Full wwPDB X-ray Structure Validation Report ⓘ

Mar 18, 2024 – 04:54 PM JST

PDB ID	:	6KSY
Title	:	Crystal structure of arginase from Zymomonas mobilis ZM4
Authors	:	Park, S.Y.; Hwangbo, S.A.
Deposited on	:	2019-08-26
Resolution	:	1.65 Å(reported)

This is a Full wwPDB X-ray Structure Validation 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 ⓘ 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 ⓘ) were used in the production of this report:

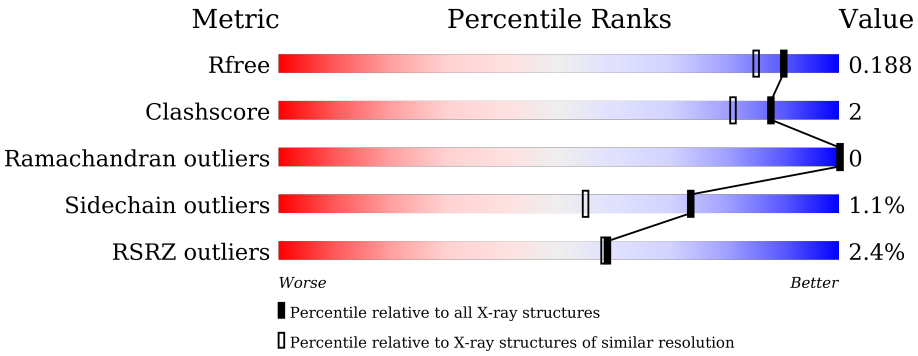
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.36
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson 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 1.65 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	3122 (1.66-1.62)
Clashscore	141614	3268 (1.66-1.62)
Ramachandran outliers	138981	3215 (1.66-1.62)
Sidechain outliers	138945	3215 (1.66-1.62)
RSRZ outliers	127900	3079 (1.66-1.62)

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	290	3% 93% 5%
1	B	290	3% 94% 5%
1	C	290	2% 96%
1	D	290	2% 93% 5%

2 Entry composition (i)

There are 3 unique types of molecules in this entry. The entry contains 9774 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 Arginase/agmatinase/formiminoglutamase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	286	Total	C	N	O	S	0	0	0
			2239	1438	380	416	5			
1	B	289	Total	C	N	O	S	0	0	0
			2259	1450	383	421	5			
1	C	284	Total	C	N	O	S	0	0	0
			2222	1428	376	413	5			
1	D	286	Total	C	N	O	S	0	0	0
			2238	1438	379	416	5			

- Molecule 2 is ZINC ION (three-letter code: ZN) (formula: Zn) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	2	Total	Zn	0	0
			2	2		
2	B	2	Total	Zn	0	0
			2	2		
2	C	2	Total	Zn	0	0
			2	2		
2	D	2	Total	Zn	0	0
			2	2		

- Molecule 3 is water.

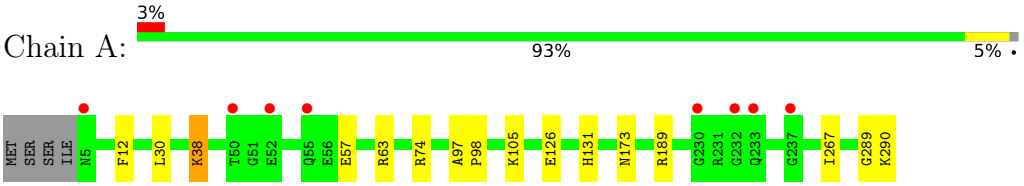
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	203	Total	O	0	0
			203	203		
3	B	204	Total	O	0	0
			204	204		
3	C	209	Total	O	0	0
			209	209		
3	D	192	Total	O	0	0
			192	192		

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.

