wwPDB X-ray Structure Validation Summary Report (i)

Nov 19, 2023 – 10:54 PM JST

PDB ID	:	7C07
Title	:	Crystal structure of yeast U2AF1 complex bound to 5'-AAGGU RNA.
Authors	:	Yoshida, H.; Park, S.Y.; Urano, T.; Obayashi, E.
Deposited on	:	2020-04-30
Resolution	:	3.20 Å(reported)

This is a wwPDB X-ray Structure Validation Summary Report for a publicly released PDB entry.

We welcome your comments at *validation@mail.wwpdb.org*
A user guide is available at
https://www.wwpdb.org/validation/2017/XrayValidationReportHelp
with specific help available everywhere you see the (i) symbol.

The types of validation reports are described at
http://www.wwpdb.org/validation/2017/FAQs#types.

The following versions of software and data (see references (1)) were used in the production of this report:

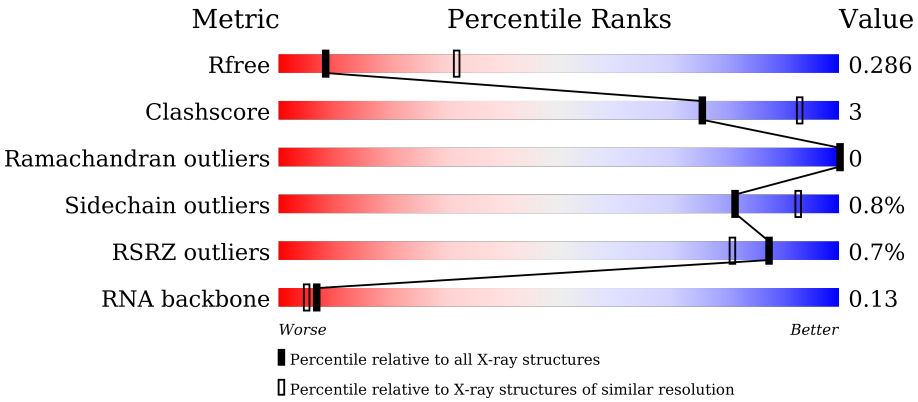
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.36
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 3.20 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



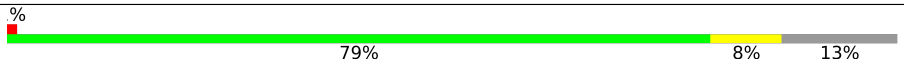
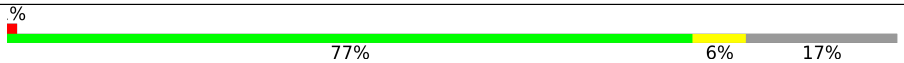
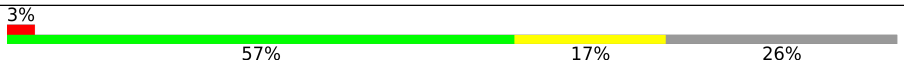
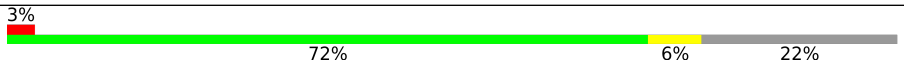
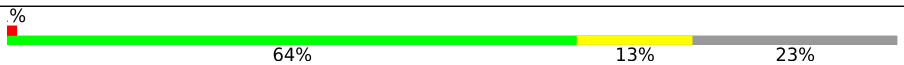
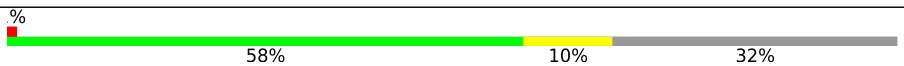
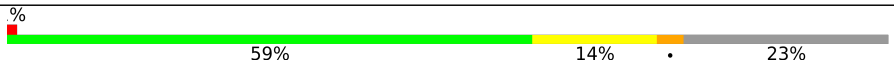
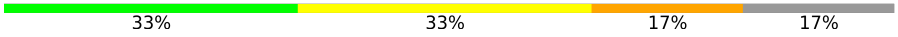
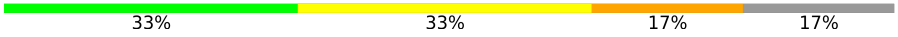
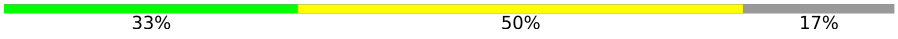
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	1133 (3.20-3.20)
Clashscore	141614	1253 (3.20-3.20)
Ramachandran outliers	138981	1234 (3.20-3.20)
Sidechain outliers	138945	1233 (3.20-3.20)
RSRZ outliers	127900	1095 (3.20-3.20)
RNA backbone	3102	1010 (3.50-2.90)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for >=3, 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions <=5% The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	216	84% 5% 11%
1	D	216	81% 5% 13%
1	G	216	81% 8% 11%
1	J	216	82% 7% 11%

Continued on next page...

Continued from previous page...

Mol	Chain	Length	Quality of chain
1	M	216	
1	P	216	
1	S	216	
1	V	216	
1	Y	216	
2	B	69	
2	E	69	
2	H	69	
2	K	69	
2	N	69	
2	Q	69	
2	T	69	
2	W	69	
2	Z	69	
3	1	6	
3	C	6	
3	F	6	
3	I	6	
3	L	6	
3	O	6	
3	R	6	
3	U	6	
3	X	6	

2 Entry composition (i)

There are 4 unique types of molecules in this entry. The entry contains 18390 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called Splicing factor U2AF 23 kDa subunit.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	193	Total	C	N	O	S	0	0	0
			1567	981	278	295	13			
1	D	187	Total	C	N	O	S	0	0	0
			1524	955	270	286	13			
1	G	193	Total	C	N	O	S	0	0	0
			1567	981	278	295	13			
1	J	193	Total	C	N	O	S	0	0	0
			1567	981	278	295	13			
1	M	179	Total	C	N	O	S	0	0	0
			1461	916	260	272	13			
1	P	189	Total	C	N	O	S	0	0	0
			1534	962	273	286	13			
1	S	188	Total	C	N	O	S	0	1	0
			1539	964	274	288	13			
1	V	174	Total	C	N	O	S	0	0	0
			1424	895	254	262	13			
1	Y	179	Total	C	N	O	S	0	0	0
			1461	916	260	272	13			

- Molecule 2 is a protein called Splicing factor U2AF 59 kDa subunit.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	56	Total	C	N	O	S	0	0	0
			455	288	82	84	1			
2	E	51	Total	C	N	O	S	0	0	0
			417	266	74	76	1			
2	H	53	Total	C	N	O	S	0	0	0
			430	274	76	79	1			
2	K	54	Total	C	N	O	S	0	0	0
			435	277	77	80	1			
2	N	52	Total	C	N	O	S	0	0	0
			426	272	75	78	1			

Continued on next page...

Continued from previous page...

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	Q	53	Total	C	N	O	S	0	0	0
			430	274	76	79	1			
2	T	47	Total	C	N	O	S	0	0	0
			383	246	67	69	1			
2	W	53	Total	C	N	O	S	0	0	0
			430	274	76	79	1			
2	Z	44	Total	C	N	O	S	0	0	0
			359	232	63	63	1			

- Molecule 3 is a RNA chain called RNA (5'-R(*U*AP*AP*GP*GP*U)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	5	Total	C	N	O	P	0	0	0
			107	49	22	32	4			
3	F	5	Total	C	N	O	P	0	0	0
			107	49	22	32	4			
3	I	5	Total	C	N	O	P	0	0	0
			107	49	22	32	4			
3	L	5	Total	C	N	O	P	0	0	0
			107	49	22	32	4			
3	O	5	Total	C	N	O	P	0	0	0
			107	49	22	32	4			
3	R	5	Total	C	N	O	P	0	0	0
			107	49	22	32	4			
3	U	5	Total	C	N	O	P	0	0	0
			107	49	22	32	4			
3	X	5	Total	C	N	O	P	0	0	0
			107	49	22	32	4			
3	1	5	Total	C	N	O	P	0	0	0
			107	49	22	32	4			

- Molecule 4 is ZINC ION (three-letter code: ZN) (formula: Zn) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	2	Total	Zn	0	0
			2	2		
4	D	2	Total	Zn	0	0
			2	2		
4	G	2	Total	Zn	0	0
			2	2		
4	J	2	Total	Zn	0	0
			2	2		

Continued on next page...

Continued from previous page...

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	M	2	Total Zn 2 2	0	0
4	P	2	Total Zn 2 2	0	0
4	S	2	Total Zn 2 2	0	0
4	V	2	Total Zn 2 2	0	0
4	Y	2	Total Zn 2 2	0	0

3 Residue-property plots (i)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density (RSRZ > 2). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

• Molecule 1: Splicing factor U2AF 23 kDa subunit

Chain A:

• Molecule 1: Splicing factor U2AF 23 kDa subunit

Chain D:

• Molecule 1: Splicing factor U2AF 23 kDa subunit

Chain G:

• Molecule 1: Splicing factor U2AF 23 kDa subunit

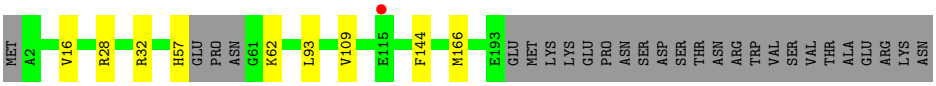
Chain J:

• Molecule 1: Splicing factor U2AF 23 kDa subunit

Chain M:

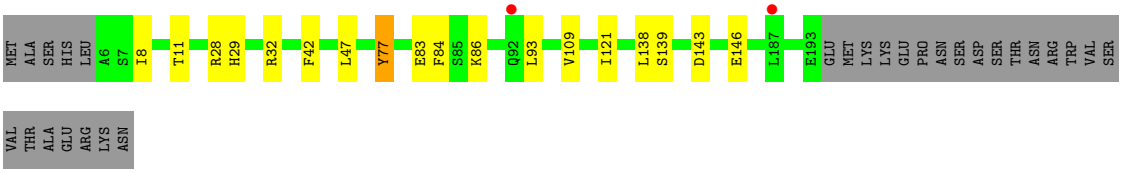
• Molecule 1: Splicing factor U2AF 23 kDa subunit

Chain P:



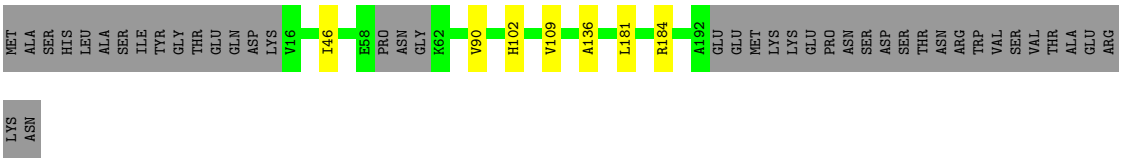
• Molecule 1: Splicing factor U2AF 23 kDa subunit

Chain S:



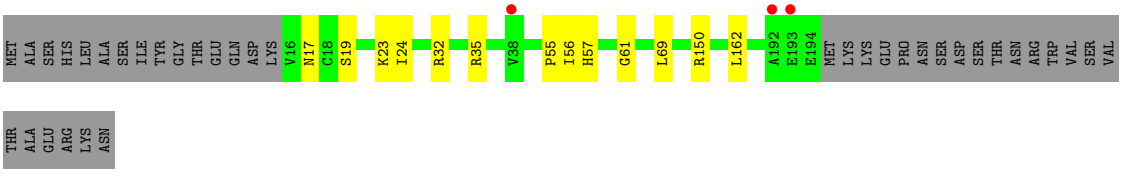
• Molecule 1: Splicing factor U2AF 23 kDa subunit

Chain V:



• Molecule 1: Splicing factor U2AF 23 kDa subunit

Chain Y:

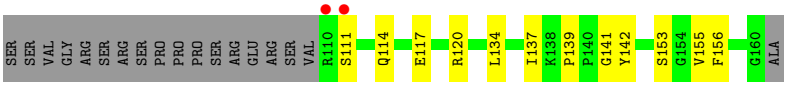


• Molecule 2: Splicing factor U2AF 59 kDa subunit

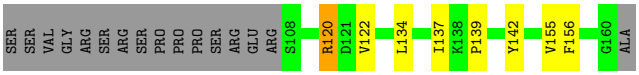
Chain B:

