a	h,62
$^{146}\text{Nd}^*$	2pF 2pF	4.970 4.993(37)	5.6541 5.867(32)	0.6321(30) 0.556(20)		0.55 - 2.97 0.22 - 0.73	He71a Ma74	h,62 i
^{148}Nd	2pF	5.002	5.6703	0.644(5)		0.55 - 2.97	He71a	h,62
$^{150}\text{Nd}^*$	2pF 2pF 2pF	5.048 5.015(37) 4.948	5.7185 5.865(35) 5.895	0.651(7) 0.571(18) 0.513		0.55 - 2.97 0.22 - 0.73 0.37 - 2.29	He71a Ma74 Hi77	g,h,62 i i,63
^{144}Sm	FB	4.947(9)				0.60 - 2.50	Mo81	†,g,i
^{148}Sm	2pF FB FB	4.989(37) 4.976(8) 5.002(6)	5.771(31)	0.596(15)		0.25 - 0.59 0.33 - 2.18 0.60 - 2.50	Ca73 Ho80 Mo81	i †,i †,g,i

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Nucleus	model	$\langle r^2 \rangle^{1/2}$ [fm]	c or a [fm]	z or α [fm]	w	q-range [fm $^{-1}$]	ref.	remarks
^{150}Sm	FB	5.045(6)				0.60 - 2.50	Mo81	†,g,i
$^{152}\text{Sm}^*$	2pF	5.0922	5.8044	0.581(15)		0.37 - 1.02	Co76	i,64
	FB	5.099(8)				0.33 - 2.18	Ho80	†,i
	FB	5.093(6)				0.60 - 2.50	Mo81	†,g,i
^{154}Sm	2pF	5.1257	5.9387	0.522(15)		0.37 - 0.97	Co76	i,64
	FB	5.126(8)				0.33 - 2.18	Ho80	†,i
^{154}Gd	FB	5.124				0.58 - 2.17	He82	†
^{156}Gd	FB	5.068	5.930	0.576		0.37 - 2.07	Hi77	i,63
^{158}Gd	2pF	5.172(6)				0.36 - 2.17	Mu84	†,i
^{165}Ho	2pF	5.23	6.18	0.57		0.54 - 1.43	Sa67	h,65
	2pF	5.19	6.12	0.57		0.50 - 1.45	Uh71	h,65
^{166}Er	2pF	5.2380	6.1610	0.4872		0.29 - 1.07	Co76	i,66
	2pF	5.303	6.186	0.269		0.37 - 2.28	Cr77	i,67
	3pF	5.237(16)	5.98(6)	0.446(31)	0.19(7)	0.29 - 2.28	Ca78	†,68
^{174}Yb	FB	5.41(3)				0.32 - 2.33	Sa79	†,i
^{176}Yb	2pF	5.3150	6.3306	0.4868		0.29 - 1.07	Co76	i,66
	2pF	5.443	6.127	0.363		0.37 - 2.28	Cr77	i,68
^{175}Lu	FB	5.37(3)				0.32 - 2.33	Sa79	†,i
$^{181}\text{Ta}^*$	2pF	5.48	6.38	0.64		0.56 - 1.42	Do57	69
^{184}W	2pF	5.42(7)	6.51(7)	0.535(36)		0.25 - 0.60	Ka73	70
^{186}W	2pF	5.40(4)	6.58(3)	0.480(23)		0.25 - 0.60	Ka73	70
^{192}Os	FB	5.413(4)				0.60 - 2.90	Re84a	†,i
^{196}Pt	FB	5.38(2)				0.34 - 2.28	Bo83	i
^{197}Au	2pF	5.33(5)	6.38(6)	0.535(27)		0.56 - 1.42	Ha56	
	2pF	5.27(9)				0.08 - 0.27	Be60	71
^{203}Tl	FB	5.463(5)				0.51 - 2.24	Eu78	†,g,i
$^{205}\text{Tl}^*$	FB	5.470(5)				0.51 - 2.24	Eu78	†,g,i
	SOG	5.479				0.51 - 2.99	Fr83	\$,g,72
^{204}Pb	FB	5.479(2)				0.51 - 2.24	Eu78	†,g,i
$^{206}\text{Pb}^*$	2pF	5.509(29)	6.61(5)	0.545(8)		0.22 - 0.88	Ja73	i
	FB	5.490(2)				0.51 - 2.24	Eu78	†,g,i
	SOG	5.490				0.51 - 2.99	Fr83	\$,g,72

TABLE I. Charge-Density-Distribution Parameters
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Nucleus	model	$\langle r^2 \rangle^{1/2}$ [fm]	c or a [fm]	z or α [fm]	w	q-range [fm $^{-1}$]	ref.	remarks
$^{207}\text{Pb}^*$	2pF	5.513(32)	6.62(6)	0.546(10)		0.22 - 0.88	Ja73	i
	FB	5.497(2)				0.51 - 2.24		†,g,i
$^{208}\text{Pb}^*$	SOG	5.503				1.7 - 3.7	Fr77a	\$,g,i,73
	FB	5.4989(7)				0.44 - 3.7		†,g,74
	FB	5.503(2)				0.51 - 2.24		†,g,i,75
$^{209}\text{Bi}^*$	2pF	5.51(5)	6.75(7)	0.468(39)	0.39(6)	0.07 - 0.53	Ni69	i,76
	3pG	5.521(2)				0.7 - 2.8		Si73b
	FB	5.519(4)				0.51 - 2.24		†,g,i,77
$^{232}\text{Th}^*$	2pF	5.7723	6.7915	0.571(15)		0.33 - 0.97	Co76	i,64
	2pF	5.645				0.46 - 2.08		i,63
$^{238}\text{U}^*$	2pF	5.84	6.8054	0.605(16)		0.37 - 0.97	Co76	i,64
	2pF	5.854				0.46 - 2.08		i,67

TABLE I. Charge-Density-Distribution Parameters
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*) Additional information can be found in the following references:	
n	: Vr64, Du65, Du66, Hu66a, Hu66b, Ch66a, Gr66a, Gr66b, St66a, Al68, Bu68, Bu69, El69, Ga71, Ba72, Be73a, Ha73a
¹ H	: Bu61, Li61, Ol61, Dr62, Le62, Be63, Ch63, Du63, Du64, Du65, Al66, Ba66, Ch66a, Ch66b, Fr66, Gr66b, Ja66, Al67, Ba67a, Be67a, Go67, Be68, Ba70, Go70, Li70, Be71a, Bi71, Ha71, Ja71, Pr71, Ak72, Ga72, Ba73, Be73a, Bo73, Ki73, Bo74, Mu74, Th74, Bo75a, Bo75b
² H	: MI57a, MI58, Be64a, Gr66a, El69, Sk73, Bu70, Ar75
³ H	: Co65, Be82
³ He	: Co65, Be72, Sz77, Ar78, Gu82, Du83
⁴ He	: B156, Ho56, MA56, Bu60, Re65, Fr67, Er68, St69, Si76, MC77, Ar78
⁶ Li	: Bu58, Be65
⁷ Li	: Be65
⁹ Be	: Ho57, Me59, Ng64, Be67b, Be67c, Be69, Be73b
¹⁰ B	: Me59
¹¹ B	: Me59
¹² C	: Fr56, Eh59, Cr66, Af67, En67, Si70b, Be71b, Ja72, Fe73a, Fe73b, Ki73, Si74a, Me76, Do79
¹⁴ N	: Me59, Bi64, Be65, Be71b, Vo78
¹⁵ N	: Pa68
¹⁶ O	: Eh59, Me59, Cr66, Be71b, Si70a, No82
¹⁹ F	: Wi78
²⁰ Ne	: Fe73a
²⁴ Mg	: He56, Sa69b, Ju70, Na72
²⁷ Al	: Be70a, Li74, Do83
²⁸ Si	: He56, Sa69c, Be70a, Mu70, Na72, Fe73a, Av74, Wh79
³¹ P	: Ko65
³² S	: He56, Lo64, Hu70, Si72, Av74, Si74a, Ry83b
³⁴ S	: Ry83b
³⁶ S	: Ry83b
³⁹ K	: Si74b
⁴⁰ Ar	: He56, Gr71, Sc71, Fe73a
⁴⁰ Ca	: Ha56, Cr61, Cr65, Be67d, Fr68, El69, He71b, Si74b, Ca80b, Em83a
⁴² Ca	: He71b
⁴⁴ Ca	: He71b
⁴⁸ Ca	: Si74b, Em83a
Ti	: En66, Fe73a
⁴⁶ Ti	: He71b
⁴⁸ Ti	: He71b
⁵⁰ Ti	: He71b
	⁵¹ V : Ha56, Cr61, Na74
	⁵⁰ Cr : Sh78c
	⁵² Cr : Be64b, Li76
	Fe : Fe73a, Th70
	⁵⁴ Fe : Be62, Li71b
	⁵⁶ Fe : Be62, He71b, Li71b
	⁵⁹ Co : Ha56, Cr61, Go63, Br66
	Ni : Fe73a, Th70
	⁵⁸ Ni : Af70, Kh70b, Li71b, Si75, Ke77, Ca80b, Wo76
	⁶⁰ Ni : Af70, Kh70b, Li72a, Ke77, Sh78a, Wo80
	⁶² Ni : Li71b, Go74
	⁶⁴ Ni : Af70
	Cu : Fe73a
	⁶³ Cu : Go74
	⁶⁵ Cu : Go74
	Zn : Th70, Fe73a
	⁶⁴ Zn : Af71, Li72a
	⁶⁶ Zn : Af71, Li72a, Ne72
	⁶⁸ Zn : Ne72
	⁸⁸ Sr : He56, Sc74
	⁸⁹ Y : Pe68
	⁹⁰ Zr : Be70b, Ph72, Dr74
	Mo : Ja76
	Cd : Ja76
	In : Ha56, Cr61, Ke63
	¹¹⁶ Sn : Ba67b, Cu69, Li72b, Ca80b
	¹¹⁸ Sn : Cu69
	¹²⁰ Sn : Cu69
	¹²⁴ Sn : Ba67a, Cu69, Ca80b
	Sb : Cr61
	¹⁴² Nd : He70b
	¹⁴⁶ Nd : Ma71
	¹⁵⁰ Nd : Ma71
	Hf : Ha56, Do57
	¹⁵² Sm : Ca78
	¹⁸¹ Ta : Ke63, Ra78
	W : Pi55, Ha56, Do57
	²⁰⁵ Tl : Ca82b
	Pb : Mi57b, Fi64, Ni69, Fe73a
	²⁰⁶ Pb : Pa79, Ca82b
	²⁰⁷ Pb : Pe65, Pa79
	²⁰⁸ Pb : Pe65, Be67e, He69, Fr72b, Ja73, Eu76a, Na71, Si73b, Be77, Eu77b, Be80, Ca80b
	²⁰⁹ Bi : Ha56, Be60, Cr61, Go63, He63, Br66
	²³² Th : Ha56, Do57
	²³⁸ U : Ha56, Do57

TABLE I. Charge-Density-Distribution Parameters
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General remarks

- †) Additional information can be found in Table IV (Fourier-Bessel coefficients).
- \$) Additional information can be found in Table V (Sum-of-Gaussians).