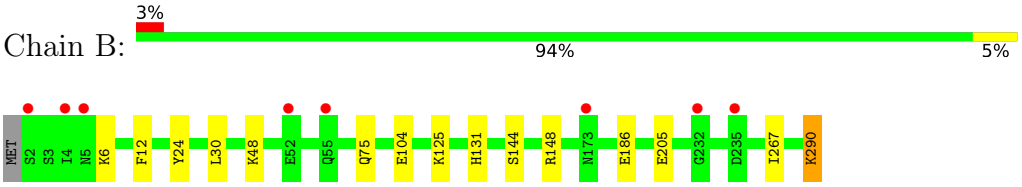
• Molecule 1: Arginase/agmatinase/formiminoglutamase

Chain A:



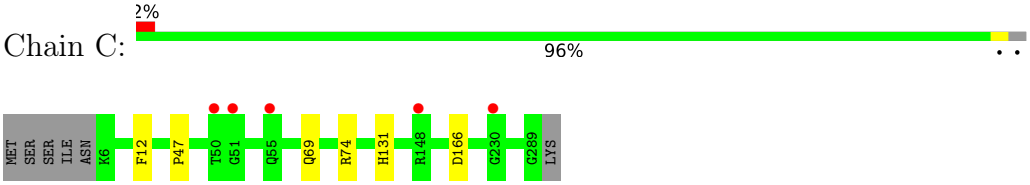
• Molecule 1: Arginase/agmatinase/formiminoglutamase

Chain B:



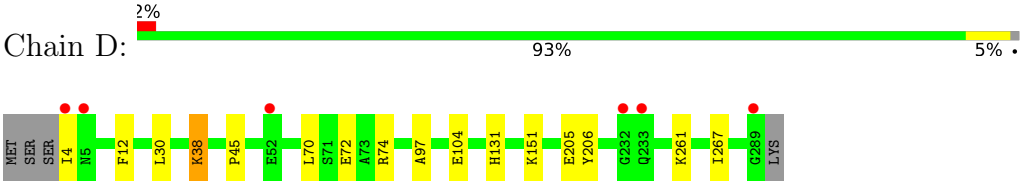
• Molecule 1: Arginase/agmatinase/formiminoglutamase

Chain C:



• Molecule 1: Arginase/agmatinase/formiminoglutamase

Chain D:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α, β, γ	160.11Å 100.32Å 79.60Å 90.00° 113.31° 90.00°	Depositor
Resolution (Å)	20.00 – 1.65 20.00 – 1.65	Depositor EDS
% Data completeness (in resolution range)	94.8 (20.00-1.65) 94.8 (20.00-1.65)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.44 (at 1.65Å)	Xtriage
Refinement program	PHENIX	Depositor
R, R_{free}	0.165 , 0.188 0.165 , 0.188	Depositor DCC
R_{free} test set	1999 reflections (1.52%)	wwPDB-VP
Wilson B-factor (Å^2)	20.1	Xtriage
Anisotropy	0.072	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.42 , 64.7	EDS
L-test for twinning[2]	$< \|L\| > = 0.47$, $< L^2 > = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.97	EDS
Total number of atoms	9774	wwPDB-VP
Average B, all atoms (Å^2)	23.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.60% 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

5.1 Standard geometry

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	0.36	0/2299	0.56	1/3133 (0.0%)
1	B	0.36	0/2319	0.54	0/3160
1	C	0.34	0/2282	0.53	0/3111
1	D	0.35	0/2298	0.56	1/3133 (0.0%)
All	All	0.35	0/9198	0.55	2/12537 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	D	38	LYS	CD-CE-NZ	-7.97	93.36	111.70
1	A	38	LYS	CA-CB-CG	-5.00	102.39	113.40

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

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	2239	0	2239	13	0
1	B	2259	0	2260	11	0
1	C	2222	0	2220	2	0
1	D	2238	0	2237	14	0
2	A	2	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	2	0	0	0	0
2	C	2	0	0	0	0
2	D	2	0	0	0	0
3	A	203	0	0	5	2
3	B	204	0	0	6	1
3	C	209	0	0	1	4
3	D	192	0	0	2	3
All	All	9774	0	8956	36	5

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 2.