• Molecule 2: Splicing factor U2AF 59 kDa subunit

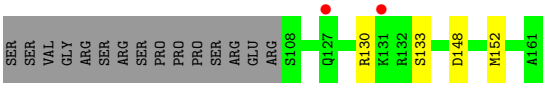
Chain E:



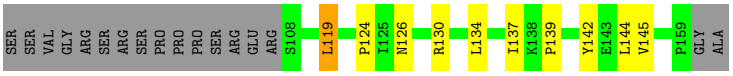
• Molecule 2: Splicing factor U2AF 59 kDa subunit



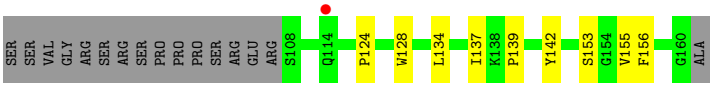
• Molecule 2: Splicing factor U2AF 59 kDa subunit



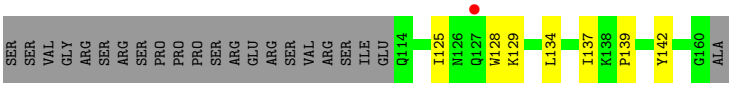
• Molecule 2: Splicing factor U2AF 59 kDa subunit



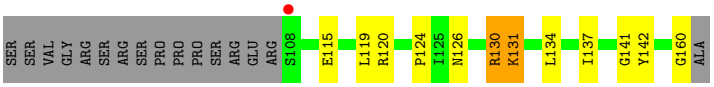
• Molecule 2: Splicing factor U2AF 59 kDa subunit



• Molecule 2: Splicing factor U2AF 59 kDa subunit



• Molecule 2: Splicing factor U2AF 59 kDa subunit

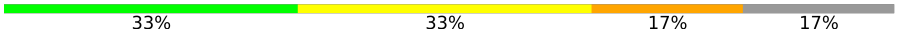


• Molecule 2: Splicing factor U2AF 59 kDa subunit

Chain Z:

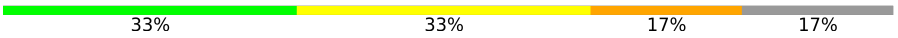
• Molecule 3: RNA (5'-R(*U*AP*AP*GP*GP*U)-3')

Chain C:



• Molecule 3: RNA (5'-R(*U*AP*AP*GP*GP*U)-3')

Chain F:

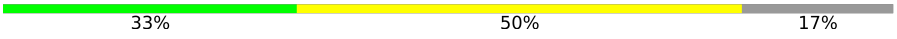


• Molecule 3: RNA (5'-R(*U*AP*AP*GP*GP*U)-3')

Chain I:

• Molecule 3: RNA (5'-R(*U*AP*AP*GP*GP*U)-3')

Chain L:



• Molecule 3: RNA (5'-R(*U*AP*AP*GP*GP*U)-3')

Chain O:

• Molecule 3: RNA (5'-R(*U*AP*AP*GP*GP*U)-3')

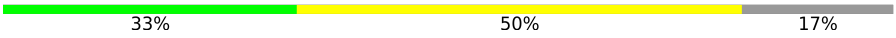
Chain R:

• Molecule 3: RNA (5'-R(*U*AP*AP*GP*GP*U)-3')

Chain U:

• Molecule 3: RNA (5'-R(*U*AP*AP*GP*GP*U)-3')

Chain X:



• Molecule 3: RNA (5'-R(*U*AP*AP*GP*GP*U)-3')

Chain 1:

4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α, β, γ	94.49Å 256.95Å 94.59Å 90.00° 100.75° 90.00°	Depositor
Resolution (Å)	48.17 – 3.20 48.17 – 3.20	Depositor EDS
% Data completeness (in resolution range)	100.0 (48.17-3.20) 100.0 (48.17-3.20)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.11 (at 3.19Å)	Xtriage
Refinement program	PHENIX 1.17.1_3660	Depositor
R, R_{free}	0.255 , 0.285 0.255 , 0.286	Depositor DCC
R_{free} test set	3656 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å^2)	73.6	Xtriage
Anisotropy	0.368	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.28 , 10.8	EDS
L-test for twinning[2]	$< \|L\| > = 0.46$, $< L^2 > = 0.29$	Xtriage
Estimated twinning fraction	0.039 for l,-k,h	Xtriage
F_o,F_c correlation	0.88	EDS
Total number of atoms	18390	wwPDB-VP
Average B, all atoms (Å^2)	80.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 14.49% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality (i)

5.1 Standard geometry (i)

Bond lengths and bond angles in the following residue types are not validated in this section: ZN

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	#$\|Z\|$ >5	RMSZ	#$\|Z\|$ >5
1	A	0.24	0/1604	0.39	0/2163
1	D	0.24	0/1560	0.40	0/2103
1	G	0.24	0/1604	0.39	0/2163
1	J	0.24	0/1604	0.39	0/2163
1	M	0.24	0/1496	0.40	0/2017
1	P	0.24	0/1569	0.38	0/2113
1	S	0.24	0/1576	0.40	0/2125
1	V	0.24	0/1457	0.38	0/1962
1	Y	0.24	0/1496	0.39	0/2017
2	B	0.23	0/465	0.37	0/629
2	E	0.24	0/427	0.39	0/578
2	H	0.24	0/440	0.40	0/596
2	K	0.23	0/445	0.38	0/603
2	N	0.23	0/436	0.43	0/591
2	Q	0.23	0/440	0.39	0/596
2	T	0.23	0/393	0.37	0/533
2	W	0.23	0/440	0.40	0/596
2	Z	0.23	0/368	0.39	0/497
3	1	0.25	0/120	0.87	0/186
3	C	0.21	0/120	0.79	0/186
3	F	0.19	0/120	0.77	0/186
3	I	0.22	0/120	0.84	0/186
3	L	0.23	0/120	0.86	0/186
3	O	0.23	0/120	0.86	0/186
3	R	0.21	0/120	0.83	0/186
3	U	0.18	0/120	0.82	0/186
3	X	0.23	0/120	0.87	0/186
All	All	0.24	0/18900	0.43	0/25719

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts (i)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1567	0	1487	7	0
1	D	1524	0	1448	8	0
1	G	1567	0	1487	16	0
1	J	1567	0	1488	9	0
1	M	1461	0	1389	10	0
1	P	1534	0	1461	8	0
1	S	1539	0	1459	13	0
1	V	1424	0	1360	5	0
1	Y	1461	0	1389	11	0
2	B	455	0	460	0	0
2	E	417	0	422	7	0
2	H	430	0	436	6	0
2	K	435	0	441	3	0
2	N	426	0	433	7	0
2	Q	430	0	436	5	0
2	T	383	0	387	4	0
2	W	430	0	436	10	0
2	Z	359	0	365	5	0
3	1	107	0	56	4	0
3	C	107	0	56	1	0
3	F	107	0	56	3	0
3	I	107	0	56	3	0
3	L	107	0	56	0	0
3	O	107	0	56	4	0
3	R	107	0	56	1	0
3	U	107	0	56	2	0
3	X	107	0	56	0	0
4	A	2	0	0	0	0
4	D	2	0	0	0	0
4	G	2	0	0	0	0
4	J	2	0	0	0	0
4	M	2	0	0	0	0
4	P	2	0	0	0	0

Continued on next page...

Continued from previous page...

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	S	2	0	0	0	0
4	V	2	0	0	0	0
4	Y	2	0	0	0	0
All	All	18390	0	17288	112	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 3.

The worst 5 of 112 close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:22:TYR:OH	1:G:184:ARG:NH2	2.22	0.71
1:D:33:CYS:HA	3:F:1:A:H62	1.57	0.70
1:Y:23:LYS:NZ	3:1:4:G:O2'	2.25	0.69
2:W:131:LYS:HD3	2:W:131:LYS:H	1.57	0.69
1:D:23:LYS:NZ	3:F:4:G:O2'	2.31	0.64

There are no symmetry-related clashes.

5.3 Torsion angles (i)

5.3.1 Protein backbone (i)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	191/216 (88%)	185 (97%)	6 (3%)	0	100	100
1	D	185/216 (86%)	181 (98%)	4 (2%)	0	100	100
1	G	191/216 (88%)	186 (97%)	5 (3%)	0	100	100
1	J	191/216 (88%)	183 (96%)	8 (4%)	0	100	100
1	M	177/216 (82%)	171 (97%)	6 (3%)	0	100	100
1	P	185/216 (86%)	179 (97%)	6 (3%)	0	100	100
1	S	187/216 (87%)	182 (97%)	5 (3%)	0	100	100

Continued on next page...

Continued from previous page...

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	V	170/216 (79%)	167 (98%)	3 (2%)	0	100	100
1	Y	177/216 (82%)	171 (97%)	6 (3%)	0	100	100
2	B	54/69 (78%)	52 (96%)	2 (4%)	0	100	100
2	E	49/69 (71%)	47 (96%)	2 (4%)	0	100	100
2	H	51/69 (74%)	48 (94%)	3 (6%)	0	100	100
2	K	52/69 (75%)	48 (92%)	4 (8%)	0	100	100
2	N	50/69 (72%)	47 (94%)	3 (6%)	0	100	100
2	Q	51/69 (74%)	49 (96%)	2 (4%)	0	100	100
2	T	45/69 (65%)	41 (91%)	4 (9%)	0	100	100
2	W	51/69 (74%)	48 (94%)	3 (6%)	0	100	100
2	Z	40/69 (58%)	36 (90%)	4 (10%)	0	100	100
All	All	2097/2565 (82%)	2021 (96%)	76 (4%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains (i)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	170/192 (88%)	170 (100%)	0	100	100
1	D	166/192 (86%)	165 (99%)	1 (1%)	86	94
1	G	170/192 (88%)	169 (99%)	1 (1%)	86	94
1	J	170/192 (88%)	168 (99%)	2 (1%)	71	88
1	M	159/192 (83%)	157 (99%)	2 (1%)	69	87
1	P	166/192 (86%)	166 (100%)	0	100	100
1	S	167/192 (87%)	166 (99%)	1 (1%)	86	94
1	V	155/192 (81%)	155 (100%)	0	100	100
1	Y	159/192 (83%)	159 (100%)	0	100	100
2	B	50/62 (81%)	49 (98%)	1 (2%)	55	80

Continued on next page...

Continued from previous page...

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
2	E	46/62 (74%)	46 (100%)	0	100 100
2	H	48/62 (77%)	47 (98%)	1 (2%)	53 79
2	K	48/62 (77%)	48 (100%)	0	100 100
2	N	48/62 (77%)	46 (96%)	2 (4%)	30 65
2	Q	48/62 (77%)	48 (100%)	0	100 100
2	T	42/62 (68%)	41 (98%)	1 (2%)	49 77
2	W	48/62 (77%)	45 (94%)	3 (6%)	18 52
2	Z	39/62 (63%)	39 (100%)	0	100 100
All	All	1899/2286 (83%)	1884 (99%)	15 (1%)	81 93

5 of 15 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	M	77	TYR
2	W	130	ARG
2	N	119	LEU
2	W	131	LYS
2	T	129	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

5.3.3 RNA (i)

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
3	1	5/6 (83%)	2 (40%)	1 (20%)
3	C	5/6 (83%)	2 (40%)	1 (20%)
3	F	4/6 (66%)	2 (50%)	0
3	I	5/6 (83%)	2 (40%)	1 (20%)
3	L	5/6 (83%)	2 (40%)	1 (20%)
3	O	5/6 (83%)	2 (40%)	1 (20%)
3	R	5/6 (83%)	1 (20%)	1 (20%)
3	U	5/6 (83%)	2 (40%)	1 (20%)
3	X	5/6 (83%)	2 (40%)	1 (20%)
All	All	44/54 (81%)	17 (38%)	8 (18%)

5 of 17 RNA backbone outliers are listed below:

Mol	Chain	Res	Type
3	C	4	G
3	C	5	U
3	F	2	A
3	F	4	G
3	I	4	G

5 of 8 RNA pucker outliers are listed below:

Mol	Chain	Res	Type
3	1	1	A
3	X	1	A
3	R	1	A
3	O	1	A
3	U	1	A

5.4 Non-standard residues in protein, DNA, RNA chains (i)

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates (i)

There are no monosaccharides in this entry.