- a) Analysis performed in the Plane Wave Born Approximation.
- b) Analysis performed in the Modified Born Approximation, using an effective q-value.
- c) Analysis performed in the High Energy Approximation (Pe66).
- d) Only statistical errors are quoted, corresponding to one standard deviation.
- e) The value of z was fixed in the analysis.
- f) A target of natural isotopic composition has been used.
- g) Muonic X-ray data have been included in the analysis.
- h) Measurement relative to ^1H .
- i) Measurement relative to ^{12}C .

Specific remarks

- 1) The tabulated value for the rms radius is formally obtained from the slope of the neutron charge form factor at $q^2 = 0$ as determined from the scattering of slow neutrons by atomic electrons. Further information on the neutron form factor, obtained from electron scattering of ^2H and ^3H , is given in the list of additional references. In ref. Ho76 a combined analysis is presented of data from scattering of slow neutrons by atomic electrons (refs. Kr73 and Ko76), from elastic deuteron scattering (refs. Dr62, Be64a, Bu70, Ga71 and Be73a) and from quasi-elastic e-D scattering (refs. Ak64, St66a, Al68, Bu68, Ba73 and Ha73a).
- 2) Result of an analysis of all available data below $q^2 = 2 \text{ fm}^{-2}$ (Bu61, Dr62, Le62, Du65, Fr66, Ja71).
- 3) Result of the analysis of the data from refs. Dr62, Du65, Fr66, Ja66, Be67a, Go70, Li70, Be71a, Pr71, Ga72, Ba73, Ki73, Bo74, Mu74, At75, Bo75a, Bo75b, St75.
- 4) Model-independent fourth order polynomial fit (ref. Bo75a) with free normalization. The analysis includes the data from refs. Du65, Be71a, Ba73, Ki73, Bo74 and Mu74.
- 5) In the analysis values of 0.336 and 0.805 fm have been used for the rms radius of the neutron and of the proton, respectively. The rms charge radius is related to the rms structure radius through $r_c^2 = r_d^2 + r_p^2 - r_n^2$.
- 6) Although the complete data set covers a momentum transfer range up to 1.99 fm^{-1} , only the low q-data were used for the determination of the rms radius via a third-order polynomial fit. The analysis included data from refs. Bu70 and Be73a. In contrast to the procedure followed in refs. Bu70 and Be73a, no additional assumption has been made about the coefficient of q^4 .
- 7) In this experiment the form factor has been determined for momentum transfer values between 0.51 and 1.72 fm^{-1} . The data points from refs. Co65 and Be82 have been included in the data analysis.
- 8) The data from refs. Co65 and Be84 have been included in the analysis.
- 9) In this experiment the recoil ^3He has been detected. A momentum transfer range of 0.94 to 1.79 fm^{-1} was covered. The data from refs. Co65, MC77, Sz77, Ar78 and Du83 were included in the analysis. The charge density was defined to be positive.
- 10) A fifth-order polynomial was fitted to the form factor.
- 11) The analysis included the data presented in ref. Fr67, Er68, MC77 and Ar78.
- 12) A fourth-order polynomial was fitted to the form factor.
- 13) The data could be described excellently by a charge distribution which is the Fourier transform of $F(q^2) = \exp(-a^2 q^2) - c^2 q^2 \exp(-b^2 q^2)$, with parameter values $a = 0.933(3) \text{ fm}$, $b = 1.30(6) \text{ fm}$ and $c = 0.45(3)$.
- 14) The data could be described out to $q^2 = 6 \text{ fm}^{-2}$ by a charge distribution as given in remark 13. A fit to the complete data set was obtained after adding an oscillatory modification: the Fourier transform of a form factor modification $\Delta F = d \exp(-(q-q_0)^2/p^2)$. The best fit results are: $a = 0.928(3) \text{ fm}$, $b = 1.26(9) \text{ fm}$, $c = 0.48(4)$, and $d = -0.00124(28)$, $p = 0.70(29) \text{ fm}^{-1}$ and $q_0 = 3.11(20) \text{ fm}^{-1}$.
- 15) In the analysis the value for the rms radius of ^{12}C of ref. Si70b was used.
- 16) The form factor has been interpreted in terms of a harmonic-oscillator shell model with a quadrupole contribution based on an undeformed p-shell model. The absolute value obtained for the quadrupole moment ($4.20 \pm 0.25 \text{ e fm}^2$) is in excellent agreement with spectroscopic measurements.
- 17) The normalization of the data of ref. Be67c has been adjusted with the value of the ^{12}C radius of ref. Si70b.
- 18) The data were analyzed with nuclear wave functions obtained by extending the Nilsson model to include single-particle orbital admixtures from higher major shells.
- 19) Combined analysis with the data from refs. Si70b and Ja72 with a free normalization for each data set. The data of this experiment cover a momentum transfer range from 0.1 to 1.0 fm^{-1} . In ref. Ca80 the 10th Fourier-Bessel coefficient is a factor 10 too high due to a typing error.

TABLE I. Charge-Density-Distribution Parameters
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- 20) Combined analysis with the data from refs. Si70b, Ja72, Fe73b and Ca80.
- 21) An exponential tail modification of a Gaussian form was added to the MHO charge distribution in order to approximate the 3pF density in the tail region. Only the amplitude of the tail modification was fitted as a free parameter, the other two parameters were taken from the ^{12}C results of ref. Si70b.
- 22) Results of a fit to the data up till the first diffraction minimum without applying corrections for elastic scattering from the M1 and C2 moments.
- 23) A Gaussian shape has been assumed for the static quadrupole moment distribution. The analysis yielded a value of $Q_2 = 3.22 \text{ e fm}^2$.
- 24) Reanalysis of the data presented in ref. Da70.
- 25) Analysis with the data presented in ref. Sc75. The present data covered a momentum transfer range from 0.35 to 3.17 fm^{-1} .
- 26) In the analysis an oscillatory modification corresponding to the Fourier transform of a damped sine wave, was added to the MHO charge distribution.
- 27) Measurement relative to the ^{16}O parameters from ref. Si70a.
- 28) Measurement relative to the ^{40}Ca parameters from refs. Fr68, Ei69 and He71b.
- 29) The data were analyzed with the uniform Gaussian model, which yielded parameter values $r = 3.13(6) \text{ fm}$ and $g = 0.96(5) \text{ fm}$.
- 30) The oscillatory modification defined in remark 14, was included in the analysis of the ^{24}Mg and the ^{32}S data. The best fit results were: $d = -0.076$, $p = 0.51 \text{ fm}^{-1}$ and $q_0 = 2.49 \text{ fm}^{-1}$ for ^{24}Mg and $d = 0.021$, $p = 0.50 \text{ fm}^{-1}$ and $q_0 = 2.83 \text{ fm}^{-1}$ for ^{32}S .
- 31) The analysis yielded the following values for the C2 and C4-moments: $24.4(+0.8,-4.0) \text{ e fm}^2$ and $15.3(+2.3, -10.0) \text{ e fm}^4$, respectively.
- 32) Reanalysis of the data from ref. Be70a.
- 33) In the 3pF analysis only the data up till 1.49 fm^{-1} were used. Data were taken up to 2.64 fm^{-1} .
- 34) In this analysis an oscillatory modification as defined in remark 14 was used, which yielded parameter values $d = -0.034(8)$, $p = 0.51(11) \text{ fm}^{-1}$ and $q_0 = 2.48(7) \text{ fm}^{-1}$.
- 35) Combined analysis with the data from ref. Si72. The present data cover a momentum transfer range from 0.3 to 2.3 fm^{-1} .
- 36) The data from ref. Me76 were included in the analysis.
- 37) Combined analysis with the data from ref. Li74b. The present data cover a momentum transfer range from 0.3 to 2.3 fm^{-1} .
- 38) For the subtraction of the C2-contribution a value for the quadrupole moment of -8.25 e fm^2 was used.
- 39) For the subtraction of the C2-contribution a value for the quadrupole moment of -6.2 e fm^2 was used.
- 40) The present experiment covered a momentum range from 0.54 to 1.26 fm^{-1} . In the analysis the data from refs. Gr71, Sc71, We74 and Fi76 were also included.
- 41) The data were analyzed with the uniform Gaussian model, which yielded parameter values $r = 3.83(8) \text{ fm}$ and $g = 0.96(5) \text{ fm}$.
- 42) The analysis included an oscillatory modification as defined in remark 14, which yielded $d = 0.086(7)$, $p = 0.43(4) \text{ fm}^{-1}$ and $q_0 = 3.14(6) \text{ fm}^{-1}$.
- 43) The analysis included an oscillatory modification as defined in remark 14, which yielded $d = 0.0814(8)$, $p = 0.43(4) \text{ fm}^{-1}$ and $q_0 = 3.14(5) \text{ fm}^{-1}$. The data of ref. Be67d and unpublished data taken by J. Heisenberg, J. McCarthy and I. Sick, were included in the analysis.
- 44) The present experiment covered a momentum range from 2.14 to 3.56 fm^{-1} . In the analysis the data from refs. Be67d, Fr68 and Si73a were also included.
- 45) The present experiment covered a momentum range from 0.35 to 2.38 fm^{-1} . In the analysis the high-q data from ref. Si79 were also included.
- 46) The analysis included an oscillatory modification as defined in remark 14, which yielded $d = 0.08$, $p = 0.5 \text{ fm}^{-1}$ and $q_0 = 3 \text{ fm}^{-1}$.
- 47) The rms radius for ^{12}C was taken from ref. En67 to be $2.42(4) \text{ fm}$.
- 48) The normalization of the data presented in ref. Th70 has been adjusted with the value of the ^{12}C radius of ref. Si70b.
- 49) The data were analyzed simultaneously with the results of optical isotope shift experiments. An oscillatory modification corresponding to the Fourier transform of a damped sine wave, was also included in the analysis. The entries presented here are a reanalysis of the original data (I. Sick, private communication, 1973).
- 50) Combined analysis with the data from refs. Cu69 and Fi72.
- 51) The normalization of the data has been adjusted with the value of the ^{12}C radius of ref. Si70b by the authors of the compilation.
- 52) Combined analysis with the data from ref. Sc74.