All (36) 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:D:74:ARG:NH1	1:D:104:GLU:OE2	2.22	0.72
1:B:186:GLU:OE1	3:B:401:HOH:O	2.08	0.71
1:B:290:LYS:O	3:B:402:HOH:O	2.09	0.71
1:A:57:GLU:OE1	3:A:402:HOH:O	2.09	0.70
1:C:74:ARG:NH1	3:C:401:HOH:O	2.26	0.69
1:A:126:GLU:O	1:B:6:LYS:HE2	1.94	0.68
1:A:189:ARG:NH2	3:A:404:HOH:O	2.29	0.65
1:D:45:PRO:HD2	1:D:72:GLU:HG2	1.85	0.59
1:B:104:GLU:OE2	3:B:404:HOH:O	2.17	0.57
1:A:173:ASN:ND2	3:A:401:HOH:O	2.02	0.56
1:D:205:GLU:HG2	1:D:206:TYR:CE2	2.41	0.55
1:B:205:GLU:OE1	3:B:405:HOH:O	2.18	0.54
1:A:290:LYS:N	1:D:38:LYS:HD2	2.24	0.53
1:D:74:ARG:O	1:D:74:ARG:HD2	2.11	0.51
1:A:30:LEU:HD21	1:A:267:ILE:HB	1.95	0.48
1:A:289:GLY:H	1:D:38:LYS:NZ	2.10	0.48
1:B:290:LYS:CA	3:B:402:HOH:O	2.61	0.48
1:A:105:LYS:NZ	3:A:412:HOH:O	2.46	0.48
1:D:205:GLU:HG2	1:D:206:TYR:CD2	2.49	0.47
1:B:24:TYR:CZ	1:B:48:LYS:HG2	2.49	0.47
1:D:30:LEU:HD21	1:D:267:ILE:HB	1.97	0.46
1:C:47:PRO:HD2	1:C:69:GLN:OE1	2.15	0.46
1:B:290:LYS:HA	3:B:402:HOH:O	2.15	0.46
1:B:144:SER:O	1:B:148:ARG:HG2	2.17	0.45
1:A:38:LYS:H	1:A:38:LYS:HG2	1.57	0.44
1:D:104:GLU:CD	1:D:151:LYS:HD2	2.38	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:38:LYS:HD3	1:D:38:LYS:N	2.32	0.44
1:D:4:ILE:N	3:D:409:HOH:O	2.50	0.43
1:A:97:ALA:HB3	1:A:98:PRO:HD3	2.01	0.43
1:A:74:ARG:HD3	3:A:421:HOH:O	2.18	0.42
1:B:75:GLN:NE2	1:B:75:GLN:HA	2.34	0.42
1:D:261:LYS:NZ	3:D:405:HOH:O	2.44	0.42
1:B:30:LEU:HD21	1:B:267:ILE:HB	2.02	0.42
1:A:57:GLU:HB2	1:A:63:ARG:HB3	2.01	0.42
1:A:289:GLY:H	1:D:38:LYS:HZ3	1.67	0.41
1:D:70:LEU:HD12	1:D:97:ALA:HB2	2.02	0.41

All (5) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:B:556:HOH:O	3:D:526:HOH:O[4_545]	2.01	0.19
3:C:585:HOH:O	3:D:561:HOH:O[3_556]	2.01	0.19
3:C:490:HOH:O	3:D:552:HOH:O[4_555]	2.03	0.17
3:A:595:HOH:O	3:C:596:HOH:O[2_656]	2.04	0.16
3:A:434:HOH:O	3:C:568:HOH:O[2_656]	2.19	0.01

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	284/290 (98%)	278 (98%)	6 (2%)	0	100	100
1	B	287/290 (99%)	281 (98%)	6 (2%)	0	100	100
1	C	282/290 (97%)	278 (99%)	4 (1%)	0	100	100
1	D	284/290 (98%)	278 (98%)	6 (2%)	0	100	100
All	All	1137/1160 (98%)	1115 (98%)	22 (2%)	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	247/251 (98%)	245 (99%)	2 (1%)	81	68
1	B	250/251 (100%)	246 (98%)	4 (2%)	62	39
1	C	245/251 (98%)	242 (99%)	3 (1%)	71	51
1	D	247/251 (98%)	245 (99%)	2 (1%)	81	68
All	All	989/1004 (98%)	978 (99%)	11 (1%)	73	55

All (11) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	12	PHE
1	A	131	HIS
1	B	12	PHE
1	B	125	LYS
1	B	131	HIS
1	B	290	LYS
1	C	12	PHE
1	C	131	HIS
1	C	166	ASP
1	D	12	PHE
1	D	131	HIS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	5	ASN
1	B	75	GLN

5.3.3 RNA

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains

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

5.5 Carbohydrates

There are no monosaccharides in this entry.