5.6 Ligand geometry (i)

Of 18 ligands modelled in this entry, 18 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers (i)

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

In the following table, the column labelled '#RSRZ> 2' contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95^{th} percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled 'Q< 0.9' lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å^2)	Q<0.9
1	A	193/216 (89%)	-0.06	0 100 100	45, 68, 108, 136	0
1	D	187/216 (86%)	-0.08	0 100 100	42, 62, 102, 125	0
1	G	193/216 (89%)	-0.06	1 (0%) 91 86	33, 56, 99, 131	0
1	J	193/216 (89%)	0.07	1 (0%) 91 86	43, 72, 106, 130	0
1	M	179/216 (82%)	-0.03	0 100 100	49, 81, 121, 138	0
1	P	189/216 (87%)	0.03	1 (0%) 91 86	42, 67, 92, 122	0
1	S	188/216 (87%)	0.15	2 (1%) 80 69	54, 84, 125, 141	0
1	V	174/216 (80%)	0.00	0 100 100	70, 94, 119, 129	0
1	Y	179/216 (82%)	0.15	3 (1%) 70 57	43, 91, 128, 142	0
2	B	56/69 (81%)	-0.10	0 100 100	48, 72, 111, 119	0
2	E	51/69 (73%)	0.03	2 (3%) 39 25	48, 65, 144, 148	0
2	H	53/69 (76%)	0.01	0 100 100	26, 60, 115, 120	0
2	K	54/69 (78%)	0.07	2 (3%) 41 26	40, 69, 112, 136	0
2	N	52/69 (75%)	0.14	0 100 100	52, 78, 139, 148	0
2	Q	53/69 (76%)	0.12	1 (1%) 66 53	45, 69, 128, 145	0
2	T	47/69 (68%)	0.29	1 (2%) 63 49	66, 89, 150, 156	0
2	W	53/69 (76%)	0.19	1 (1%) 66 53	69, 94, 144, 148	0
2	Z	44/69 (63%)	0.03	0 100 100	48, 71, 124, 137	0
3	1	5/6 (83%)	0.02	0 100 100	105, 115, 126, 146	0
3	C	5/6 (83%)	-0.09	0 100 100	65, 80, 90, 93	0
3	F	5/6 (83%)	0.19	0 100 100	66, 68, 83, 118	0
3	I	5/6 (83%)	-0.13	0 100 100	55, 60, 82, 95	0
3	L	5/6 (83%)	-0.20	0 100 100	66, 87, 92, 101	0
3	O	5/6 (83%)	0.11	0 100 100	100, 108, 130, 142	0

Continued on next page...

Continued from previous page...

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2			OWAB(Å^2)	Q<0.9
3	R	5/6 (83%)	0.15	0	100	100	58, 60, 83, 90	0
3	U	5/6 (83%)	0.18	0	100	100	71, 72, 83, 145	0
3	X	5/6 (83%)	-0.47	0	100	100	105, 107, 112, 118	0
All	All	2183/2619 (83%)	0.03	15 (0%)	87	81	26, 76, 125, 156	0

The worst 5 of 15 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	J	8	ILE	3.5
1	Y	192	ALA	3.2
2	Q	114	GLN	3.2
2	W	108	SER	3.1
1	Y	193	GLU	2.9

6.2 Non-standard residues in protein, DNA, RNA chains (i)

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates (i)

There are no monosaccharides in this entry.

6.4 Ligands (i)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
4	ZN	J	301	1/1	0.96	0.20	67,67,67,67	0
4	ZN	G	301	1/1	0.98	0.24	48,48,48,48	0
4	ZN	A	302	1/1	0.98	0.22	55,55,55,55	0
4	ZN	J	302	1/1	0.98	0.21	54,54,54,54	0
4	ZN	S	301	1/1	0.98	0.23	59,59,59,59	0
4	ZN	S	302	1/1	0.98	0.21	73,73,73,73	0
4	ZN	A	301	1/1	0.99	0.18	60,60,60,60	0
4	ZN	M	301	1/1	0.99	0.20	70,70,70,70	0
4	ZN	M	302	1/1	0.99	0.18	65,65,65,65	0

Continued on next page...

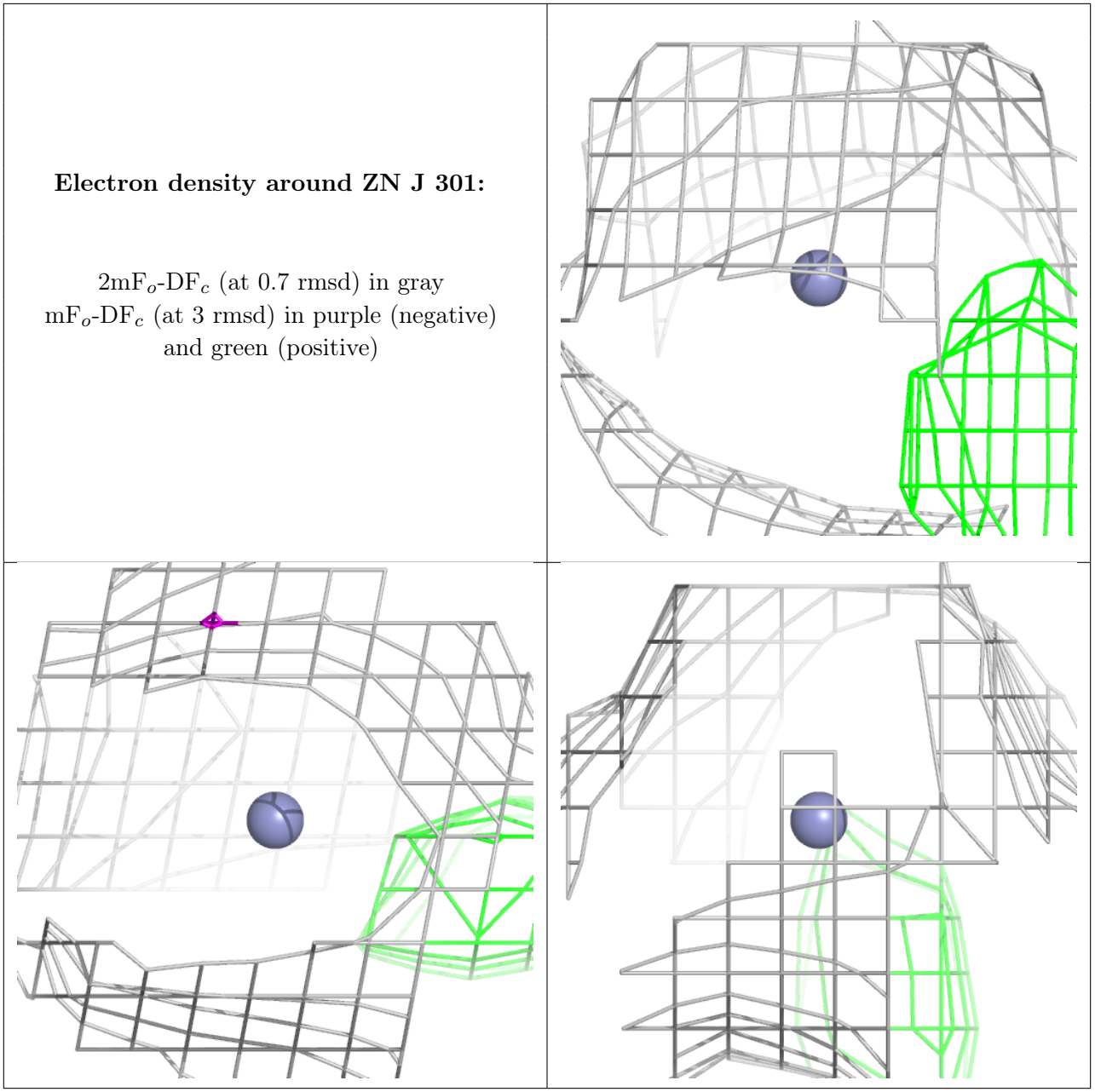
Continued from previous page...

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
4	ZN	P	301	1/1	0.99	0.23	44,44,44,44	0
4	ZN	P	302	1/1	0.99	0.23	45,45,45,45	0
4	ZN	G	302	1/1	0.99	0.22	54,54,54,54	0
4	ZN	D	302	1/1	0.99	0.20	49,49,49,49	0
4	ZN	V	301	1/1	0.99	0.23	72,72,72,72	0
4	ZN	V	302	1/1	0.99	0.17	71,71,71,71	0
4	ZN	Y	301	1/1	0.99	0.19	80,80,80,80	0
4	ZN	Y	302	1/1	0.99	0.17	70,70,70,70	0
4	ZN	D	301	1/1	1.00	0.23	45,45,45,45	0

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

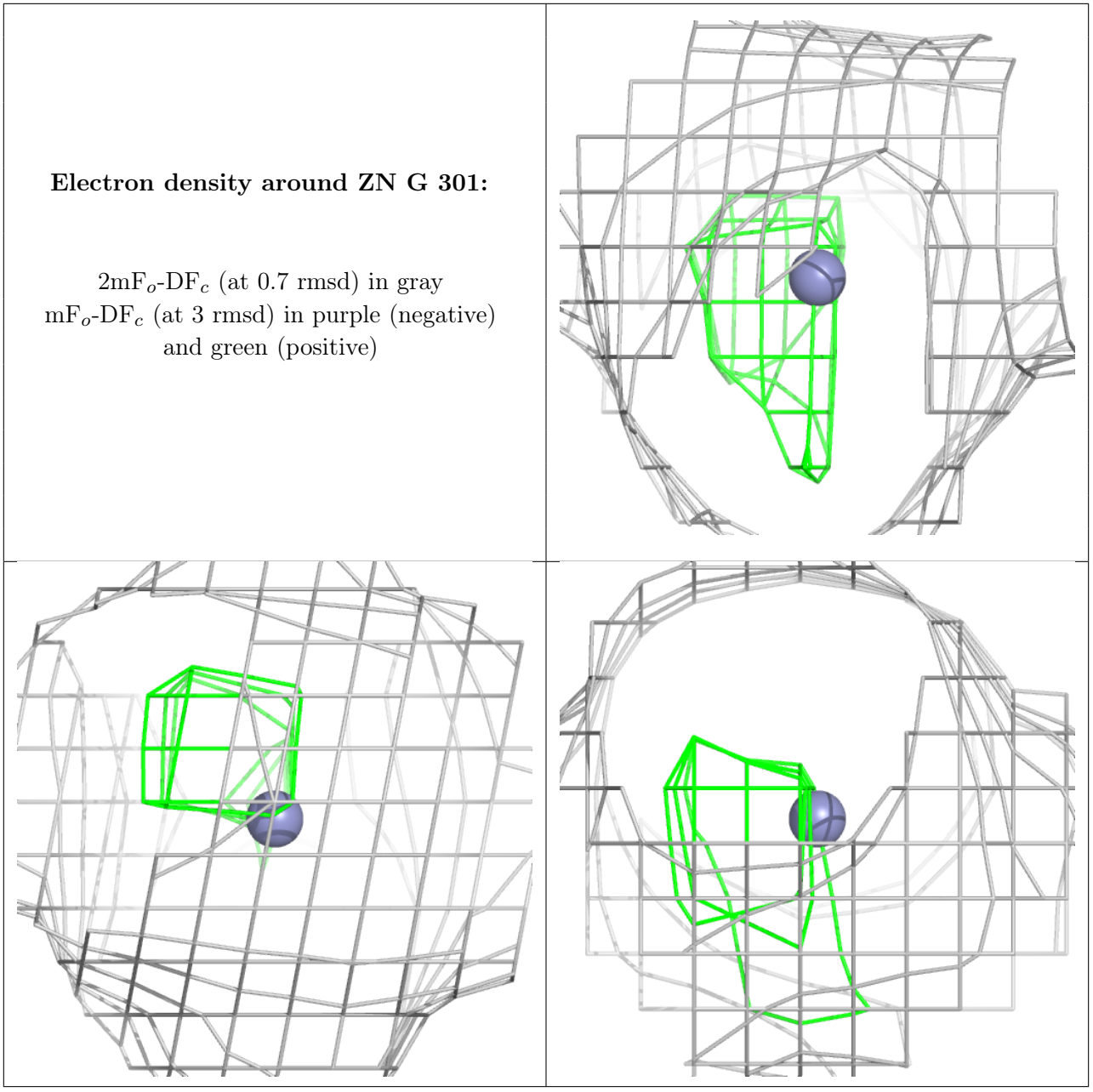
Electron density around ZN J 301:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)