TABLE I. Charge-Density-Distribution Parameters

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- 53) In a private communication the authors gave preference to the results obtained with the 3pG model for ease of comparison with the other tabulated results. The value of w was fixed in the analysis. The errors were obtained by assuming the same percentage errors as yielded by the analysis with the 2pF model. The entries presented in ref. Si73c have been calculated with a faulty version of the phase shift code. The entries listed in this table have been recalculated by the original authors.
- 54) Measurements relative to the Zr parameters from ref. Fa71.
- 55) The entries presented in ref. Gi75 have been calculated with a faulty version of the phase shift code. The entries listed in this table have been recalculated by the original authors.
- 56) Measurements relative to the Sn parameters from ref. Fi72.
- 57) The data were analyzed simultaneously with the results of optical isotope shift experiments. An oscillatory modification corresponding to the Fourier transform of a damped sine wave, was included in the analysis.
- 58) The present data cover a momentum transfer range between 1.4 and 3.6 fm⁻¹. The data were analyzed simultaneously with the data from refs. Cu69 and Fi72.
- 59) In the analysis an oscillatory modification as defined in remark 14, was included, which yielded $d = 0.13$, $p = 0.25$ fm⁻¹ and $q_0 = 2.55$ fm⁻¹.
- 60) The analysis included an oscillatory modification.
- 61) Reanalysis of the data presented in ref. Ma71, omitting the 15 MeV points. The normalization of the data has been adjusted with the value of the ¹²C radius of ref. Si70b.
- 62) In the analysis for ^{144,146,148}Nd K _{α} X-ray data relative to ¹⁵⁰Nd were used as constraints.
- 63) The elastic electron scattering data were analyzed with a deformed Fermi distribution for the ground state. The values found for the deformation parameters β_2 , β_4 and β_6 are 0.2693, 0.0795 and 0.0161(¹⁵⁰Nd); 0.3147, 0.0648 and 0.0020(¹⁵⁶Gd) and 0.2122, 0.0607 and -0.2490(²³²Th), respectively.
- 64) The elastic electron scattering data were analyzed simultaneously with data for electro-excitation of the ground-state rotational band with a deformed Fermi distribution. The values for the rms radius and for the deformation parameters β_2 and β_6 were taken from other types of experiments, which left β_4 and z as free parameters. The values for β_2 , β_4 and β_6 either used in or yielded by the analysis, are 0.287(3), 0.070(3) and -0.012(¹⁵²Sm); 0.311(3), 0.087(2) and -0.018(¹⁵⁴Sm); 0.238(2), 0.101(3) and 0.0(²³²Th) and 0.261(2), 0.087(3) and 0.0(²³⁸U), respectively.
- 65) Cross sections were measured for electron scattering from randomly oriented and from aligned ¹⁶⁵Ho nuclei. The tabulated values present the results of an analysis of the data for randomly oriented nuclei with the 2pF distribution, after subtraction of the scattering from the quadrupole moment.
- 66) The elastic electron scattering data were analyzed with a deformed Fermi distribution for the ground state. The values for the deformation parameters β_2 , β_4 and β_6 were kept fixed at 0.3266, 0.0 and -0.018(¹⁶⁶Er) and 0.3100, -0.054 and -0.006(¹⁷⁶Yb), respectively.
- 67) The elastic electron scattering data were analyzed with a deformed Fermi distribution for the ground state. The values for the deformation parameters β_2 , β_4 and β_6 found are 0.4874, -0.0259 and 0.0423(¹⁶⁶Er); 0.4987, -0.0525 and 0.1451(¹⁷⁶Yb) and 0.2802, -0.0035 and -0.1107(²³⁸U), respectively.
- 68) Combined analysis with the data from refs. Co76 and Cr77. No good fit could be obtained with either a deformed Fermi distribution or a modified Gaussian distribution.
- 69) Analysis of the data presented in ref. Ha56, with a deformed 2pF distribution. The tabulated values are for a 2pF distribution which gives a good approximation to the monopole term of the deformed distribution.
- 70) Measurements relative to ⁶Li, for which the parameters of ref. Li71a have been used.
- 71) Analysis with a uniform charge distribution.
- 72) The low-q data from ref. Eu78 were included in the analysis. The data from ref. Eu78 taken at 289 MeV were excluded due to inconsistencies with the present data set.
- 73) The data covered a momentum range from 1.7 to 3.7 fm⁻¹. The data were analyzed simultaneously with those from refs. He69, Eu76a, Ni69 and muonic X-ray experiments. The data from ref. Eu78 in the momentum range from 1.8 to 2.3 fm⁻¹ were excluded because they were incompatible with the other data.
- 74) Reanalysis of the data from refs. He69, Ni69, Na71, Fr72b, Eu76 and Fr77a.
- 75) The high-q (2.35 to 2.73 fm⁻¹) data from ref. He69 were also included in the analysis.
- 76) Also ⁹Be has been used as a comparison nucleus. The normalization of the data has been adjusted with the values of the ⁹Be radius of ref. Fe73a and the ¹²C radius of ref. Si70b by the authors of the compilation.
- 77) The high-q (2.11 to 2.62 fm⁻¹) data from ref. Si73b were also included in the analysis.

TABLE II. Differences in Charge-Density-Distribution Parameters between Isotopes
See page 500 for Explanation of Tables

Isotope pair	model	$\Delta \langle r^2 \rangle^{1/2}$ [fm]	Δc [fm]	Δz [fm]	Δw	q-range [fm $^{-1}$]	ref.	remarks
4- ^3He	MI	-0.271(15)				0.12 - 0.53	Gu82	
7- ^6Li	MI	-0.08(2)				0.51 - 1.27	Be65	
	MI	-0.13(2)				0.35 - 0.71	Su67	1.2
	MI	-0.003(20)				0.36 - 0.78	Ni71	1
13- $^{12}\text{C}^*$	MHO	-0.023(10)				0.30 - 1.50	He70a	
	HO	-0.06(5)				0.26 - 1.23	Ya71	1,3
	HO	-0.012(19)				0.26 - 0.55	Be71b	
14- ^{12}C	MHO	0.12(5)				1.04 - 2.16	KI73	
15- ^{14}N	HO	0.040(12)				0.22 - 0.48	Sc75	
17- $^{16}\text{O}^*$	HO	-0.013(15)				0.39 - 0.99	Si70a	
	HO	0.004(15)				0.46 - 1.21	Ki78	
	FB	-0.008(7)				0.50 - 2.60	Mi79	
18- $^{16}\text{O}^*$	HO	0.053(12)				0.47 - 0.99	Si70a	
	HO	0.070(11)				0.22 - 0.48	Sc75	
	FB	0.074(5)				0.50 - 2.60	Mi79	
22- ^{20}Ne	2pF	-0.073(27)				0.22 - 1.04	Mo71	
	2pF	-0.053(9)	0.098(10)	-0.042(6)		0.21 - 1.30	Kn81	
25- ^{24}Mg	2pF	0.016(32)	-0.063(36)	0.021(23)		0.20 - 1.15	Le76	
	UG	-0.05(4)				0.69 - 1.40	Kh70a	5
	2pF	-0.019(32)	0.055(36)	-0.021(23)		0.20 - 1.15	Le76	
29- ^{28}Si	2pF	-0.052(25)	0.03(3)	-0.030(20)		0.16 - 1.1	Br77	
30- ^{28}Si	2pF	0.03(+15,-7)	0.04(37)	0.(-18,-08)		0.16 - 1.1	Br77	
34- ^{32}S	FB	0.032(10)				0.50 - 2.60	Ry83a	
	FB	-0.007(12)				0.50 - 2.60	Ry83a	
	FB	0.034(10)				0.50 - 2.60	Ry83a	
42- ^{40}Ca	3pF	0.030	0.052	0.006	-0.014	0.55 - 1.70	Fr68	
44- ^{40}Ca	3pF	0.028	0.072	-0.014	0.007	0.70 - 1.79	Fr68	
48- $^{40}\text{Ca}^*$	3pF	-0.0107	0.069(17)	-0.060(18)	0.072(13)	0.49 - 2.53	Fr68	
	2pF	-0.04(7)	0.16(11)	-0.065		0.13 - 0.59	Ei69	6
	FB	-0.0014(26)				0.35 - 3.35	Em83b	
48- ^{46}Ti	2pF	-0.005(27)	-0.01(4)	0.0		0.22 - 0.57	Th67	
	2pF	-0.015(10)	0.055(20)	-0.025(10)		0.55 - 1.11	Ro71	
	3pF	0.003(15)	-0.0076	-0.0293	0.0624	0.55 - 2.42	He72	
50- ^{46}Ti	2pF	0.003(21)	0.005(33)	0.0		0.22 - 0.57	Th67	
	2pF	-0.030(15)	0.090(25)	-0.045(15)		0.55 - 1.11	Ro71	
	3pF	-0.013(15)	-0.0137	-0.0509	0.1004		He72	
50- ^{48}Ti	2pF	-0.020(10)	0.045(15)	-0.025(10)		0.55 - 1.11	Ro71	
	3pF	-0.016(15)	-0.0061	-0.0216	0.0380	0.55 - 2.49	He72	
52- ^{50}Cr	2pF	-0.037(11)	0.064(28)	-0.042(15)		0.15 - 0.74	La76	
	2pF	-0.028(12)	0.047(8)	-0.0289(27)		0.97 - 1.62	Sh78b	7
53- ^{50}Cr	2pF	0.027(12)	0.051(8)	-0.0034(27)		0.97 - 1.62	Sh78b	7

TABLE II. Differences in Charge-Density-Distribution Parameters between Isotopes
See page 500 for Explanation of Tables