5.6 Ligand geometry

Of 8 ligands modelled in this entry, 8 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

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	286/290 (98%)	-0.12	8 (2%)	53	51	13, 20, 37, 48	0
1	B	289/290 (99%)	-0.07	8 (2%)	53	51	15, 21, 37, 46	0
1	C	284/290 (97%)	-0.16	5 (1%)	68	69	14, 20, 32, 51	0
1	D	286/290 (98%)	-0.17	6 (2%)	63	64	15, 21, 35, 49	0
All	All	1145/1160 (98%)	-0.13	27 (2%)	59	58	13, 21, 36, 51	0

All (27) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	4	ILE	5.8
1	B	2	SER	4.7
1	B	5	ASN	4.1
1	C	50	THR	3.8
1	D	232	GLY	3.5
1	A	230	GLY	3.4
1	D	5	ASN	3.3
1	D	233	GLN	3.1
1	A	237	GLY	2.9
1	A	52	GLU	2.9
1	A	55	GLN	2.7
1	B	4	ILE	2.6
1	C	55	GLN	2.5
1	D	52	GLU	2.5
1	A	5	ASN	2.5
1	C	230	GLY	2.4
1	C	51	GLY	2.3
1	A	232	GLY	2.3
1	A	50	THR	2.2
1	B	173	ASN	2.2
1	C	148	ARG	2.1

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Mol	Chain	Res	Type	RSRZ
1	B	55	GLN	2.1
1	B	232	GLY	2.1
1	B	52	GLU	2.1
1	B	235	ASP	2.1
1	D	289	GLY	2.1
1	A	233	GLN	2.0

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)

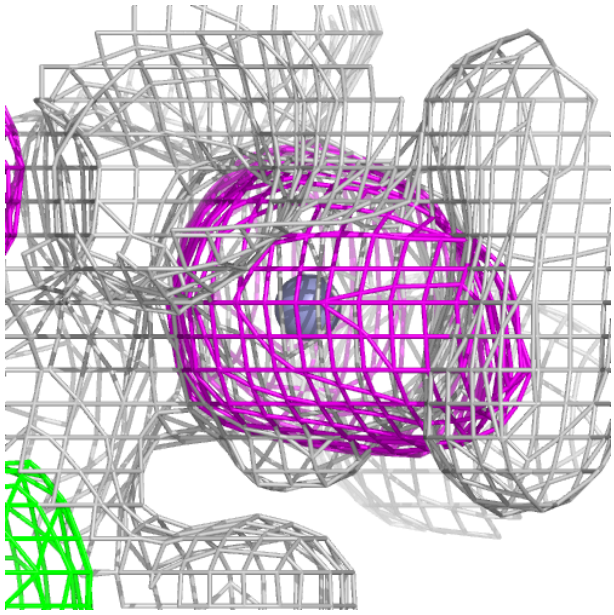
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95^{th} percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9'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

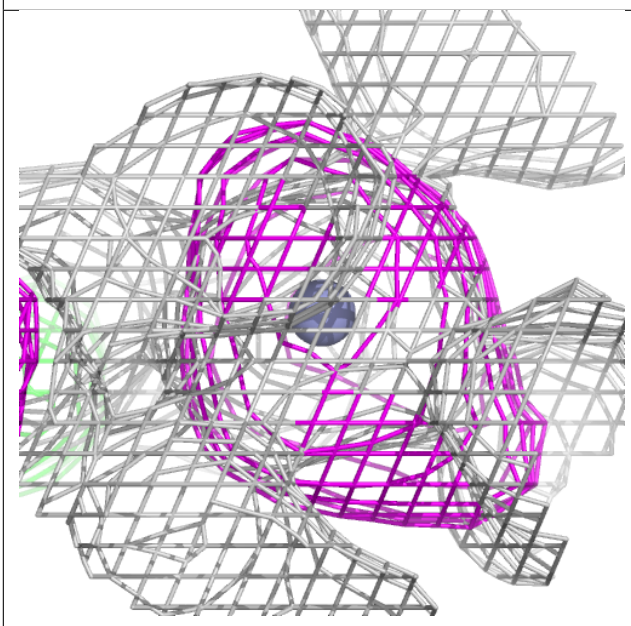
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
2	ZN	A	302	1/1	0.99	0.12	29,29,29,29	0
2	ZN	D	301	1/1	0.99	0.05	23,23,23,23	0
2	ZN	B	301	1/1	1.00	0.13	31,31,31,31	0
2	ZN	B	302	1/1	1.00	0.07	25,25,25,25	0
2	ZN	C	301	1/1	1.00	0.05	21,21,21,21	0
2	ZN	C	302	1/1	1.00	0.10	27,27,27,27	0
2	ZN	A	301	1/1	1.00	0.03	21,21,21,21	0
2	ZN	D	302	1/1	1.00	0.12	29,29,29,29	0