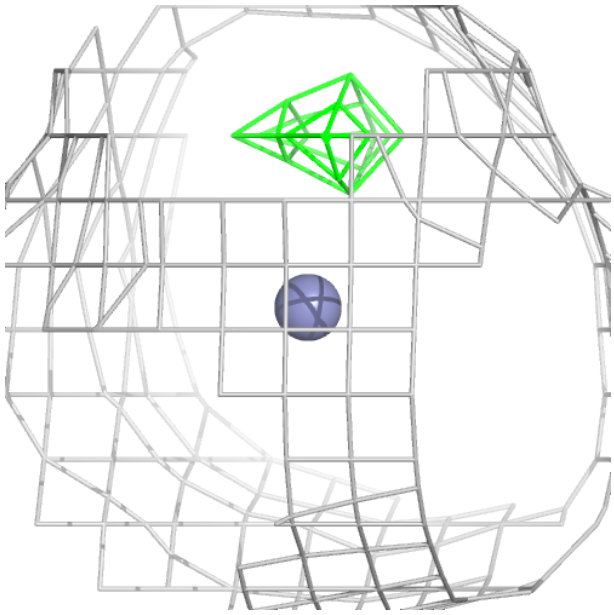
Electron density around ZN G 301:

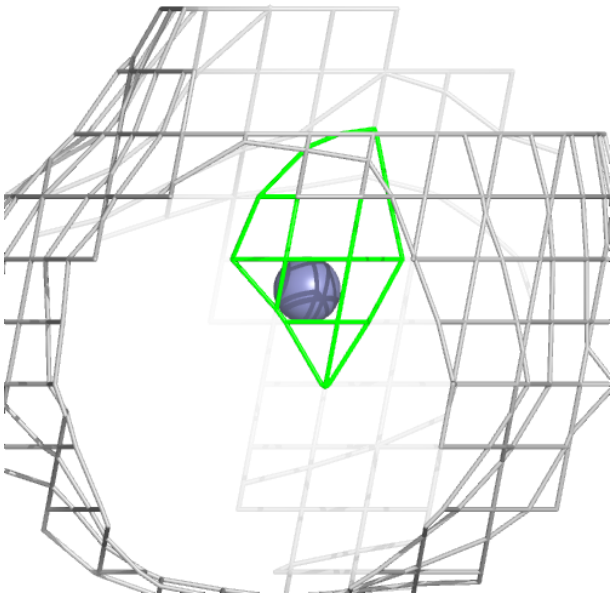
$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

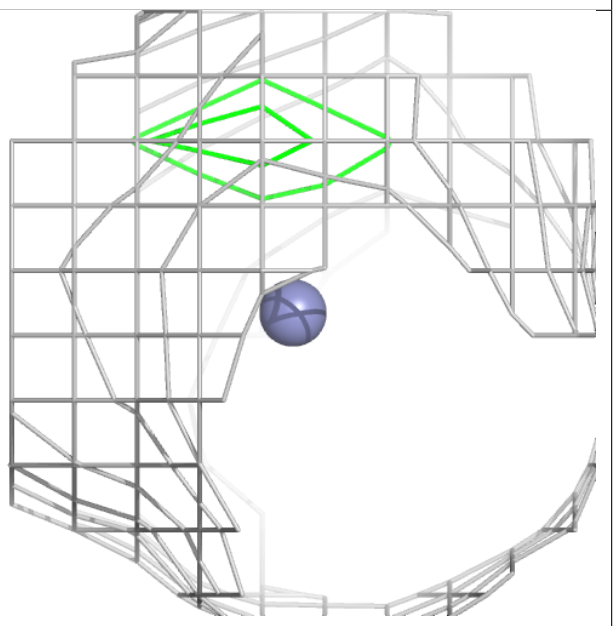


Electron density around ZN A 302:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

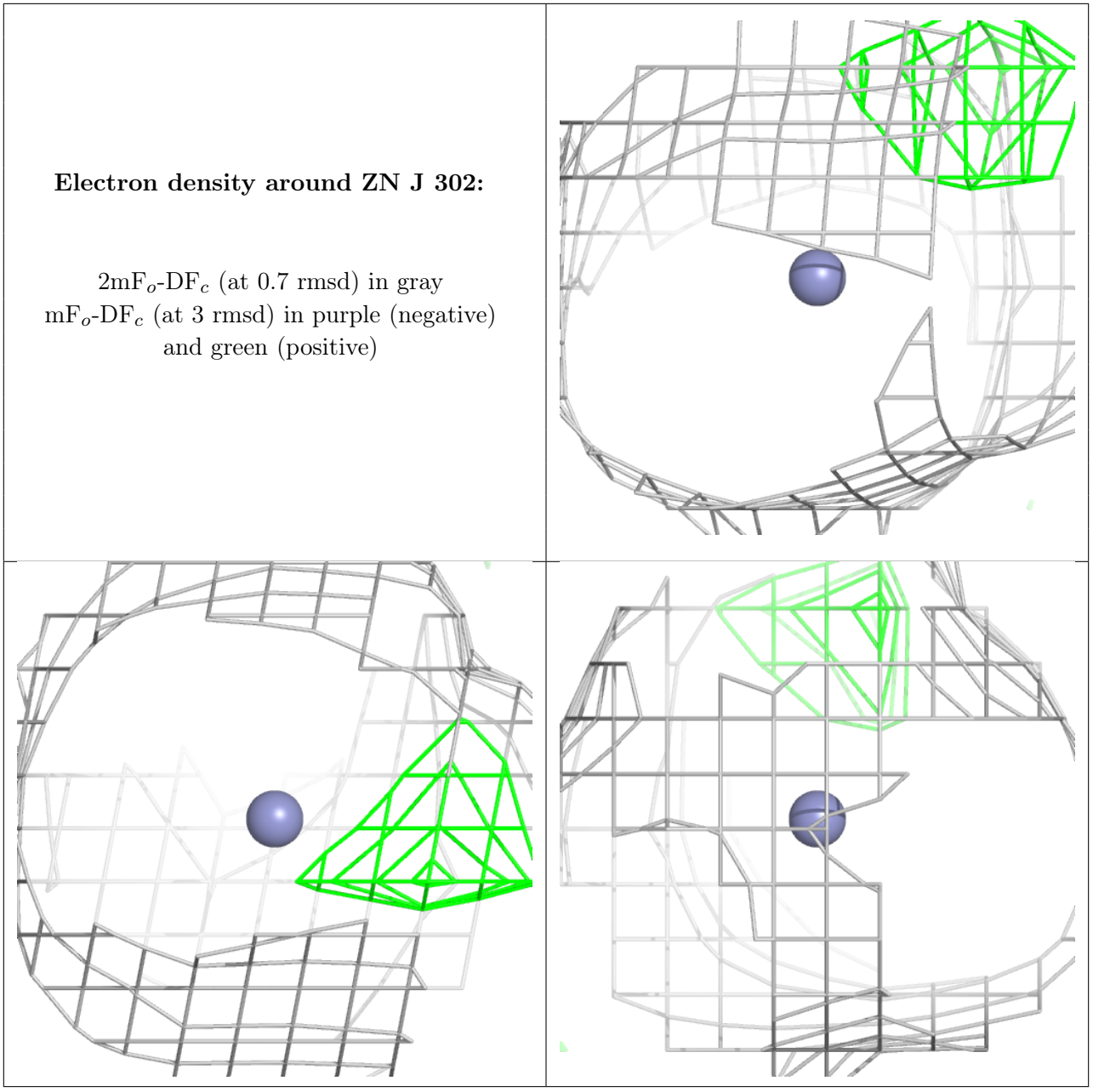






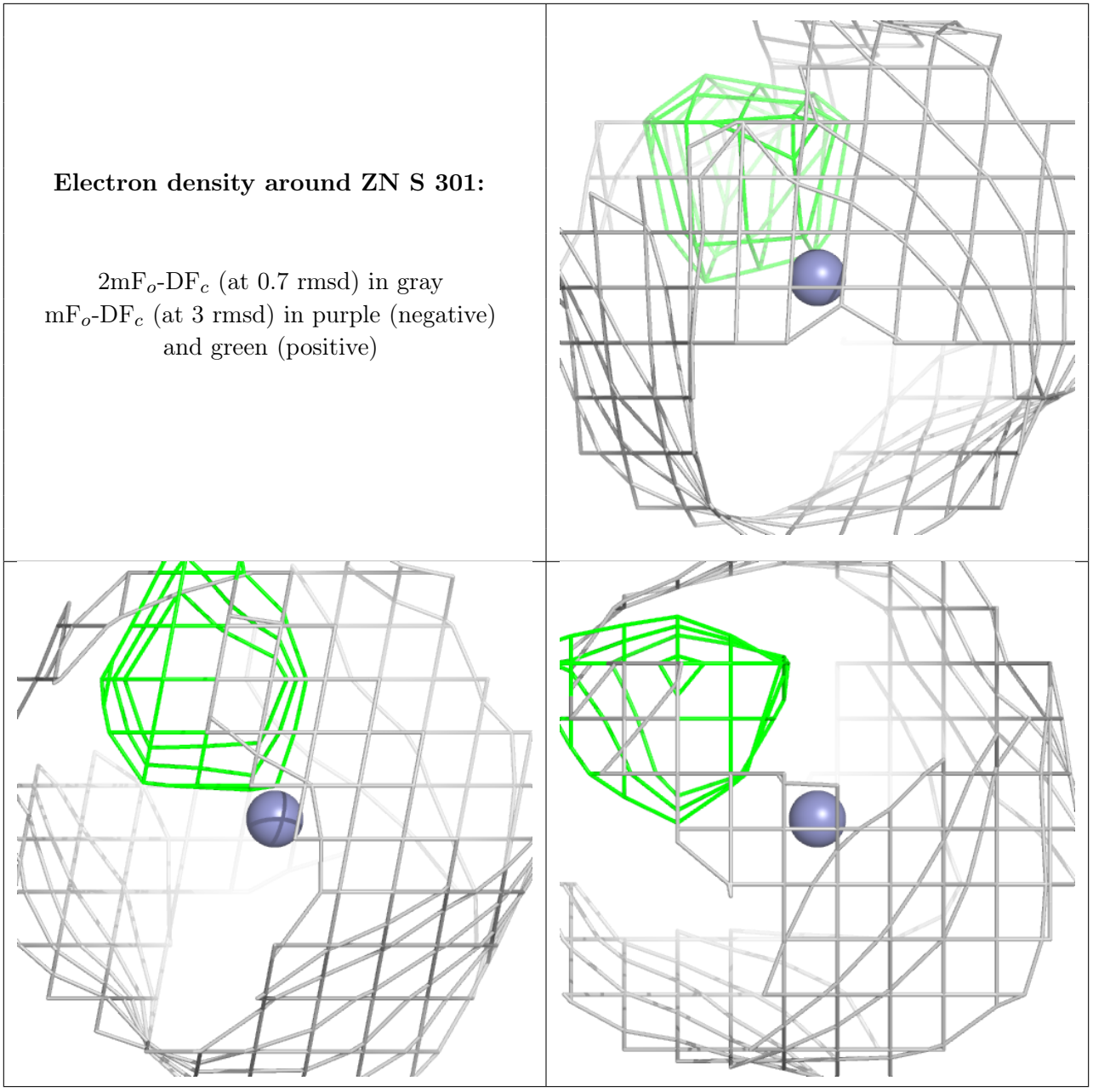
Electron density around ZN J 302:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)