Isotope pair	model	$\Delta \langle r^2 \rangle^{1/2}$ [fm]	Δc [fm]	Δz [fm]	Δw	q-range [fm $^{-1}$]	ref.	remarks
54- 50Cr	2pF	0.033(6)	0.044(15)	0.002(8)		0.15 - 0.74	La76	
	2pF	0.068(12)	0.078(9)	0.0082(27)		0.97 - 1.62	Sh78b	7
54- 52Cr	2pF	0.048(9)	0.015(25)	0.035(13)		0.15 - 0.74	La76	
56- 54Fe*	2pF	0.048(9)	0.015(25)	0.020(13)		0.15 - 0.74	La76	
	FB	0.0445(8)				0.51 - 2.22	Wo80	8
58- 56Fe	2pF	0.070(9)	0.031(6)	0.0241(20)		0.97 - 1.62	Sh78b	7
	2pF	0.004(18)	0.092(20)	-0.026(6)		1.02 - 1.77	Li71b	7
	FB	0.0382(10)				0.51 - 2.22	Wo80	8
60- 58Ni*	2pF	0.040	0.062(18)	0.000(5)		1.05 - 1.61	Kh70b	7
	FB	0.0373(11)				0.51 - 2.22	Wo80	8
	FB	0.027(7)				0.25 - 2.65	Be83	
62- 58Ni	2pF	0.061(8)	0.081(9)	0.006(3)		1.02 - 1.93	Li71b	7
	2pF	0.079(6)	0.065(13)	0.018(8)		0.15 - 0.79	Ke77	
62- 60Ni	2pF	0.034(6)	0.033(13)	0.006(8)		0.15 - 0.79	Ke77	
	FB	0.0286(8)				0.51 - 2.22	Wo80	8
64- 58Ni	2pF	0.082	0.095(18)	0.010(5)		1.05 - 1.61	Kh70b	7
64- 60Ni	2pF	0.043	0.044(16)	0.006(5)		1.05 - 1.61	Kh70b	7
64- 62Ni	FB	0.0181(6)				0.51 - 2.22	Wo80	8
65- 63Cu*	2pF	0.024(9)	0.055(14)	-0.006(8)		0.29 - 0.88	Go74	
	FB	0.0202(6)				0.51 - 2.22	Sc77	8
	2pF	0.022(7)	0.047(8)	-0.0041(23)		0.96 - 1.61	Sh78a	7
66- 64Zn*	2pF	0.025(13)	0.048(13)	-0.003(3)		0.98 - 1.61	Li72a	7
	2pF	0.018(11)	0.030(21)	-0.001(12)		0.15 - 0.79	Ke77	
	FB	0.0225(30)				0.51 - 2.22	Wo80	8
68- 66Zn*	2pF	-0.016(9)	0.059(5)	-0.029(5)		0.30 - 1.09	Ne72	
	2pF	0.007(13)	0.054(24)	0.015(15)		0.15 - 0.79	Ke77	
	FB	0.0144(30)				0.51 - 2.22	Wo80	8
70- 68Zn	2pF	0.047(15)	0.032(10)	0.014(8)		0.30 - 1.09	Ne72	
	2pF	0.026(16)	0.049(31)	0.006(4)		0.15 - 0.79	Ke77	
	FB	0.0173(25)				0.51 - 2.22	Wo80	8
72- 70Ge	FB	0.0168(24)				0.35 - 2.90	Ma84	
74- 72Ge	FB	0.0151(24)				0.35 - 2.90	Ma84	
76- 74Ge	FB	0.0064(24)				0.35 - 2.90	Ma84	
76- 70Ge	FB	0.0383(24)				0.35 - 2.90	Ma84	
94- 92Mo	FB	0.040(5)				0.56 - 1.96	Dr75	
96- 94Mo	FB	0.028(5)				0.56 - 1.96	Dr75	
98- 96Mo	FB	0.028(5)				0.56 - 1.96	Dr75	
100- 98Mo	FB	0.039(5)				0.56 - 1.96	Dr75	
112-110Cd	2pF	0.030(10)				0.25 - 1.07	Gi75	
114-112Cd	2pF	0.021(11)				0.25 - 1.07	Gi75	
116-114Cd	2pF	0.010(11)				0.25 - 1.07	Gi75	
116-110Cd	2pF	0.061(12)				0.25 - 1.07	Gi75	
118-112Sn	2pF	0.019	0.021(20)	0.000(7)		0.82 - 1.40	Kh70b	7
118-116Sn	2pF	-0.004(6)	0.025(6)	-0.013(2)		0.84 - 1.75	Li72b	7

TABLE II. Differences in Charge-Density-Distribution Parameters between Isotopes
See page 500 for Explanation of Tables

Isotope pair	model	$\Delta \langle r^2 \rangle^{1/2}$ [fm]	Δc [fm]	Δz [fm]	Δw	q-range [fm $^{-1}$]	ref.	remarks
124-116Sn	2pF	0.015(7)	0.069(7)	-0.020(7)		0.84 - 1.75	Li72b	7
124-118Sn	2pF	0.019(8)	0.045(9)	-0.007(3)		0.84 - 1.75	Li72b	7
146-142Nd	2pF	-0.10				0.24 - 0.59	Ca73	9
	2pF	0.0527(36)	0.018(26)	0.025(11)		0.22 - 0.73	Ma74	10
150-142Nd	2pF	0.08				0.24 - 0.59	Ca73	9
	2pF	0.115(5)	0.078(27)	0.041(14)		0.22 - 0.73	Ma74	10
150-146Nd	2pF	0.067(4)	0.046(27)	0.023(14)		0.22 - 0.73	Ma74	10
150-148Sm	2pF	0.03(7)				0.25 - 0.59	Ca73	
152-148Sm	2pF	0.17(7)				0.25 - 0.59	Ca73	
	FB	0.12(2)				0.33 - 2.18	Ho80	
154-152Sm	FB	0.03(2)				0.33 - 2.18	Ho80	
207-206Pb*	2pF	0.005(7)	0.008(16)	0.001(4)		0.22 - 0.88	Ja73	
208-206Pb*	2pF	0.013(4)	0.010(9)	0.0039(25)		0.22 - 0.88	Ja73	
208-207Pb*	2pF	0.008(7)	0.001(15)	0.005(4)		0.22 - 0.88	Ja73	

*) Additional information can be found in the following references:

13 - ^{12}C : Cr67	66 - ^{64}Zn : Ne72
17 - ^{16}O : No82	68 - ^{66}Zn : Li73
18 - ^{16}O : La61, No82	205 - ^{203}Tl : Eu78
48 - ^{40}Ca : Em83a	208 - ^{206}Pb : Eu78
56 - ^{54}Fe : Li71b	208 - ^{207}Pb : Eu78
60 - ^{58}Ni : Ha57, Ke77	207 - ^{206}Pb : Pe65, Eu78
65 - ^{63}Cu : Ke77	

Remarks

- 1) Analysis performed in Plane Wave Born Approximation.
- 2) No correction applied for scattering from the C2 or higher charge multipole moments.
- 3) No correction applied for the magnetic contribution to the elastic scattering.
- 4) The scattering from the C2 and the C4 distribution has been subtracted.
- 5) Analysis performed in the Modified Born Approximation. The parameters obtained for the Uniform Gaussian model were $\Delta r = 0.03(2)$ fm and $\Delta g = -0.07(4)$ fm.
- 6) The value of Δz was fixed in the analysis.
- 7) Analysis performed in the High Energy Approximation (Pe66).
- 8) Data analyzed simultaneously with muonic X-ray data.
- 9) Reanalysis of the data presented in ref. Ma71.
- 10) Data analyzed simultaneously with K_{α} X-ray data.

TABLE III. Differences in Charge-Density-Distribution Parameters between Neighboring Nuclei (Not Isotopes)
See page 500 for Explanation of Tables

Nucleus pair	model	$\Delta \langle r^2 \rangle^{1/2}$ [fm]	Δc [fm]	Δz [fm]	Δw	q-range [fm $^{-1}$]	ref.	remarks
$^{14}\text{N} - ^{12}\text{C}$	FB	0.098(19)				0.26 - 0.55	Be71b	
$^{16}\text{O} - ^{12}\text{C}$	FB	0.271(22)				0.26 - 0.55	Be71b	
$^{16}\text{O} - ^{14}\text{N}$	HO	0.166(19)				0.29 - 0.48	Sc75	
$^{20}\text{Ne} - ^{16}\text{O}$	2pF	0.279(20)				0.22 - 0.48	Fr72a	
$^{46}\text{Ti} - ^{45}\text{Sc}$	2pF	0.04(5)	0.06(8)	0.0		0.29 - 0.59	Th70	1
$^{48}\text{Ti} - ^{40}\text{Ca}$	3pF	0.097	0.179	-0.022	0.026	0.42 - 1.79	Fr68	
$^{50}\text{Cr} - ^{48}\text{Ti}$	2pF	-0.009(12)	0.104(10)	-0.0357(32)		1.03 - 1.62	Sh78c	
$^{51}\text{V} - ^{50}\text{Ti}$	2pF	0.06(5)	0.10(8)	0.0		0.29 - 0.59	Th70	1
$^{54}\text{Fe} - ^{54}\text{Cr}$	2pF	-0.002(12)	0.069(28)	-0.025(16)		0.15 - 0.74	La76	
	2pF	-0.058(12)	0.051(8)	-0.0448(27)		0.97 - 1.62	Sh78b	
$^{56}\text{Fe} - ^{50}\text{Cr}$	2pF	0.081(12)	0.153(8)	-0.0093(27)		0.97 - 1.62	Sh78b	2
$^{56}\text{Fe} - ^{54}\text{Cr}$	2pF	0.014(12)	0.085(9)	-0.0200(27)		0.97 - 1.62	Sh78b	2
$^{59}\text{Co} - ^{58}\text{Fe}$	FB	0.0120(23)				0.51 - 2.22	Sc77	3
$^{58}\text{Ni} - ^{54}\text{Fe}$	2pF	0.050(8)	0.064(8)	0.002(3)		1.02 - 1.77	Li71b	2
$^{58}\text{Ni} - ^{56}\text{Fe}^*$	2pF	-0.022(13)	0.107(14)	-0.046(4)		1.02 - 1.98	Li71b	2
	FB	0.0386(10)				0.51 - 2.22	Wo80	3
$^{58}\text{Ni} - ^{58}\text{Fe}$	2pF	-0.026(12)	0.005(14)	-0.016(4)		1.02 - 1.77	Li71b	2
$^{60}\text{Ni} - ^{58}\text{Fe}$	FB	0.0382(23)				0.51 - 2.22	Wo80	3
$^{60}\text{Ni} - ^{59}\text{Co}$	2pF	0.019(7)	0.042(6)	0.0009(23)		0.96 - 1.61	Sh78a	
	FB	0.0252(23)				0.51 - 2.22	Sc77	3
$^{63}\text{Cu} - ^{62}\text{Ni}$	FB	0.0419(7)				0.51 - 2.22	Sc77	3
$^{65}\text{Cu} - ^{60}\text{Ni}$	2pF	0.097(15)	0.093(11)	0.016(4)		0.96 - 1.61	Sh78a	2
$^{65}\text{Cu} - ^{64}\text{Ni}$	FB	0.0440(20)				0.51 - 2.22	Sc77	3
$^{64}\text{Zn} - ^{60}\text{Ni}$	2pF	0.132(20)	0.033(20)	0.052(5)		0.98 - 1.61	Li72a	2
	FB	0.0893(12)				0.51 - 2.22	Wo80	3
$^{64}\text{Zn} - ^{62}\text{Ni}$	2pF	0.094(7)	0.024(14)	0.041(8)		0.15 - 0.79	Ke77	
	FB	0.0893(12)				0.51 - 2.22	Wo80	3
$^{64}\text{Zn} - ^{63}\text{Cu}$	2pF	-0.024(11)	0.168(22)	-0.063(11)		0.15 - 0.79	Ke77	
	FB	0.0473(7)				0.51 - 2.22	Sc77	3
$^{66}\text{Zn} - ^{60}\text{Ni}$	2pF	0.157(20)	0.082(19)	0.049(4)		0.98 - 1.61	Li72a	2
$^{66}\text{Zn} - ^{64}\text{Ni}$	FB	0.0938(50)				0.51 - 2.22	Wo80	3
$^{66}\text{Zn} - ^{65}\text{Cu}$	FB	0.0497(20)				0.51 - 2.22	Sc77	3
$^{90}\text{Zr} - ^{88}\text{Sr}$	2pF	0.046(12)	0.030(23)	0.014(7)		0.24 - 1.08	Si73c	4
	3pG	0.050(4)	0.091(18)	0.043(4)	-0.074(27)	0.51 - 2.05	Sc74	
	FB	0.055(7)				0.51 - 2.05	Sc74	