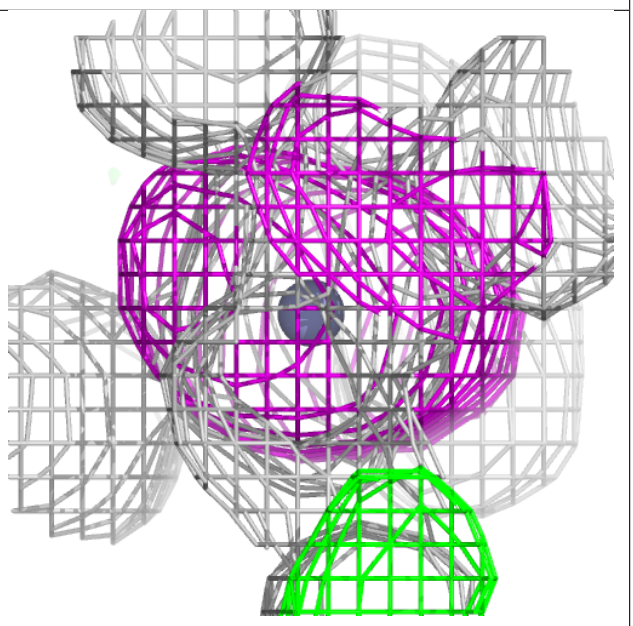
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around ZN 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)