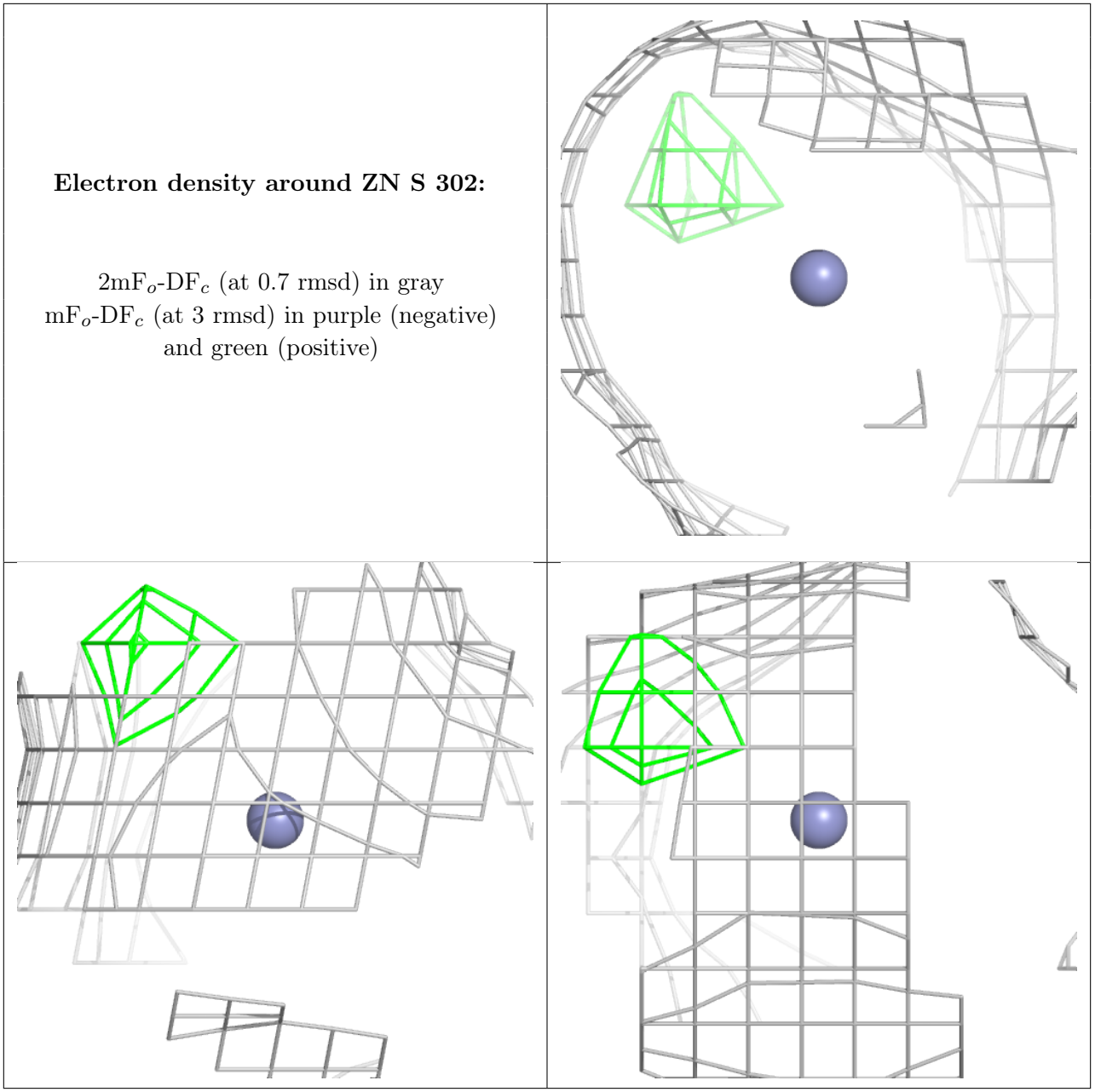
Electron density around ZN S 301:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)



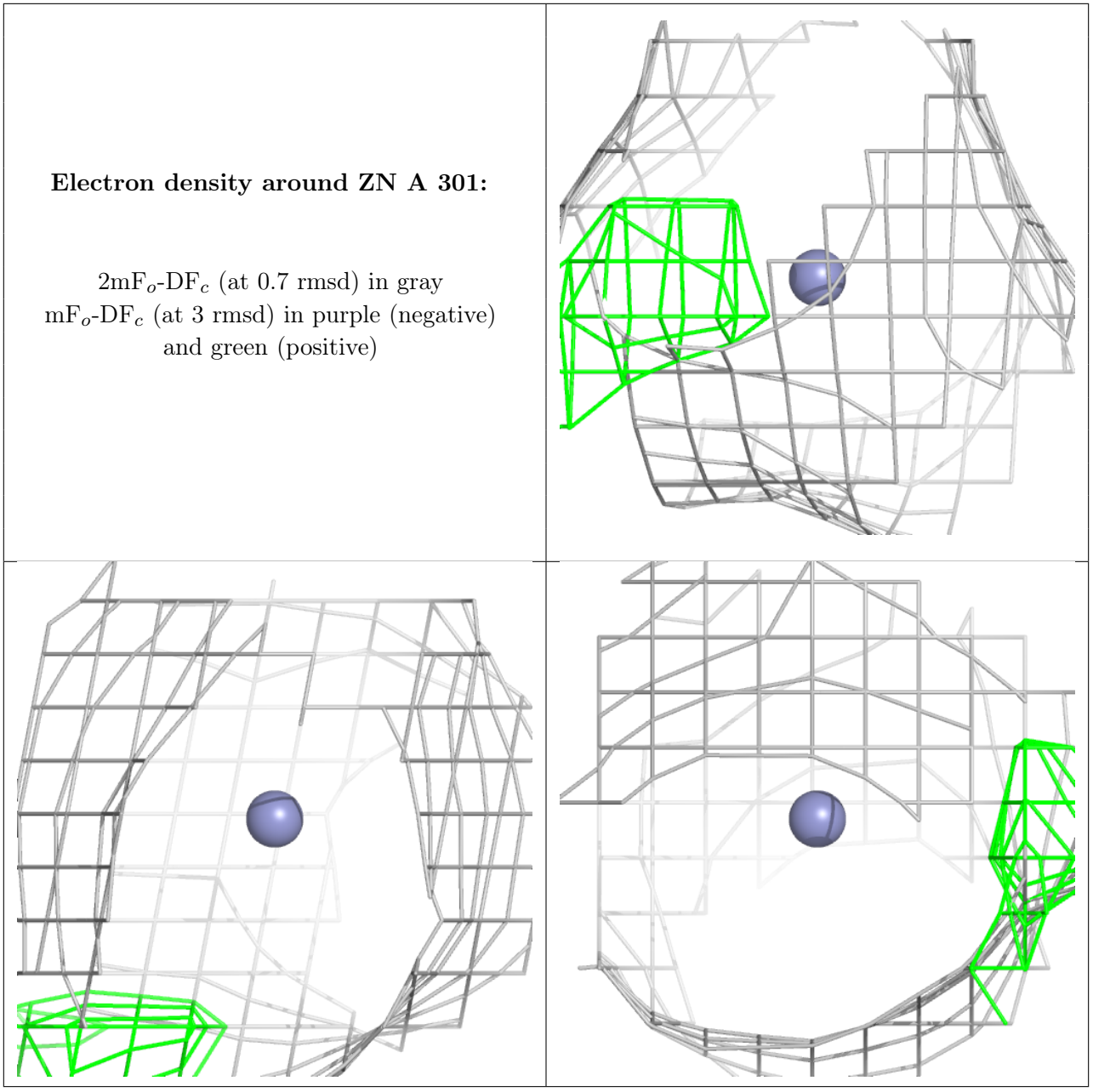
Electron density around ZN S 302:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)



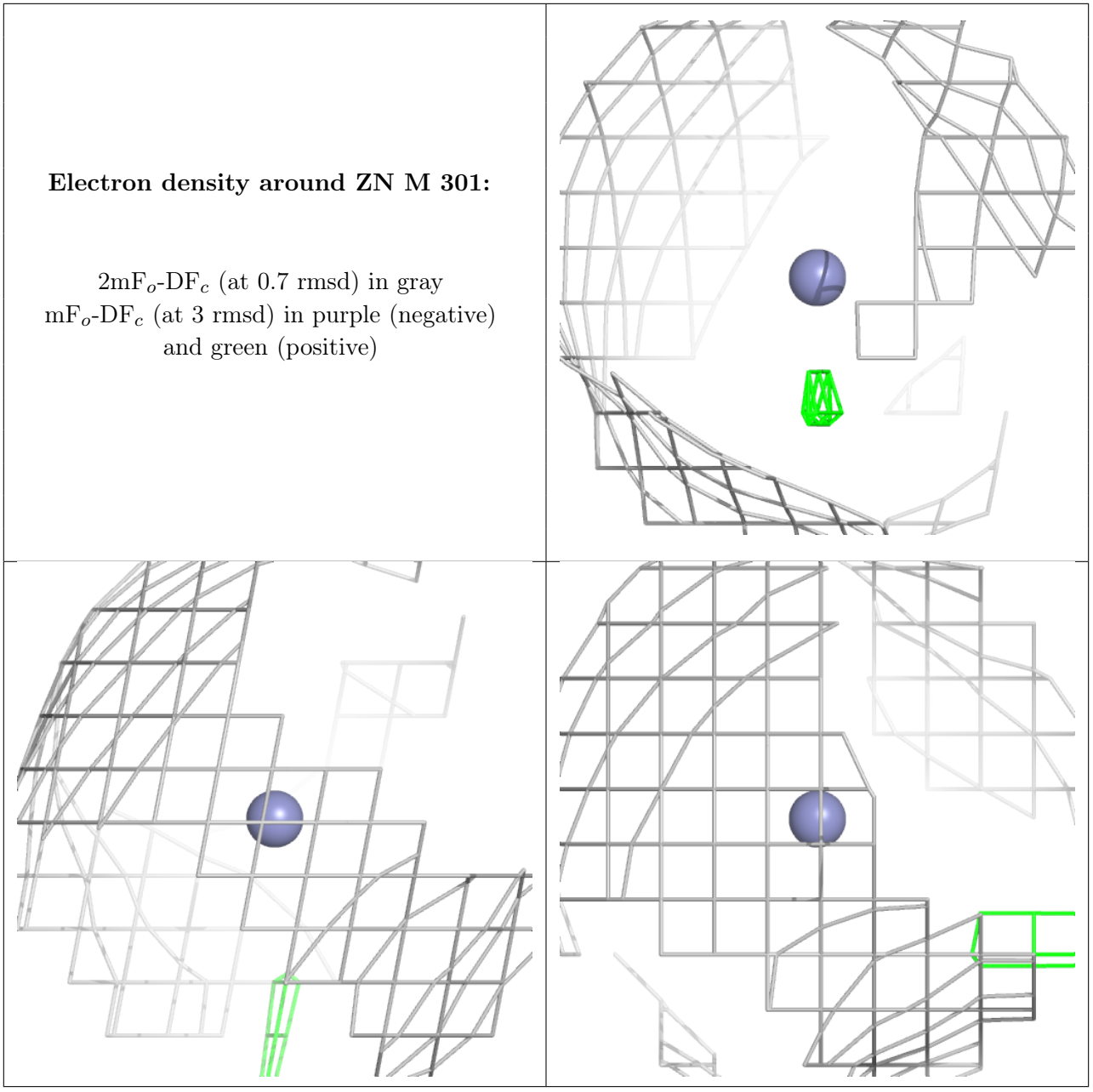
Electron density around ZN A 301:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)