TABLE III. Differences in Charge-Density-Distribution Parameters between Neighboring Nuclei (Not Isotopes)
See page 500 for Explanation of Tables

Nucleus pair	model	$\Delta \langle r^2 \rangle^{1/2}$ [fm]	Δc [fm]	Δz [fm]	Δw	q-range [fm $^{-1}$]	ref.	remarks
^{92}Mo - ^{90}Zr	3pG	0.054(22)	0.04(5)	0.01(3)	0.07(5)	0.64 - 1.86	Ph72	
	FB	0.050(5)				0.56 - 1.96	Dr75	
^{204}Pb - $^{203}\text{Tl}^*$	FB	0.015(12)				0.51 - 1.49	Eu78	
^{206}Pb - $^{205}\text{Tl}^*$	FB	0.029(12)				0.51 - 1.49	Eu78	
^{209}Bi - $^{208}\text{Pb}^*$	3pG	0.021(2)	0.029	-0.007	0.026	0.7 - 2.8	Si73b	3,5
	FB	0.021(12)				0.51 - 1.49	Eu78	

*) Additional information can be found in the following references:

^{40}Ca - ^{40}Ar : We74
 ^{40}Ca - ^{39}K : Si73a
 ^{58}Ni - ^{56}Fe : Ha57

^{204}Pb - ^{203}Tl : Eu76a
 ^{206}Pb - ^{205}Tl : Eu76a
 ^{209}Bi - ^{208}Pb : Eu77b

Remarks

- 1) The value of Δz was fixed in the analysis.
- 2) Analysis performed in the High Energy Approximation (Pe66).
- 3) Data analyzed simultaneously with muonic X-ray data.
- 4) Only statistical errors are quoted.
- 5) The difference in the 0.8th moment of the charge distributions from muonic X-ray data was used as a constraint. A slightly better fit to the data was obtained by adding a $1h_{9/2}$ shell-model wave function to the 3pG distribution for ^{209}Bi .

TABLE IV. Fourier-Bessel Coefficients
See page 500 for Explanation of Tables

nucleus	^3H	^3He	^{12}C	^{12}C	^{15}N
rms [fm]	1.68(3)	1.877(19)	2.472(15)	2.464(12)	2.611(9)
a1	0.25182e-1	0.20020e-1	0.15721e-1	0.15737e-1	0.25491e-1
a2	0.34215e-1	0.41934e-1	0.38732e-1	0.38897e-1	0.50618e-1
a3	0.15257e-1	0.36254e-1	0.36808e-1	0.37085e-1	0.29822e-1
a4		0.17941e-1	0.14671e-1	0.14795e-1	-0.55196e-2
a5		0.46608e-2	-0.43277e-2	-0.44831e-2	-0.15913e-1
a6		0.46834e-2	-0.97752e-2	-0.10057e-1	-0.76184e-2
a7		0.52042e-2	-0.68908e-2	-0.68696e-2	-0.23992e-2
a8		0.38280e-2	-0.27631e-2	-0.28813e-2	-0.47940e-3
a9		0.25661e-2	-0.63568e-3	-0.77229e-3	
a10		0.14182e-2	0.71809e-4	0.66908e-4	
a11		0.61390e-3	0.18441e-3	0.10636e-3	
a12		0.22929e-3	0.75066e-4	-0.36864e-4	
a13			0.51069e-4	-0.50135e-5	
a14			0.14308e-4	0.94550e-5	
a15			0.23170e-5	-0.47687e-5	
a16			0.68465e-6		
a17					
ref. q-range [fm $^{-1}$] data- sets	Be84 0.51- 2.83	Re84b 0.18-10.10	Ca80 0.10- 4.01	Re82 0.25- 2.75	Vr86 0.22- 3.17
	Co65,Be82,Be84	Co65,MC77,Sz77, Ar78,Du83,Re84b	Si70b,Ja71,Ca80	Re82	Sc75,Vr86
R [fm]	3.5	5.0	8.0	8.0	7.0

nucleus	^{16}O	^{27}Al	^{28}Si	^{29}Si	^{30}Si
rms [fm]	2.737(8)	3.035(2)	3.085(17)	3.080(17)	3.173(25)
a1	0.20238e-1	0.43418e-1	0.33495e-1	0.33521e-1	0.28397e-1
a2	0.44793e-1	0.60298e-1	0.59533e-1	0.59679e-1	0.54163e-1
a3	0.33533e-1	0.28950e-2	0.20979e-1	0.20593e-1	0.25167e-1
a4	0.35030e-2	-0.23522e-1	-0.16900e-1	-0.18646e-1	-0.12858e-1
a5	-0.12293e-1	-0.79791e-2	-0.14998e-1	-0.16550e-1	-0.17592e-1
a6	-0.10329e-1	0.23010e-2	-0.93248e-3	-0.11922e-2	-0.46722e-2
a7	-0.34036e-2	0.10794e-2	0.33266e-2	0.28025e-2	0.24804e-2
a8	-0.41627e-3	0.12574e-3	0.59244e-3	-0.67353e-4	0.14760e-2
a9	-0.94435e-3	-0.13021e-3	-0.40013e-3	-0.34619e-3	-0.30168e-3
a10	-0.25771e-3	0.56563e-4	0.12242e-3	0.17611e-3	0.48346e-4
a11	0.23759e-3	-0.18011e-4	-0.12994e-4	-0.57173e-5	0.00000e0
a12	-0.10603e-3	0.42869e-5	-0.92784e-5	0.12371e-4	-0.51570e-5
a13	0.41480e-4		0.72595e-5		0.30261e-5
a14			-0.42096e-5		
a15					
a16					
a17					
ref. q-range [fm $^{-1}$] data- sets	La82 0.29- 2.77	Ro86 0.47-2.70	Mi82 0.25- 2.64	Mi82 0.25- 2.64	Mi82 0.25- 2.64
	La82	Ro86	Mi82	Mi82	Mi82
R [fm]	8.0	7.0	8.0	8.0	8.5

TABLE IV. Fourier-Bessel Coefficients
See page 500 for Explanation of Tables

nucleus	31P	32S	34S	36S	40Ar
rms [fm]	3.187(10)	3.248(4)	3.281(4)	3.278(6)	3.423(14)
a1	0.35305e-1	0.37251e-1	0.37036e-1	0.37032e-1	0.30451e-1
a2	0.59642e-1	0.60248e-1	0.58506e-1	0.57939e-1	0.55337e-1
a3	0.17274e-1	0.14748e-1	0.12082e-1	0.10049e-1	0.20203e-1
a4	-0.19303e-1	-0.18352e-1	-0.19022e-1	-0.19852e-1	-0.16765e-1
a5	-0.13545e-1	-0.10347e-1	-0.83421e-2	-0.67176e-2	-0.13578e-1
a6	0.63209e-3	0.30461e-2	0.45434e-2	0.61882e-2	-0.43204e-4
a7	0.35462e-2	0.35277e-2	0.28346e-2	0.37795e-2	0.91988e-3
a8	0.83653e-3	-0.39834e-4	-0.52304e-3	-0.55272e-3	-0.41205e-3
a9	-0.47904e-3	-0.97177e-4	0.27754e-4	-0.12904e-3	0.11971e-3
a10	0.19099e-3	0.92279e-4	0.59403e-4	0.15845e-3	-0.19801e-4
a11	-0.69611e-4	-0.51931e-4	-0.42794e-4	-0.84063e-4	-0.43204e-5
a12	0.23196e-4	0.22958e-4	0.20407e-4	0.34010e-4	0.61205e-5
a13	-0.77780e-5	-0.86609e-5	-0.79934e-5	-0.11663e-4	-0.37803e-5
a14	0.28879e-5	0.27354e-5	0.35204e-5	0.18001e-5	
a15		-0.86632e-6	-0.83914e-6	-0.95135e-6	-0.77407e-6
a16					
a17					
ref. q-range [fm ⁻¹] data- sets	Mi82 0.25- 2.64	Ry83b 0.47- 2.56	Ry83b 0.47- 2.56	Ry83b 0.47- 2.56	Ot82 0.29- 1.81
R [fm]	8.0	8.0	8.0	8.0	9.0

nucleus	40Ca	48Ca	48Ti	50Ti	50Cr
rms [fm]	3.450(10)	3.451(9)	3.597(1)	3.572(2)	3.662(4)
a1	0.44846e-1	0.44782e-1	0.27850e-1	0.31818e-1	0.39174e-1
a2	0.61326e-1	0.59523e-1	0.55432e-1	0.58556e-1	0.61822e-1
a3	-0.16818e-2	-0.74148e-2	0.26369e-1	0.19637e-1	0.68550e-2
a4	-0.26217e-1	-0.29466e-1	-0.17091e-1	-0.24309e-1	-0.30170e-1
a5	-0.29725e-2	-0.28350e-3	-0.21798e-1	-0.18748e-1	-0.98745e-2
a6	0.85534e-2	0.10829e-1	-0.24889e-2	0.33741e-2	0.87944e-2
a7	0.35322e-2	0.30465e-2	0.76631e-2	0.89961e-2	0.68502e-2
a8	-0.48258e-3	-0.10237e-2	0.34554e-2	0.37954e-2	-0.93609e-3
a9	-0.39346e-3	-0.17830e-3	-0.67477e-3	-0.41238e-3	-0.24962e-2
a10	0.20338e-3	0.55391e-4	0.10764e-3	0.12540e-3	-0.15361e-2
a11	0.25461e-4	-0.22644e-4	-0.16564e-5		-0.73687e-3
a12	-0.17794e-4	0.82671e-5	-0.55566e-5		
a13	0.67394e-5	-0.27343e-5			
a14	-0.21033e-5	0.82461e-6			
a15		-0.22780e-6			
a16					
a17					
ref. q-range [fm ⁻¹] data- sets	Em83b 0.35- 3.55	Em83b 0.35- 3.55	Se85 0.61- 2.20	Se85 0.61- 2.20	Li83c 0.15- 2.59
R [fm]	8.0	8.0	10.0	9.5	9.0