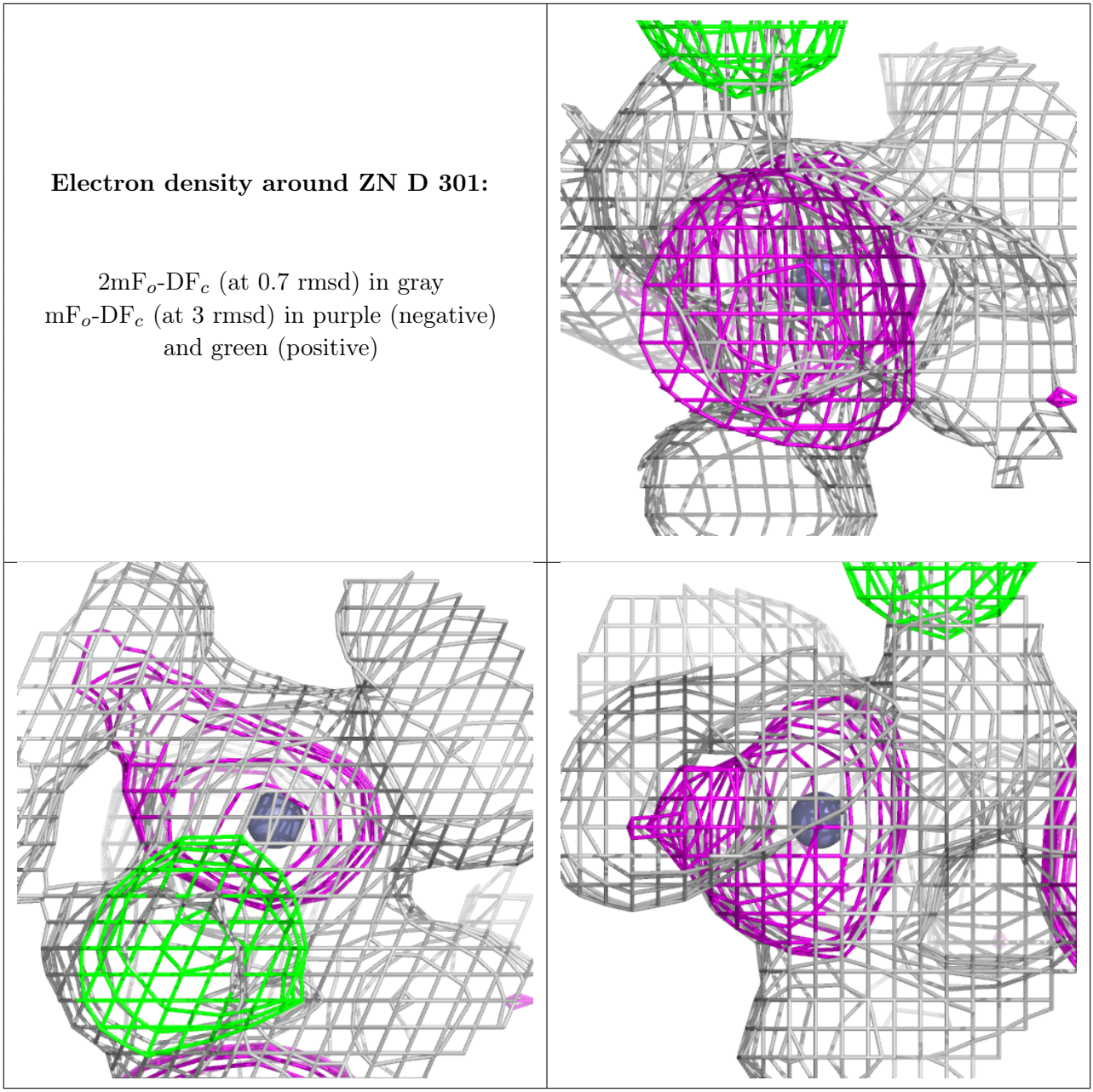






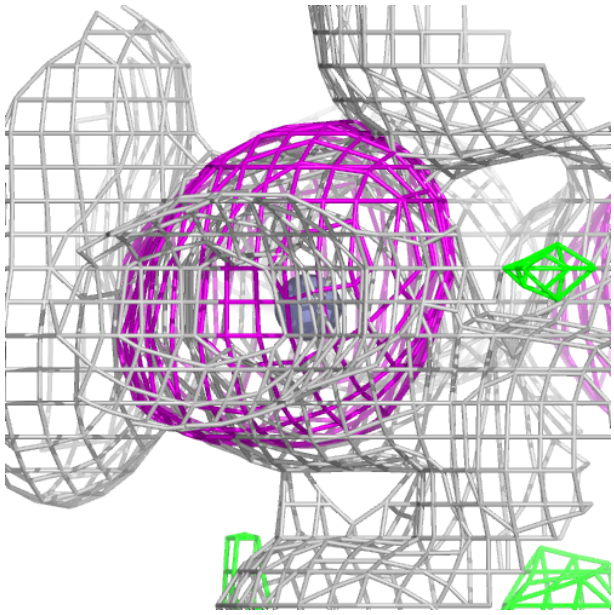
Electron density around ZN D 301:

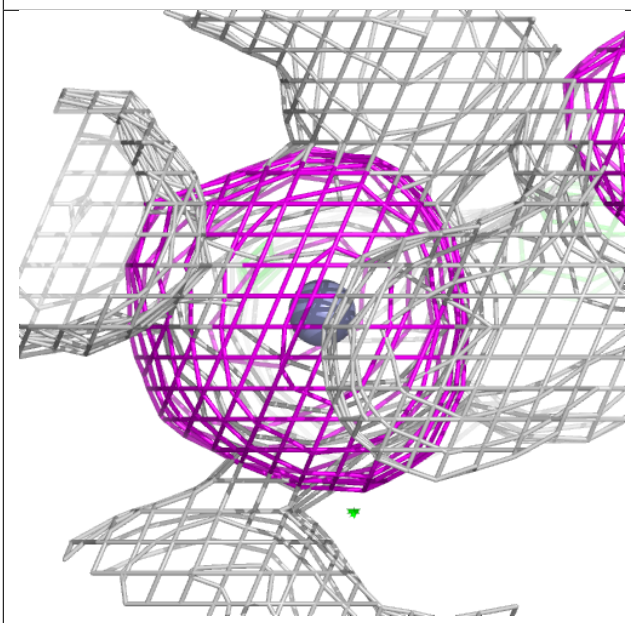
$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