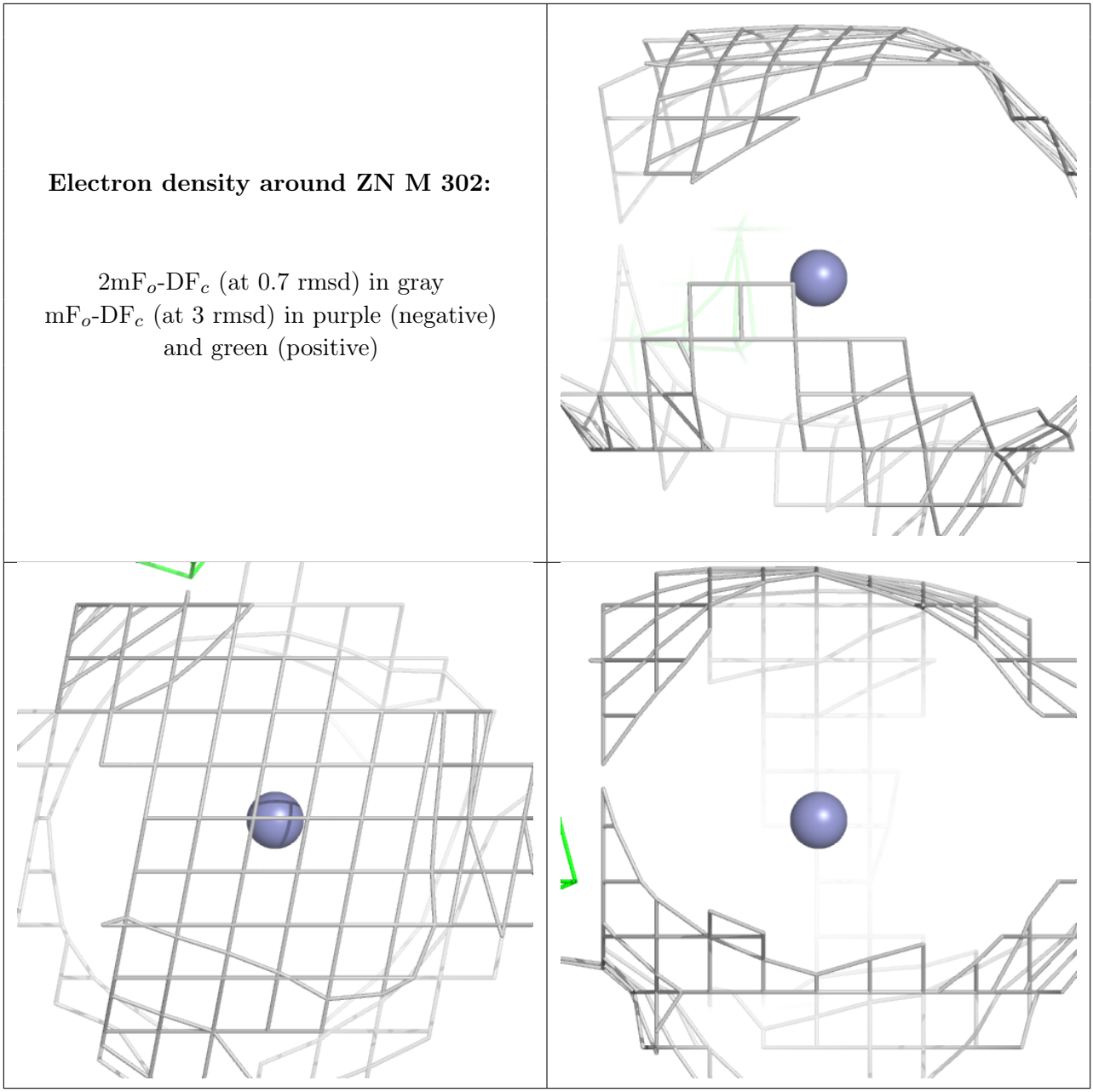
Electron density around ZN M 301:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)



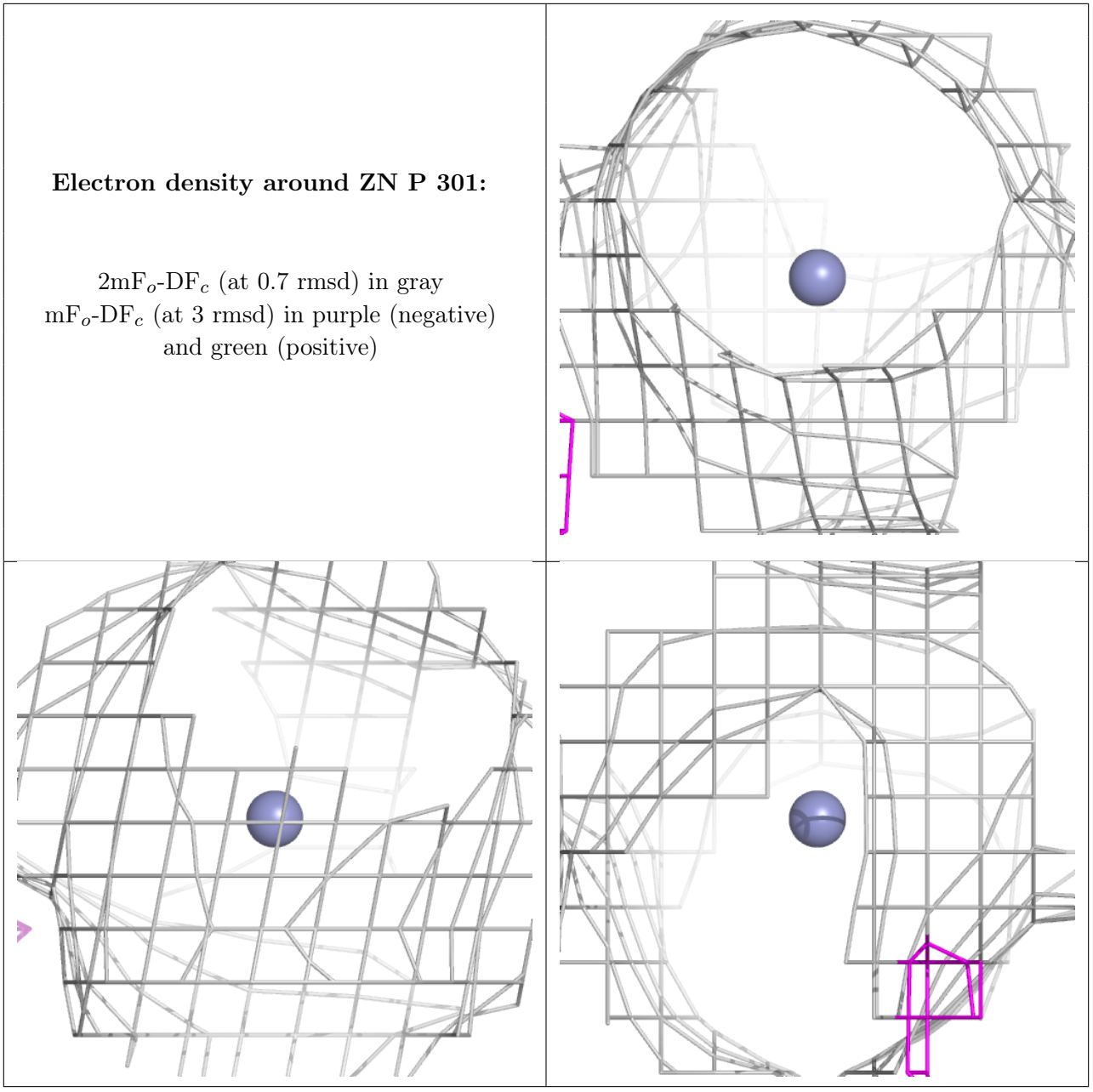
Electron density around ZN M 302:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)



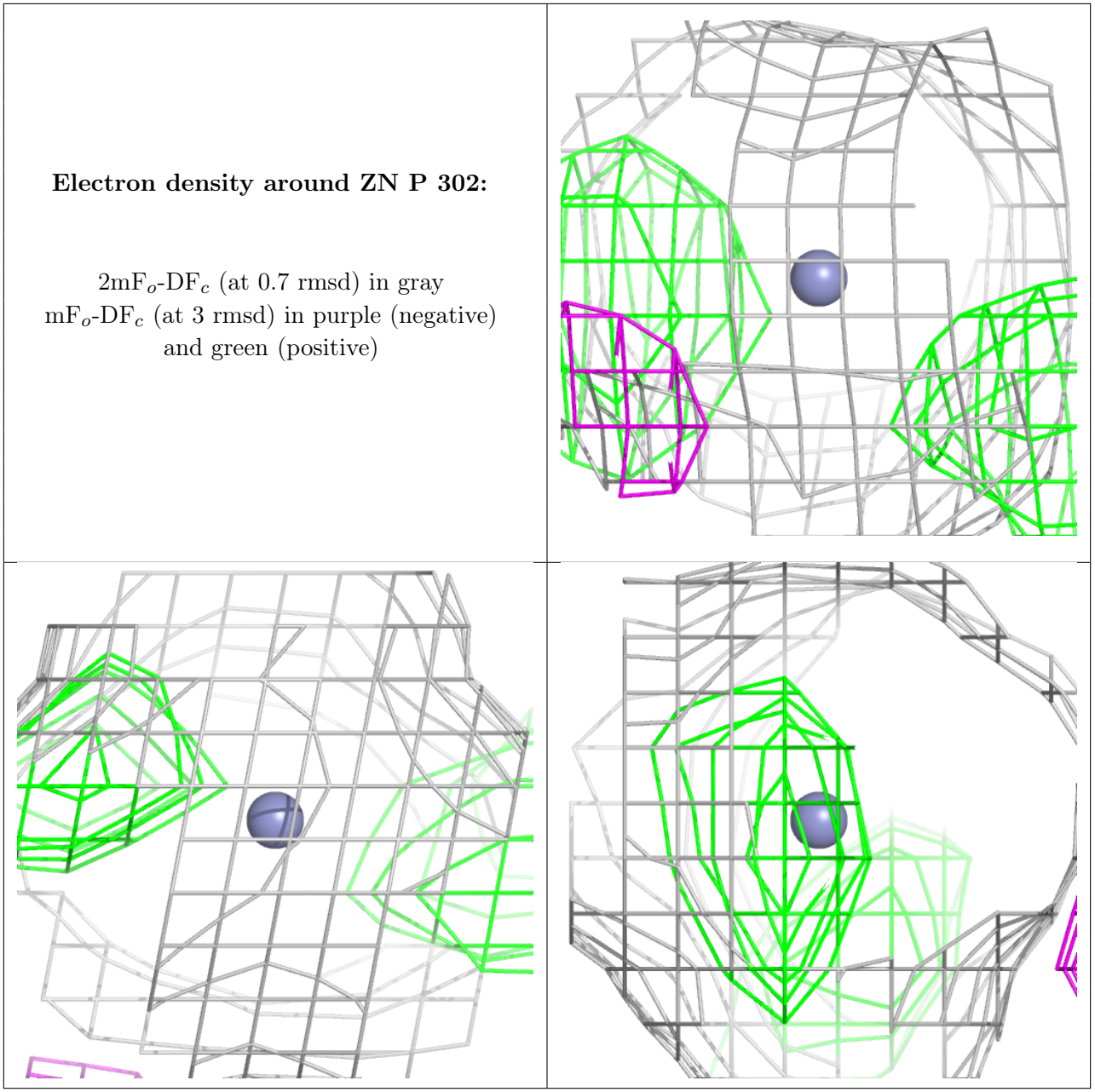
Electron density around ZN P 301:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)



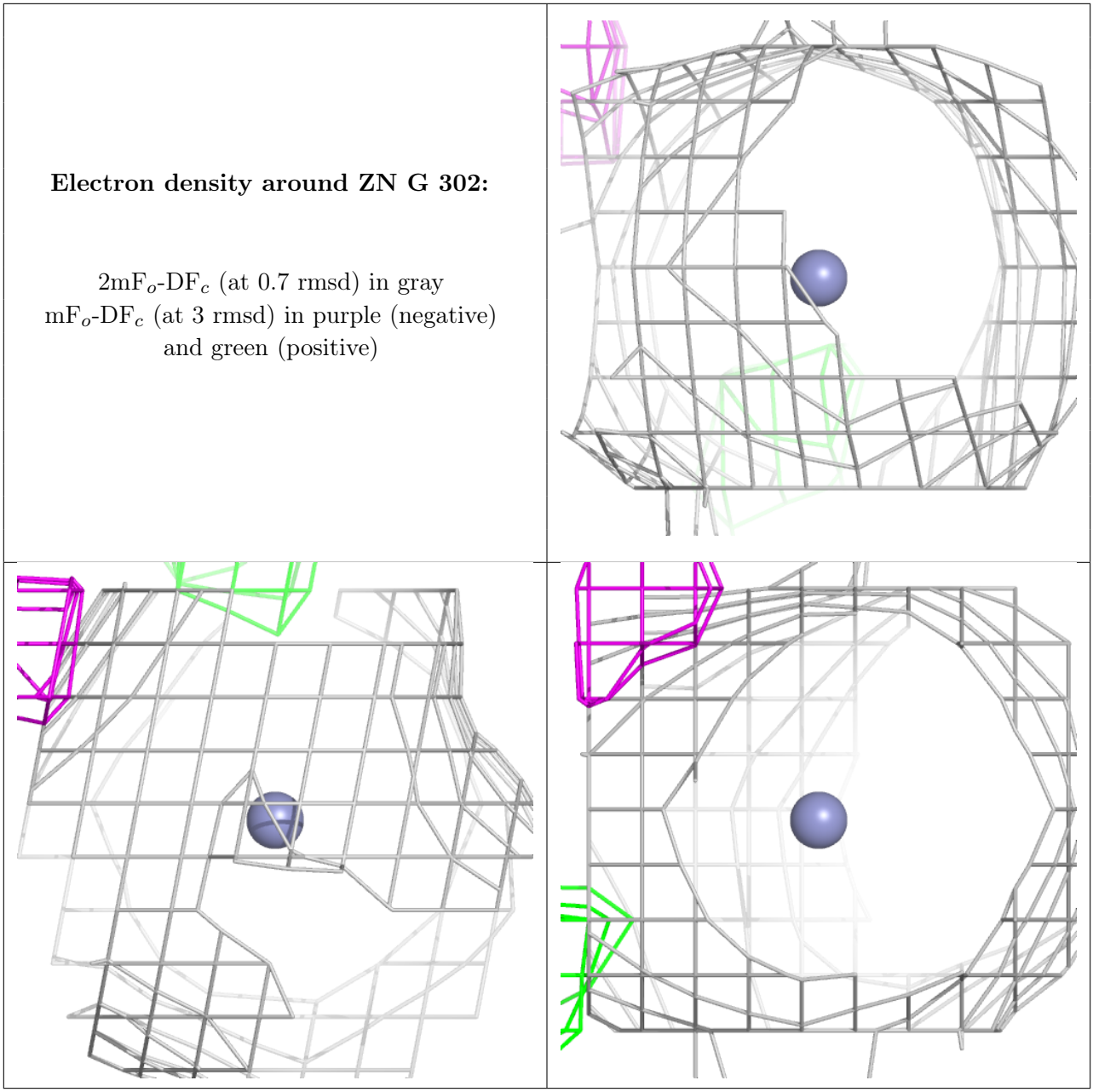
Electron density around ZN P 302:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)