TABLE IV. Fourier-Bessel Coefficients
See page 500 for Explanation of Tables

nucleus	52Cr	54Cr	54Fe	56Fe	58Fe
rms [fm]	3.643(3)	3.689(4)	3.663(25)	3.714(24)	3.746(25)
a1	0.39287e-1	0.39002e-1	0.42339e-1	0.42018e-1	0.41791e-1
a2	0.62477e-1	0.60305e-1	0.64428e-1	0.62337e-1	0.60524e-1
a3	0.62482e-2	0.45845e-2	0.15840e-2	0.23995e-3	-0.14978e-2
a4	-0.32885e-1	-0.30723e-1	-0.35171e-1	-0.32776e-1	-0.31183e-1
a5	-0.10648e-1	-0.91355e-2	-0.10116e-1	-0.79941e-2	-0.58013e-2
a6	0.10520e-1	0.93251e-2	0.12069e-1	0.10844e-1	0.10611e-1
a7	0.85478e-2	0.60583e-2	0.62230e-2	0.49123e-2	0.41629e-2
a8	-0.24003e-3	-0.15602e-2	-0.12045e-2	-0.22144e-2	-0.29045e-5
a9	-0.20499e-2	-0.76809e-3	0.13561e-3	-0.18146e-3	0.54106e-3
a10	-0.12001e-2	0.76809e-3	0.10428e-4	0.37261e-3	-0.38689e-3
a11	-0.56649e-3	-0.34804e-3	-0.16980e-4	-0.23296e-3	0.20514e-3
a12			0.91817e-5	0.11494e-3	-0.95237e-4
a13			-0.39988e-5	-0.50596e-4	0.40707e-4
a14			0.15731e-5	0.20652e-4	-0.16346e-4
a15			-0.57862e-6	-0.79428e-5	0.62233e-5
a16			0.20186e-6	0.28986e-5	-0.22568e-5
a17			-0.67892e-7	-0.10075e-5	0.78077e-6
ref. q-range [fm ⁻¹] data- sets	Li83c 0.15- 2.59 La76,Li83c μ	Li83c 0.15- 2.59 La76,Li83c μ	Wo76 0.51- 2.22 Wo76	Wo76 0.51- 2.22 Wo76	Wo76 0.51- 2.22 Wo76
R [fm]	9.0	9.0	9.0	9.0	9.0

nucleus	59Co	58Ni	58Ni	60Ni	60Ni
rms [fm]	3.788(5)	3.769(13)	3.742(24)	3.797(13)	3.764(24)
a1	0.43133e-1	0.44880e-1	0.45030e-1	0.44668e-1	0.44855e-1
a2	0.61249e-1	0.64756e-1	0.65044e-1	0.63072e-1	0.63476e-1
a3	-0.32523e-2	-0.27899e-2	-0.32843e-2	-0.42797e-2	-0.51001e-2
a4	-0.32681e-1	-0.37016e-1	-0.36241e-1	-0.34806e-1	-0.34496e-1
a5	-0.49583e-2	-0.71915e-2	-0.67442e-2	-0.48625e-2	-0.43132e-2
a6	0.11494e-1	0.13594e-1	0.13146e-1	0.12794e-1	0.12767e-1
a7	0.55428e-2	0.66331e-2	0.50903e-2	0.54401e-2	0.49935e-2
a8	0.31398e-3	-0.14095e-2	-0.20787e-2	-0.14075e-2	-0.92940e-3
a9	-0.70578e-4	-0.10141e-2	0.12901e-3	-0.76976e-3	0.28281e-3
a10	0.53725e-5	0.38616e-3	0.14828e-3	0.33487e-3	-0.76557e-4
a11	-0.74650e-6	-0.13871e-3	-0.11530e-3	-0.13141e-3	0.18677e-4
a12	0.19793e-5	0.47788e-4	0.60881e-4	0.52132e-4	0.36855e-5
a13	-0.28059e-5	-0.15295e-4	-0.27676e-4	-0.20394e-4	-0.32276e-6
a14	0.27183e-5	0.59131e-5	0.11506e-4	0.59131e-5	0.19843e-6
a15	-0.19454e-5	-0.67880e-5	0.44764e-5	-0.67880e-5	0.16275e-6
a16	0.10963e-5		0.16468e-5		-0.82891e-7
a17	-0.51114e-6		-0.57496e-6		-0.34896e-7
ref. q-range [fm ⁻¹] data- sets	Sc77 0.51- 2.22 Sc77 μ	Be83 0.25- 2.65 Be83	Wo76 0.51- 2.22 Wo76	Be83 0.25- 2.65 Be83	Wo76 0.51- 2.22 Wo76
R [fm]	9.0	9.0	9.0	9.0	9.0

TABLE IV. Fourier-Bessel Coefficients
See page 500 for Explanation of Tables

nucleus	60Ni	62Ni	64Ni	63Cu	65Cu
rms [fm]	3.782(24)	3.802(24)	3.821(24)	3.885(5)	3.905(5)
a1	0.44742e-1	0.44581e-1	0.44429e-1	0.45598e-1	0.45444e-1
a2	0.62987e-1	0.61478e-1	0.60116e-1	0.60706e-1	0.59544e-1
a3	-0.49864e-2	-0.69425e-2	-0.92003e-2	-0.78616e-2	-0.94968e-2
a4	-0.34306e-1	-0.33126e-1	-0.33452e-1	-0.31638e-1	-0.31561e-1
a5	-0.44060e-2	-0.24964e-2	-0.52856e-3	-0.14447e-2	0.22898e-3
a6	0.12810e-1	0.12674e-1	0.13156e-1	0.10953e-1	0.11189e-1
a7	0.46914e-2	0.37148e-2	0.35152e-2	0.42578e-2	0.37360e-2
a8	-0.84373e-3	-0.20881e-2	-0.21671e-2	-0.24224e-3	-0.64873e-3
a9	0.36928e-3	0.30193e-3	0.46497e-4	-0.30067e-3	-0.51133e-3
a10	-0.15003e-3	0.57573e-4	0.25366e-3	0.23903e-3	0.43765e-3
a11	0.59665e-4	-0.77965e-4	-0.18438e-2	-0.12910e-3	-0.24276e-3
a12	-0.23215e-4	0.46906e-4	0.96874e-4	0.60195e-4	0.11507e-3
a13	0.88005e-5	-0.22724e-4	-0.44224e-4	-0.25755e-4	-0.49761e-4
a14	-0.32305e-5	0.98243e-5	0.18493e-4	0.10332e-4	0.20140e-4
a15	0.11496e-5	-0.39250e-5	-0.72361e-5	-0.39330e-5	-0.76945e-5
a16	-0.39658e-6	0.14732e-5	0.26740e-5	0.14254e-5	0.28055e-5
a17	0.13145e-6	0.52344e-6	-0.93929e-6	-0.49221e-6	-0.97411e-6
ref. q-range [fm ⁻¹] data- sets	Wo80 0.51- 2.22	Wo76 0.51- 2.22	Wo76 0.51- 2.22	Sc77 0.51- 2.22	Sc77 0.51- 2.22
R [fm]	9.0	9.0	9.0	9.0	9.0

nucleus	64Zn	66Zn	68Zn	70Zn	70Ge
rms [fm]	3.899(23)	3.903(25)	3.948(24)	3.987(31)	4.043(2)
a1	0.47038e-1	0.46991e-1	0.46654e-1	0.46362e-1	0.38182e-1
a2	0.61536e-1	0.60995e-1	0.58827e-1	0.57130e-1	0.60306e-1
a3	-0.90045e-2	-0.96693e-2	-0.12283e-1	-0.13877e-1	0.64346e-2
a4	-0.30669e-1	-0.30457e-1	-0.29865e-1	-0.30030e-1	-0.29427e-1
a5	-0.78705e-3	-0.53435e-3	0.25669e-2	0.35341e-2	-0.95888e-2
a6	0.10034e-1	0.97083e-2	0.10235e-1	0.10113e-1	0.87849e-2
a7	0.14053e-2	0.14091e-2	0.31861e-2	0.41029e-2	0.49187e-2
a8	-0.20640e-2	-0.70813e-3	-0.17351e-3	0.76469e-3	-0.15189e-2
a9	0.35105e-3	0.20809e-3	-0.42979e-3	-0.10138e-2	-0.17385e-2
a10	0.27303e-4	-0.48275e-4	0.33700e-3	0.60837e-3	-0.16794e-3
a11	-0.63811e-4	0.72680e-5	-0.18435e-3	-0.29929e-3	-0.11746e-3
a12	0.40893e-4	0.91369e-6	0.87043e-4	0.13329e-3	0.65768e-4
a13	-0.20311e-4	-0.14874e-5	-0.37612e-4	-0.55502e-4	-0.30691e-4
a14	0.88986e-5	0.88831e-6	0.15220e-4	0.21893e-4	0.13051e-5
a15	-0.35849e-5	-0.41689e-6	-0.58282e-5	-0.82286e-5	-0.52251e-5
a16	0.13522e-5	0.17283e-6	0.21230e-5	0.29559e-5	
a17	-0.38635e-6	-0.65968e-7	-0.73709e-6	-0.10148e-5	
ref. q-range [fm ⁻¹] data- sets	Wo76 0.51- 2.22	Wo76 0.51- 2.22	Wo76 0.51- 2.22	Wo76 0.51- 2.22	Ma84 0.35- 2.90
R [fm]	9.0	9.0	9.0	9.0	10.0