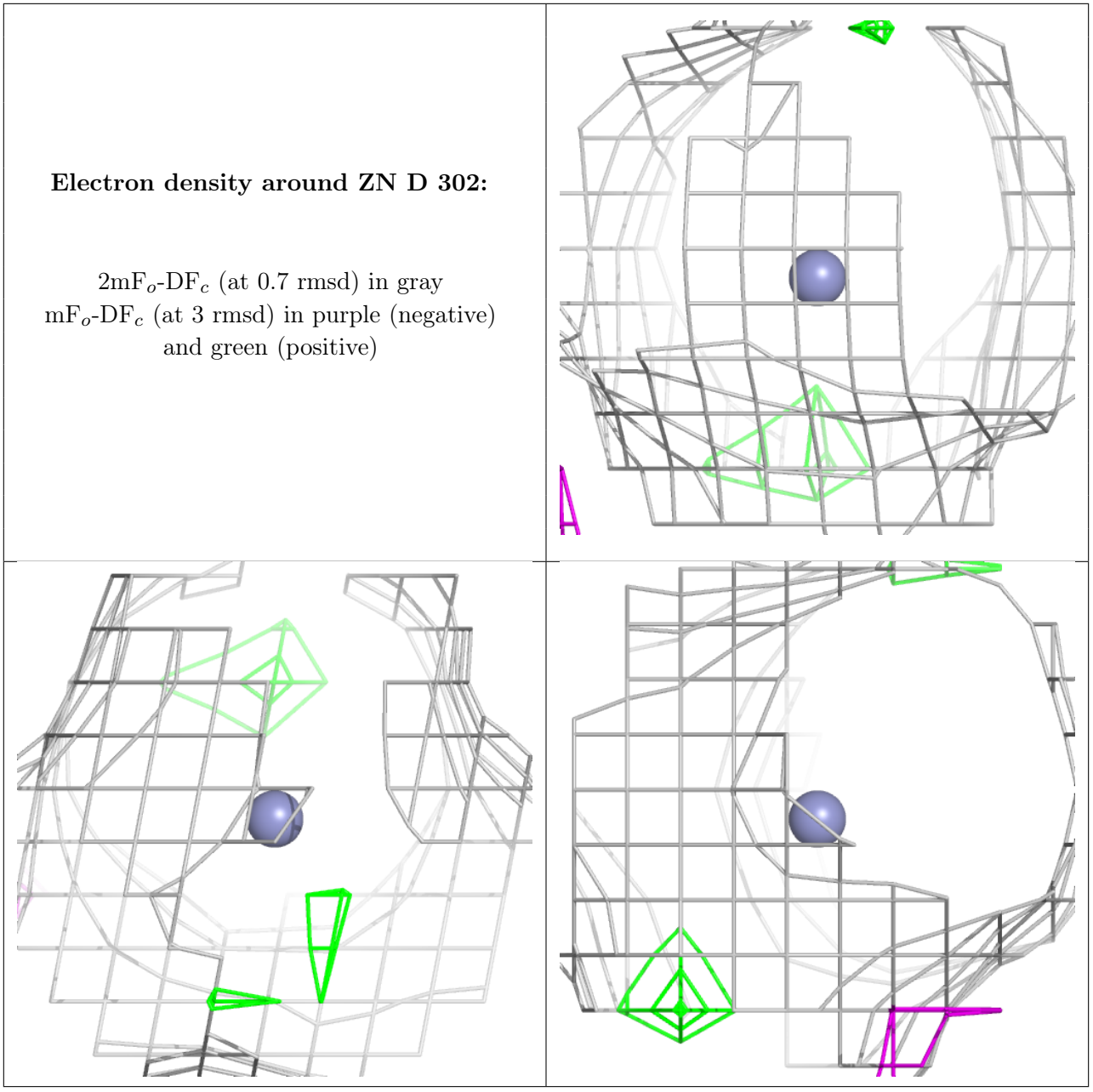
Electron density around ZN G 302:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)



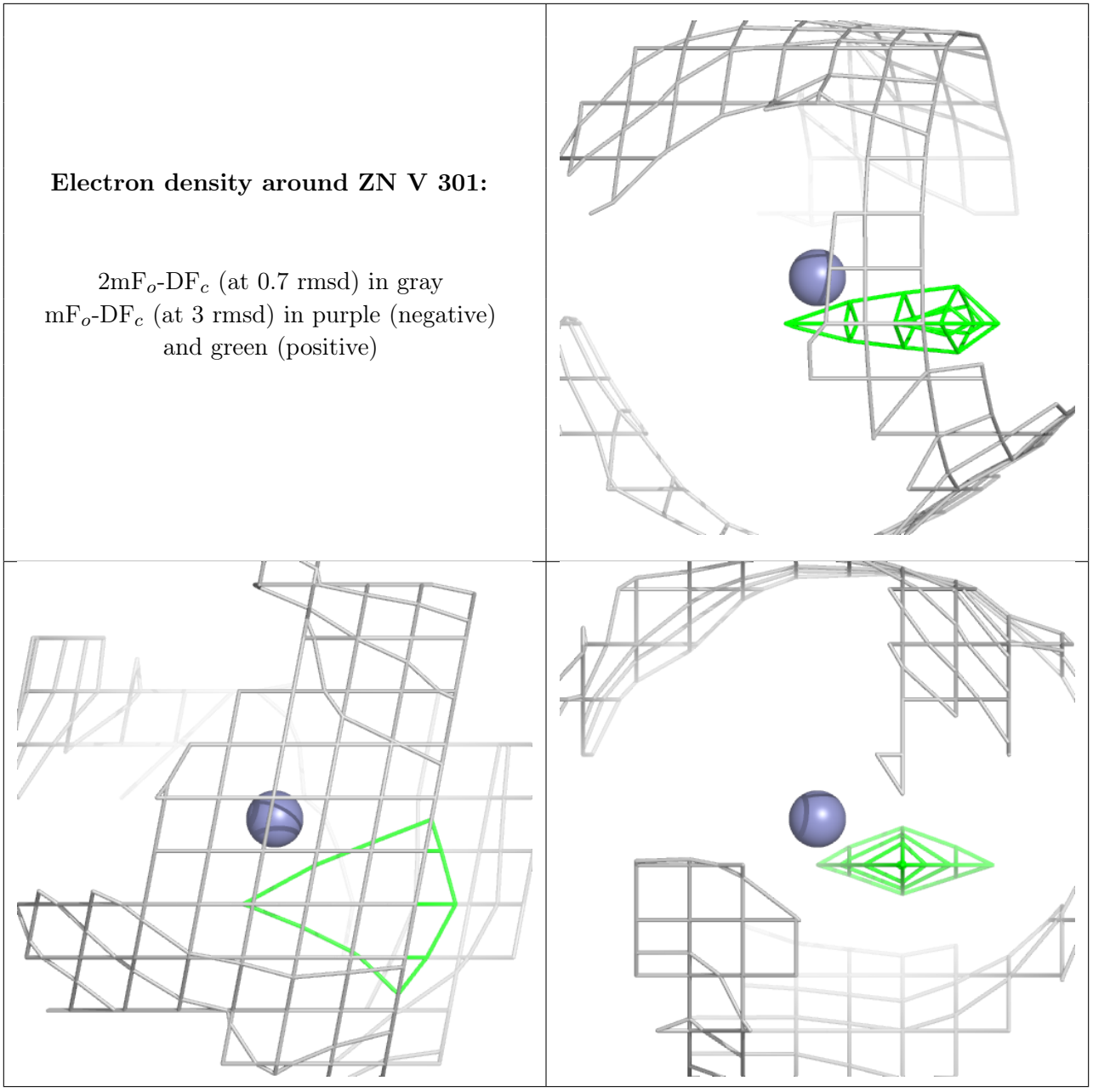
Electron density around ZN D 302:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)



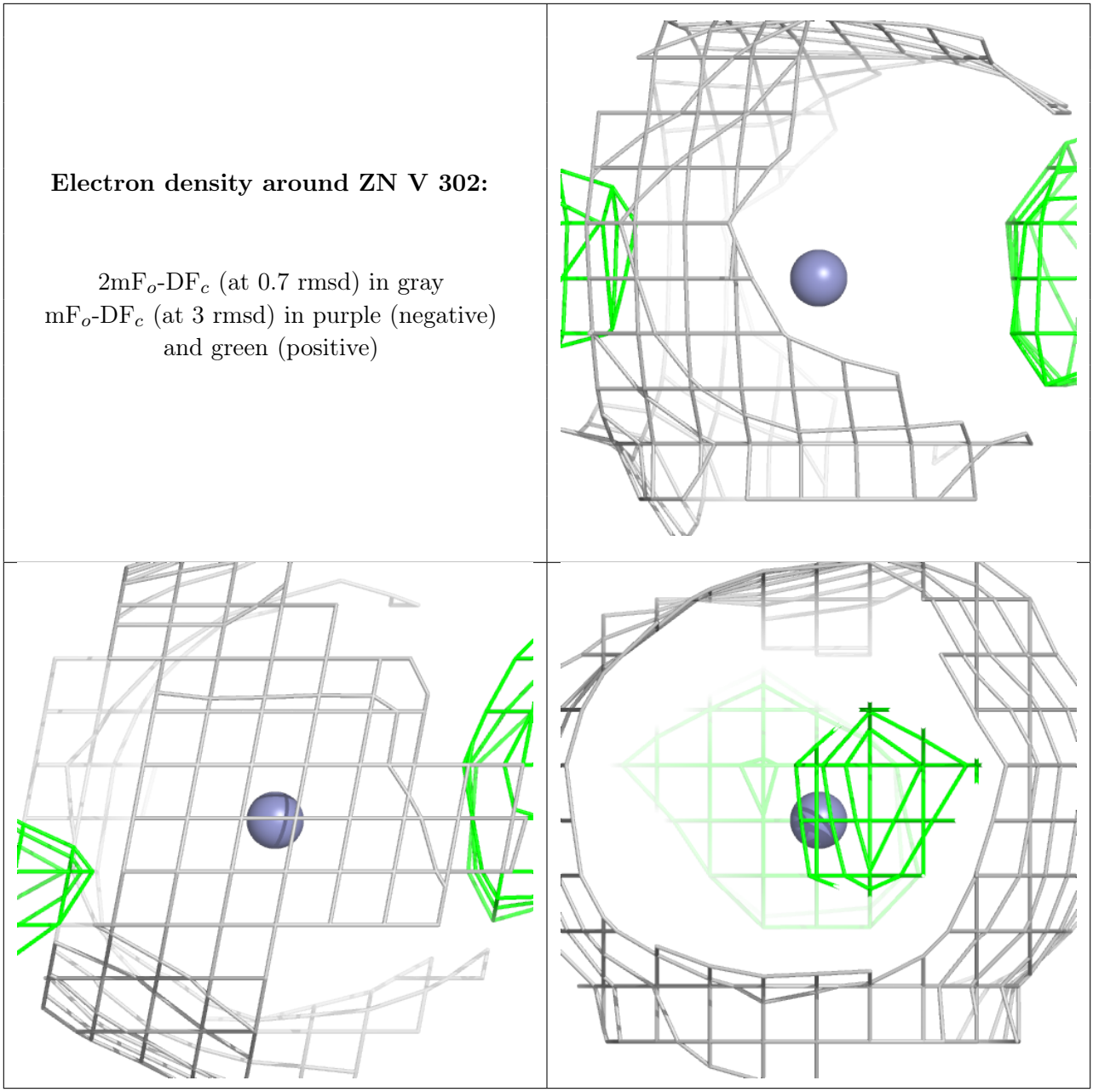
Electron density around ZN V 301:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)