TABLE IV. Fourier-Bessel Coefficients
See page 500 for Explanation of Tables

nucleus	72Ge	74Ge	76Ge	88Sr	90Zr
rms [fm]	4.060(2)	4.075(2)	4.081(2)	4.197(6)	4.258(8)
a1	0.38083e-1	0.37989e-1	0.37951e-1	0.56435e-1	0.46188e-1
a2	0.59342e-1	0.58298e-1	0.57876e-1	0.55072e-1	0.61795e-1
a3	0.47718e-2	0.27406e-2	0.15303e-2	-0.33363e-1	-0.12315e-1
a4	-0.29953e-1	-0.30666e-1	-0.31822e-1	-0.26061e-1	-0.36915e-1
a5	-0.88476e-2	-0.81505e-2	-0.76875e-2	0.15749e-1	0.25175e-2
a6	0.96205e-2	0.10231e-1	0.11237e-1	0.75233e-2	0.15234e-1
a7	0.47901e-2	0.49382e-2	0.50780e-2	-0.55044e-2	-0.55146e-3
a8	-0.16869e-2	-0.16270e-2	-0.17293e-2	-0.23643e-2	-0.60631e-2
a9	-0.15406e-2	-0.13937e-2	-0.15523e-2	0.39362e-3	-0.12198e-2
a10	-0.97230e-4	0.15476e-3	0.72439e-4	-0.22733e-3	0.36200e-3
a11	-0.47640e-4	0.14396e-3	0.16560e-3	0.12519e-3	-0.16466e-3
a12	-0.15669e-5	-0.73075e-4	-0.86631e-4	-0.61176e-4	0.53305e-4
a13	0.67076e-5	0.31998e-4	0.39159e-4	0.27243e-4	-0.50873e-5
a14	-0.44500e-5	-0.12822e-4	-0.16259e-4	-0.11285e-4	-0.85658e-5
a15	0.22158e-5	0.48406e-5	0.63681e-5	0.43997e-5	0.86095e-5
a16				-0.16248e-5	
a17				0.57053e-6	
ref. q-range [fm ⁻¹] data- sets	Ma84 0.35- 2.90	Ma84 0.35- 2.90	Ma84 0.35- 2.90	St76 0.50- 2.50	Ro76 0.50- 2.50
R [fm]	10.0	10.0	10.0	9.0	10.0

nucleus	92Zr	94Zr	92Mo	94Mo	96Mo
rms [fm]	4.295(11)	4.315(10)	4.294(16)	4.333(16)	4.364(16)
a1	0.45939e-1	0.45798e-1	0.30782e-1	0.30661e-1	0.30564e-1
a2	0.60104e-1	0.59245e-1	0.59896e-1	0.58828e-1	0.58013e-1
a3	-0.13341e-1	-0.13389e-1	0.22016e-1	0.20396e-1	0.19255e-1
a4	-0.35106e-1	-0.33252e-1	-0.28945e-1	-0.28830e-1	-0.28372e-1
a5	0.31760e-2	0.39888e-2	-0.26707e-1	-0.25077e-1	-0.23304e-1
a6	0.13753e-1	0.12750e-1	0.40426e-2	0.44768e-2	0.49894e-2
a7	-0.82682e-3	-0.15793e-2	0.14429e-1	0.13127e-1	0.12126e-1
a8	-0.53001e-2	-0.56692e-2	0.31696e-2	0.19548e-2	0.10496e-2
a9	-0.97579e-3	-0.15698e-2	-0.63061e-2	-0.61403e-2	-0.62592e-2
a10	0.26489e-3	0.54394e-4	-0.45119e-2	-0.35825e-2	-0.32814e-2
a11	-0.15873e-3	-0.24032e-4	0.46236e-3	0.73790e-3	0.89668e-3
a12	0.69301e-4	0.38401e-4	0.94909e-3	0.61882e-3	0.50636e-3
a13	-0.22278e-4	-0.31690e-4	-0.38930e-3	-0.40556e-3	-0.43412e-3
a14	0.39533e-5	0.18481e-4	-0.14808e-3	-0.55748e-5	0.71531e-4
a15	0.10609e-5	-0.85367e-5	0.19622e-3	-0.12453e-3	0.76745e-4
a16			-0.40197e-4	-0.57812e-4	-0.54316e-4
a17			-0.71949e-4	-0.21657e-4	0.23386e-6
ref. q-range [fm ⁻¹] data- sets	Ro76 0.50- 2.50	Ro76 0.50- 2.50	Dr75 0.56- 1.96	Dr75 0.56- 1.96	Dr75 0.56- 1.96
R [fm]	10.0	10.0	12.0	12.0	12.0

TABLE IV. Fourier-Bessel Coefficients

See page 500 for Explanation of Tables

nucleus	98Mo	100Mo	104Pd	106Pd	108Pd
rms [fm]	4.388(16)	4.430(16)	4.437(10)	4.467(11)	4.524(10)
a1	0.30483e-1	0.30353e-1	0.41210e-1	0.41056e-1	0.40754e-1
a2	0.57207e-1	0.56087e-1	0.62846e-1	0.61757e-1	0.59460e-1
a3	0.17888e-1	0.16057e-1	-0.21202e-2	-0.29891e-2	-0.54077e-2
a4	-0.28388e-1	-0.28767e-1	-0.38359e-1	-0.37356e-1	-0.36305e-1
a5	-0.21778e-1	-0.20683e-1	-0.44693e-2	-0.35348e-2	-0.21987e-2
a6	0.56780e-2	0.62429e-2	0.16656e-1	0.16085e-1	0.15418e-1
a7	0.11236e-1	0.11058e-1	0.36873e-2	0.28502e-2	0.25927e-2
a8	0.82176e-3	0.11502e-2	-0.57534e-2	-0.55764e-2	-0.52781e-2
a9	-0.50390e-2	-0.39395e-2	-0.32499e-2	-0.15433e-2	-0.19757e-2
a10	-0.23877e-2	-0.14978e-2	0.69844e-3	0.22281e-2	0.10339e-2
a11	0.71492e-3	0.76350e-3	0.16304e-2	0.13160e-2	0.22891e-3
a12	0.29839e-3	0.10554e-3	0.59882e-3	0.16508e-4	-0.33464e-3
a13	-0.31408e-3	-0.25658e-3			
a14	0.80177e-3	0.10964e-3			
a15	0.43682e-4	0.10015e-4			
a16	-0.51394e-4	-0.40341e-4			
a17	0.22293e-4	0.25744e-4			
ref. q-range [fm ⁻¹] data- sets	Dr75 0.56- 1.96	Dr75 0.56- 1.96	La86 0.29- 2.38	La86 0.29- 2.38	La86 0.29- 2.38
R [fm]	12.0	12.0	11.0	11.0	11.0

nucleus	110Pd	144Sm	148Sm	148Sm	150Sm
rms [fm]	4.540(10)	4.943(9)	4.977(8)	5.002(6)	5.042(6)
a1	0.40668e-1	0.74734e-1	0.70491e-1	0.73859e-1	0.73338e-1
a2	0.58793e-1	0.26145e-1	0.32601e-1	0.24023e-1	0.24626e-1
a3	-0.61375e-2	-0.63832e-1	-0.55421e-1	-0.59437e-1	-0.52773e-1
a4	-0.35983e-1	0.10432e-1	0.50111e-2	0.10761e-1	0.10582e-1
a5	-0.17447e-2	0.19183e-1	0.20216e-1	0.17022e-1	0.15353e-1
a6	0.14998e-1	-0.12572e-1	-0.85944e-2	-0.11401e-1	-0.95624e-2
a7	0.19994e-2	-0.39707e-2	-0.40106e-2	-0.18102e-2	-0.18804e-2
a8	-0.53170e-2	-0.18703e-2	0.19303e-2	0.93011e-3	-0.79019e-3
a9	-0.14289e-2	0.12602e-2	-0.49689e-3	0.98012e-3	0.10102e-2
a10	0.16033e-2	-0.11902e-2	-0.17040e-3	-0.12601e-2	-0.26606e-2
a11	0.31574e-3	-0.15703e-2		-0.17402e-2	-0.18304e-2
a12	-0.42195e-3				
a13					
a14					
a15					
a16					
a17					
ref. q-range [fm ⁻¹] data- sets	La86 0.29- 2.40	Mo81 0.60- 2.50	Ho80 0.33- 2.18	Mo81 0.60- 2.50	Mo81 0.60- 2.50
R [fm]	11.0	9.25	9.5	9.25	9.25

TABLE IV. Fourier-Bessel Coefficients
See page 500 for Explanation of Tables

nucleus	152Sm	152Sm	154Sm	154Gd	158Gd
rms [fm]	5.099(8)	5.087(6)	5.126(8)	5.124	5.172(6)
a1	0.56097e-1	0.72646e-1	0.55859e-1	0.63832e-1	0.57217e-1
a2	0.45123e-1	0.21824e-1	0.44002e-1	0.36983e-1	0.43061e-1
a3	-0.40310e-1	-0.54112e-1	-0.40342e-1	-0.48193e-1	-0.41996e-1
a4	-0.18171e-1	0.98321e-2	-0.17989e-1	-0.51046e-2	-0.17203e-1
a5	0.20515e-1	0.16213e-1	0.19817e-1	0.19805e-1	0.19933e-1
a6	0.49023e-2	-0.65614e-2	0.51643e-2	-0.82574e-3	0.51060e-2
a7	-0.67674e-2	0.53611e-2	-0.60212e-2	-0.46942e-2	-0.73665e-2
a8	-0.18927e-2	-0.14103e-2	-0.23127e-2		-0.20926e-2
a9	0.15333e-2	-0.99022e-3	0.47024e-3		0.21883e-2
a10		-0.23005e-2			
a11					
a12					
a13					
a14					
a15					
a16					
a17					
ref. q-range [fm ⁻¹] data- sets	Ho80 0.33- 2.18	Mo81 0.60- 2.50	Ho80 0.33- 2.18	He82 0.58- 2.17	Mu84 0.36- 2.17
R [fm]	10.5	9.25	10.5	10.0	10.5

nucleus	166Er	174Yb	175Lu	192Os	196Pt
rms [fm]	5.227(20)	5.415(30)	5.366(30)	5.412(4)	5.383(20)
a1	0.54426e-1	0.54440e-1	0.55609e-1	0.59041e-1	0.50218e-1
a2	0.47165e-1	0.40034e-1	0.42243e-1	0.41498e-1	0.53722e-1
a3	-0.38654e-1	-0.45606e-1	-0.45028e-1	-0.49900e-1	-0.35015e-1
a4	-0.19672e-1	-0.20932e-1	-0.19491e-1	-0.10183e-1	-0.34588e-1
a5	0.22092e-1	0.20455e-1	0.22514e-1	0.29067e-1	0.23564e-1
a6	0.78708e-2	0.27061e-2	0.38982e-2	-0.57382e-2	0.14340e-1
a7	-0.53005e-2	-0.60489e-2	-0.72395e-2	-0.92348e-2	-0.13270e-1
a8	0.50005e-3	-0.15918e-3	0.31822e-3	0.36170e-2	-0.51212e-2
a9	0.52005e-3	0.11938e-2	-0.23866e-3	0.28736e-2	0.56088e-2
a10	-0.35003e-3			0.24194e-3	0.14890e-2
a11	0.12001e-3			-0.16766e-2	-0.10928e-2
a12				0.73610e-3	0.55662e-3
a13					-0.50557e-4
a14					-0.19708e-3
a15					0.24016e-3
a16					
a17					
ref. q-range [fm ⁻¹] data- sets	Ca78 0.29- 2.28	Sa79 0.32- 2.33	Sa79 0.32- 2.33	Re84 0.60- 2.90	Bo83 0.34- 2.28
R [fm]	11.0	11.0	11.0	11.0	12.0

TABLE IV. Fourier-Bessel Coefficients
See page 500 for Explanation of Tables

nucleus	203Tl	205Tl	204Pb	206Pb	207Pb
rms [fm]	5.463(5)	5.470(5)	5.479(2)	5.490(2)	5.497(2)
a1	0.51568e-1	0.51518e-1	0.52102e-1	0.52019e-1	0.51981e-1
a2	0.51562e-1	0.51165e-1	0.51786e-1	0.51190e-1	0.51059e-1
a3	-0.39299e-1	-0.39559e-1	-0.39188e-1	-0.39459e-1	-0.39447e-1
a4	-0.30826e-1	-0.30118e-1	-0.29242e-1	-0.28405e-1	-0.28428e-1
a5	0.27491e-1	0.27600e-1	0.28992e-1	0.28862e-1	0.28988e-1
a6	0.10795e-1	0.10412e-1	0.11040e-1	0.10685e-1	0.10329e-1
a7	-0.15922e-1	-0.15725e-1	-0.14591e-1	-0.14550e-1	-0.14029e-1
a8	-0.25527e-2	-0.26546e-2	-0.94917e-3	-0.13519e-2	-0.46728e-3
a9	0.58548e-2	0.70184e-2	0.71349e-2	0.77624e-2	0.67984e-2
a10	0.19324e-3	0.82116e-3	0.24780e-3	-0.41882e-4	0.56905e-3
a11	-0.17925e-3	-0.51805e-3	-0.61656e-3	-0.97010e-3	-0.50430e-3
a12	0.14307e-3	0.32560e-3	0.42335e-3	0.69611e-3	0.32796e-3
a13	-0.91669e-4	-0.18670e-3	-0.25250e-3	-0.42410e-3	-0.19157e-3
a14	0.53497e-4	0.10202e-3	0.14106e-3	0.23857e-3	0.10565e-3
a15	-0.29492e-4	-0.53857e-4	-0.75446e-4	-0.12828e-3	-0.56200e-4
a16	0.15625e-4	0.27672e-4	0.39143e-4	0.66663e-4	0.29020e-4
a17	-0.80141e-5	-0.13873e-4	-0.19760e-4	-0.33718e-4	-0.14621e-4
ref. q-range [fm ⁻¹] data- sets	Eu78 0.51- 2.24				
R [fm]	12.0	12.0	12.0	12.0	12.0

nucleus	208Pb	208Pb	209Bi
rms [fm]	5.503(2)	5.499(1)	5.518(4)
a1	0.51936e-1	0.62732e-1	0.52448e-1
a2	0.50768e-1	0.38542e-1	0.50400e-1
a3	-0.39646e-1	-0.55105e-1	-0.41014e-1
a4	-0.28218e-1	-0.26990e-2	-0.27927e-1
a5	0.28916e-1	0.31016e-1	0.29587e-1
a6	0.98910e-2	-0.99486e-2	0.98017e-2
a7	-0.14388e-1	-0.93012e-2	-0.14930e-1
a8	-0.98262e-3	0.76653e-2	-0.31967e-3
a9	0.72578e-2	0.20885e-2	0.77252e-2
a10	0.82318e-3	-0.17840e-2	0.57533e-3
a11	-0.14823e-2	0.74876e-4	-0.82529e-3
a12	0.13245e-3	0.32278e-3	0.25728e-3
a13	-0.84345e-4	-0.11353e-3	-0.11043e-3
a14	0.48417e-4		0.51930e-4
a15	-0.26562e-4		-0.24767e-4
a16	0.14035e-4		0.11863e-4
a17	-0.71863e-5		-0.56554e-5
ref. q-range [fm ⁻¹] data- sets	Eu78 0.51- 2.24	Fr77b 0.44- 3.70	Eu78 0.51- 2.24
R [fm]	12.0	11.0	12.0

TABLE V. Sum-of-Gaussians Parameters
See page 500 for Explanation of Tables

nucleus	^3H		^3He		^4He		^{12}C		^{16}O	
rms [fm]	1.76(4)		1.83(5)		1.676(8)		2.469(6)		2.711	
i	R _i	Q _i	R _i	Q _i	R _i	Q _i	R _i	Q _i	R _i	Q _i
1	0.0	0.035952	0.0	0.000029	0.2	0.034724	0.0	0.016690	0.4	0.057056
2	0.2	0.027778	0.6	0.606482	0.6	0.430761	0.4	0.050325	1.1	0.195701
3	0.5	0.131291	1.0	0.066077	0.9	0.203166	1.0	0.128621	1.9	0.311188
4	0.8	0.221551	1.3	0.000023	1.4	0.192986	1.3	0.180515	2.2	0.224321
5	1.2	0.253691	1.8	0.204417	1.9	0.083866	1.7	0.219097	2.7	0.059946
6	1.6	0.072905	2.3	0.115236	2.3	0.033007	2.3	0.278416	3.3	0.135714
7	2.0	0.152243	2.7	0.000001	2.6	0.014201	2.7	0.058779	4.1	0.000024
8	2.5	0.051564	3.2	0.006974	3.1	0.000000	3.5	0.057817	4.6	0.013961
9	3.0	0.053023	4.1	0.000765	3.5	0.006860	4.3	0.007739	5.3	0.000007
10					4.2	0.000000	5.4	0.002001	5.6	0.000002
11					4.9	0.000438	6.7	0.000007	5.9	0.002096
12					5.2	0.000000			6.4	0.000002
ref. q-range [fm ⁻¹] data- sets	Ju85 0.55- 4.79		MC77 0.59- 4.47		Si82 0.14- 7.70		Si82 0.13-3.70		Si70b 0.29- 3.97	
RP [fm]	0.80		1.10		1.00		1.20		1.30	

nucleus	^{24}Mg		^{28}Si		^{32}S		^{39}K		^{40}Ca	
rms [fm]	3.027		3.121		3.258		3.427		3.480(3)	
i	R _i	Q _i	R _i	Q _i	R _i	Q _i	R _i	Q _i	R _i	Q _i
1	0.1	0.007372	0.4	0.033149	0.4	0.045356	0.4	0.043308	0.4	0.042870
2	0.6	0.061552	1.0	0.106452	1.1	0.067478	0.9	0.036283	1.2	0.056020
3	1.1	0.056984	1.9	0.206866	1.7	0.172560	1.7	0.110517	1.8	0.167853
4	1.5	0.035187	2.4	0.286391	2.5	0.324870	2.1	0.147676	2.7	0.317962
5	1.9	0.291692	3.2	0.250448	3.2	0.254889	2.6	0.189541	3.2	0.155450
6	2.6	0.228920	3.6	0.056944	4.0	0.101799	3.2	0.274173	3.6	0.161897
7	3.2	0.233532	4.1	0.016829	4.6	0.022166	3.7	0.117691	4.3	0.053763
8	4.1	0.074086	4.6	0.039630	5.0	0.002081	4.2	0.058273	4.6	0.032612
9	4.7	0.000002	5.1	0.000002	5.5	0.005616	4.7	0.000006	5.4	0.004803
10	5.2	0.010876	5.5	0.000938	6.3	0.000020	5.5	0.021380	6.3	0.004541
11	6.1	0.000002	6.0	0.000002	7.3	0.000020	5.9	0.000002	6.6	0.000015
12	7.0	0.000002	6.9	0.002366	7.7	0.003219	6.9	0.001145	8.1	0.002218
ref. q-range [fm ⁻¹] data- sets	Li74 0.74- 3.64		Li74 0.74- 3.71		Li74 0.74- 3.71		Si74 0.64- 3.43		Si79 0.53- 3.56	
RP [fm]	1.25		1.30		1.35		1.45		1.45	

TABLE V. Sum-of-Gaussians Parameters
See page 500 for Explanation of Tables

nucleus	^{48}Ca		^{58}Ni		^{116}Sn		^{124}Sn		^{205}Tl	
rms [fm]	3.460		3.772(4)		4.627(1)		4.677(1)		5.479	
i	R _i	Q _i	R _i	Q _i	R _i	Q _i	R _i	Q _i	R _i	Q _i
1	0.6	0.063035	0.5	0.035228	0.1	0.005727	0.1	0.004877	0.6	0.007818
2	1.1	0.011672	1.4	0.065586	0.7	0.009643	0.7	0.010685	1.1	0.022853
3	1.7	0.064201	2.2	0.174552	1.3	0.038209	1.3	0.030309	2.1	0.000084
4	2.1	0.203813	3.0	0.199916	1.8	0.009466	1.8	0.015857	2.6	0.105635
5	2.9	0.259070	3.4	0.232360	2.3	0.096665	2.3	0.088927	3.1	0.022340
6	3.4	0.307899	3.9	0.118496	3.1	0.097840	3.1	0.091917	3.8	0.059933
7	4.3	0.080585	4.2	0.099325	3.8	0.269373	3.8	0.257379	4.4	0.235874
8	5.2	0.008498	4.6	0.029860	4.8	0.396671	4.8	0.401877	5.0	0.000004
9	5.7	0.000025	5.2	0.044912	5.5	0.026390	5.5	0.053646	5.7	0.460292
10	6.2	0.000005	5.9	0.000232	6.1	0.048157	6.1	0.043193	6.8	0.081621
11	6.5	0.000004	6.6	0.000002	7.1	0.001367	7.1	0.001319	7.2	0.002761
12	7.4	0.001210	7.9	0.000010	8.1	0.000509	8.1	0.000036	8.6	0.000803
ref. q-range [fm ⁻¹] data- sets	Si74 0.49- 3.37		Ca80b 0.58- 3.80		Ca82a 0.36- 3.60		Ca82a 0.36- 3.60		Fr83 0.51- 2.99	
	Be67d,Fr68, Si73a μ		Fi70,Si75 μ		Fi72,Ca82a μ		Fi72,Ca82a μ		Eu78,Fr83 μ	
RP [fm]	1.45		1.45		1.60		1.60		1.70	

nucleus	^{206}Pb		^{208}Pb	
rms [fm]	5.490		5.503(2)	
i	R _i	Q _i	R _i	Q _i
1	0.6	0.010615	0.1	0.003845
2	1.1	0.021108	0.7	0.009724
3	2.1	0.000060	1.6	0.033093
4	2.6	0.102206	2.1	0.000120
5	3.1	0.023476	2.7	0.083107
6	3.8	0.065884	3.5	0.080869
7	4.4	0.226032	4.2	0.139957
8	5.0	0.000005	5.1	0.260892
9	5.7	0.459690	6.0	0.336013
10	6.8	0.086351	6.6	0.033637
11	7.2	0.004589	7.6	0.018729
12	8.6	0.000011	8.7	0.000020
ref. q-range [fm ⁻¹] data- sets	Fr83 0.51- 2.99		Fr77a 0.44- 3.70	
	Eu78,Fr83 μ		He69,Ni69, Eu76a,Fr77a μ	
RP [fm]	1.70		1.70	

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