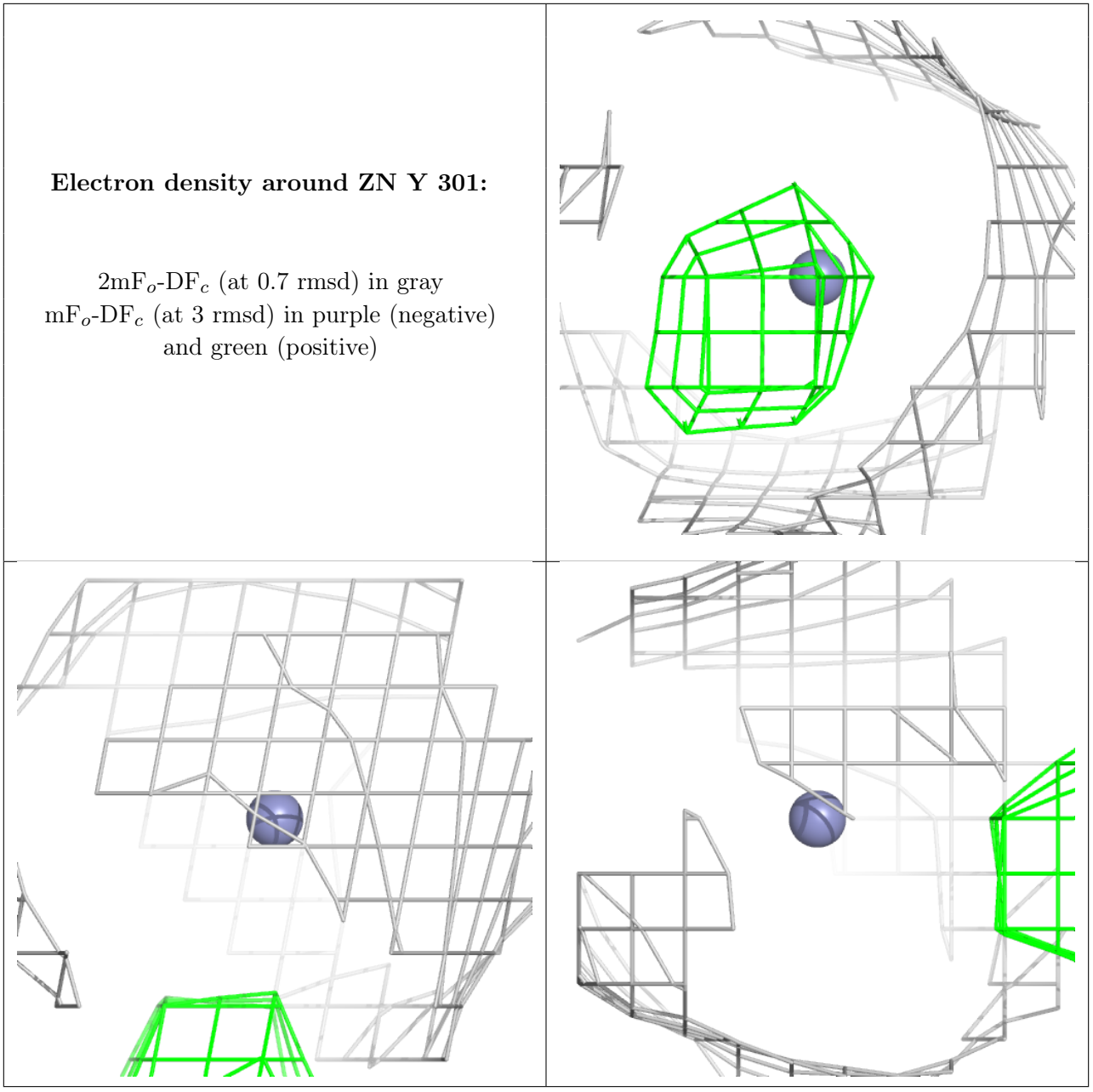
Electron density around ZN V 302:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)



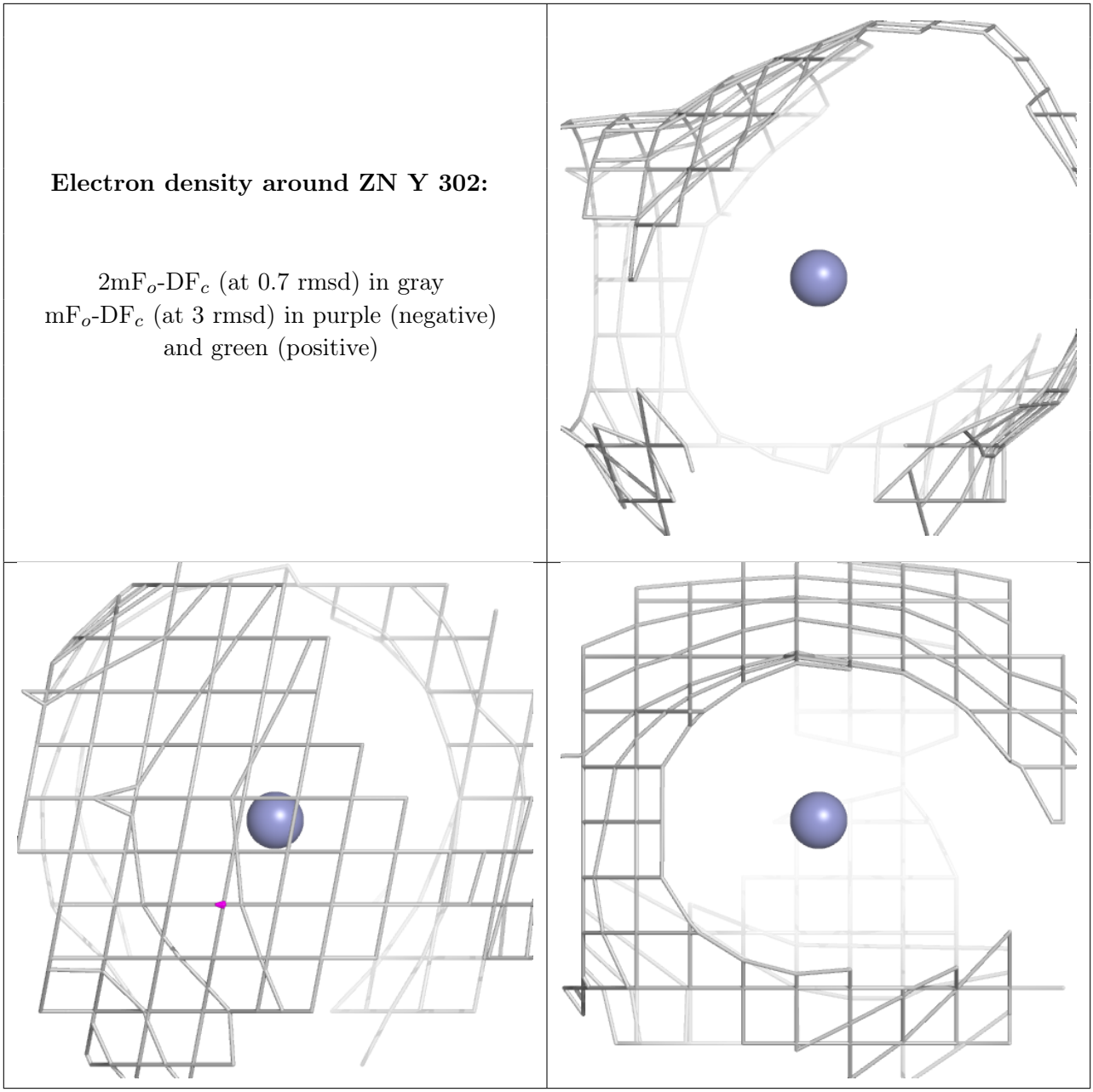
Electron density around ZN Y 301:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)



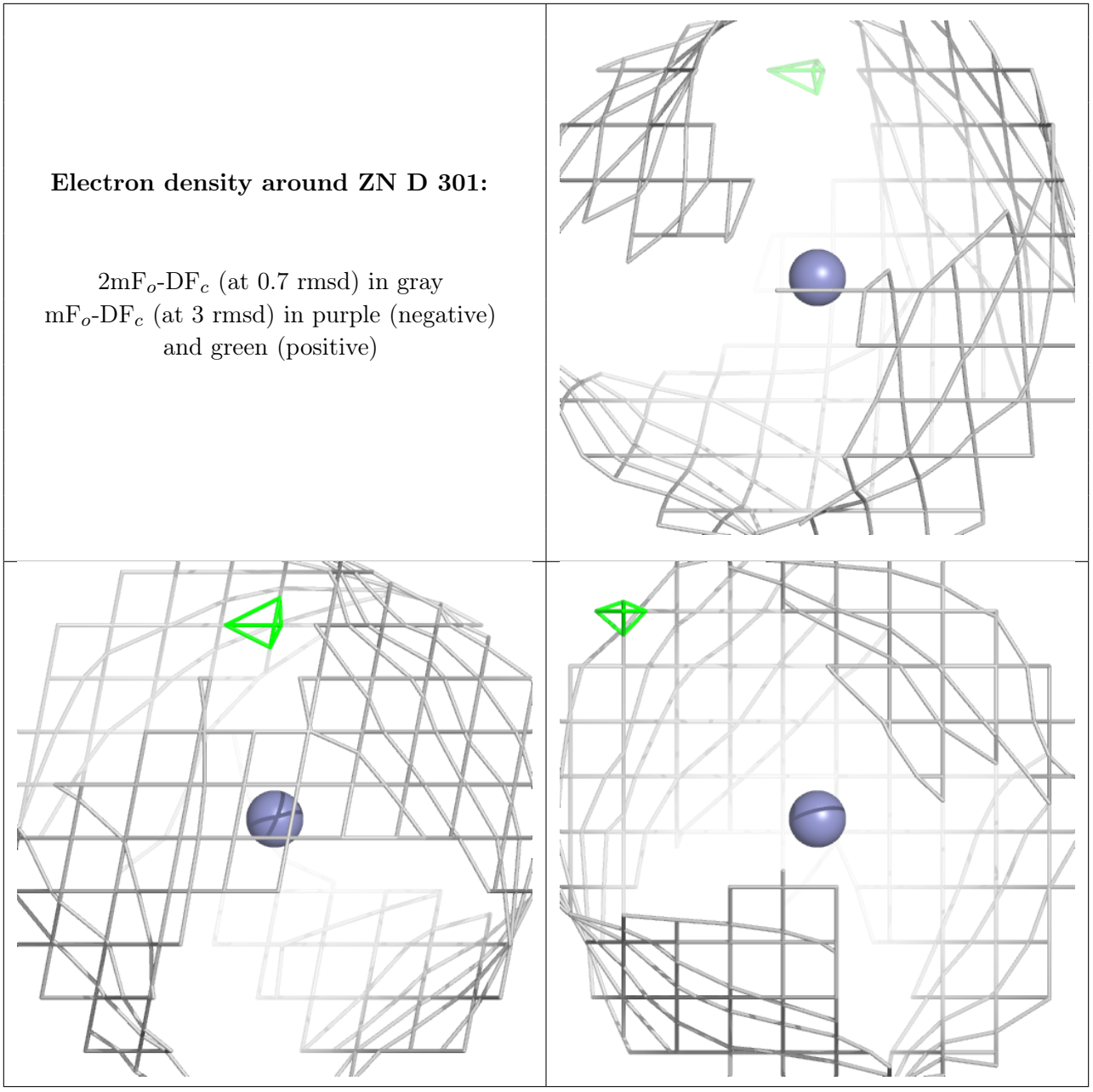
Electron density around ZN Y 302:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)



Electron density around ZN D 301:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)



6.5 Other polymers

There are no such residues in this entry.