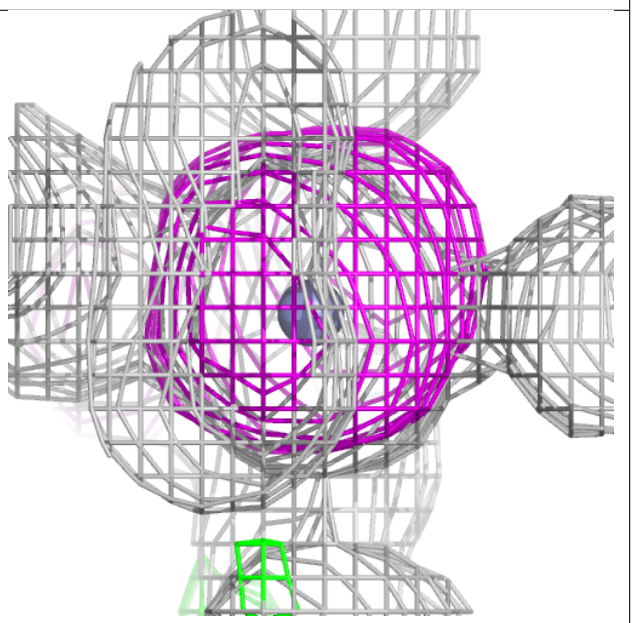


Electron density around ZN B 301:

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

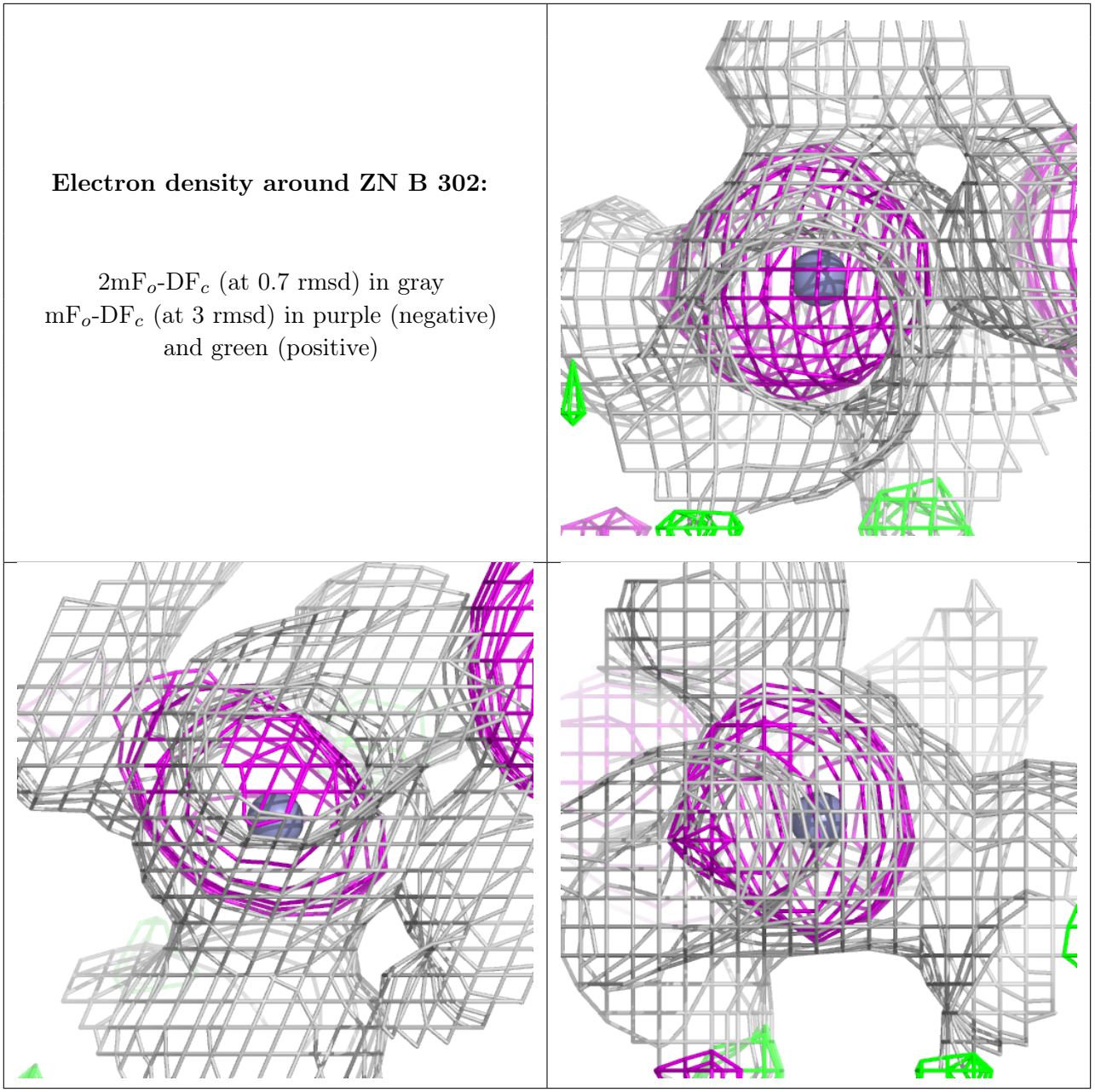






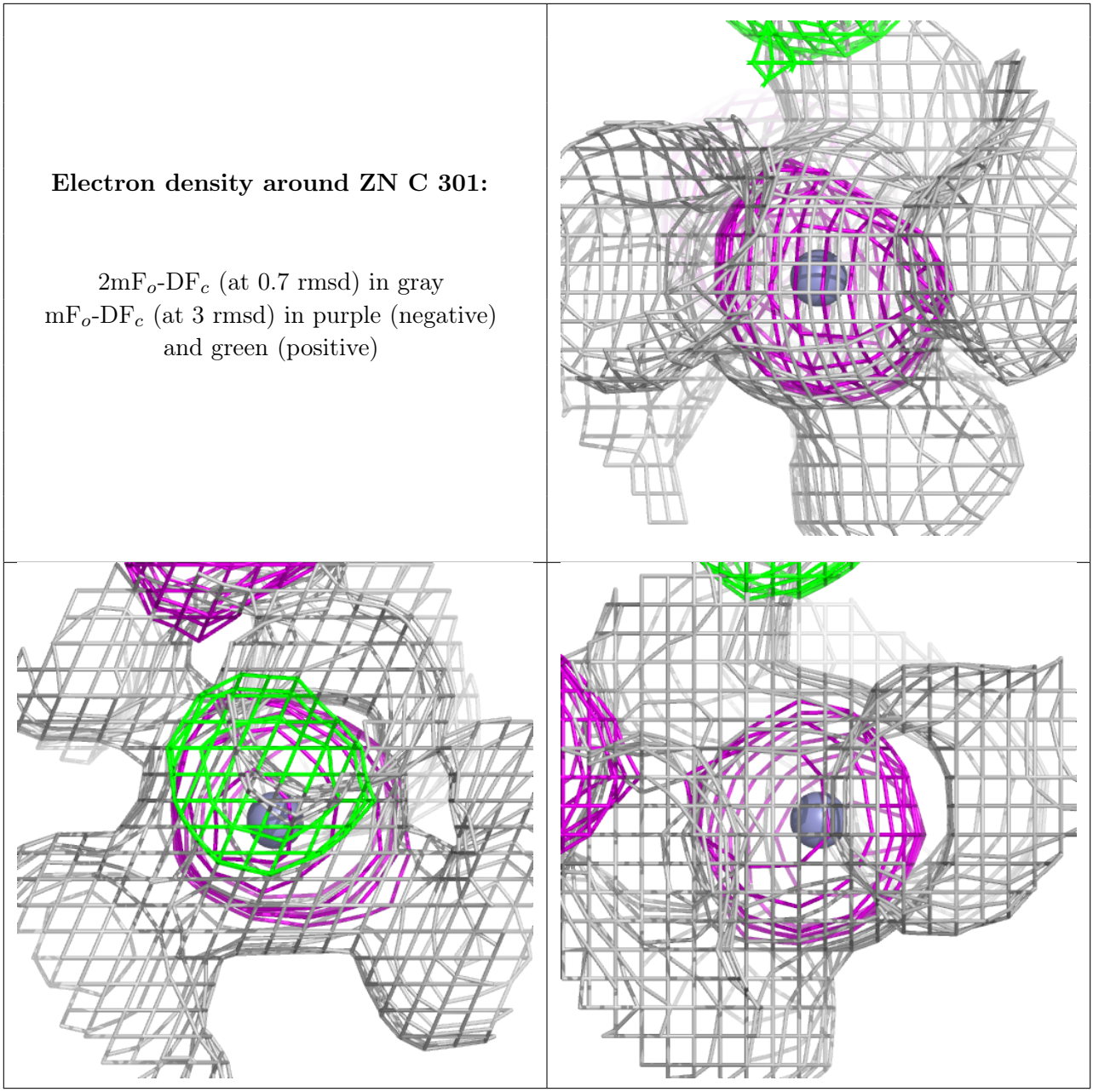
Electron density around ZN B 302:

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



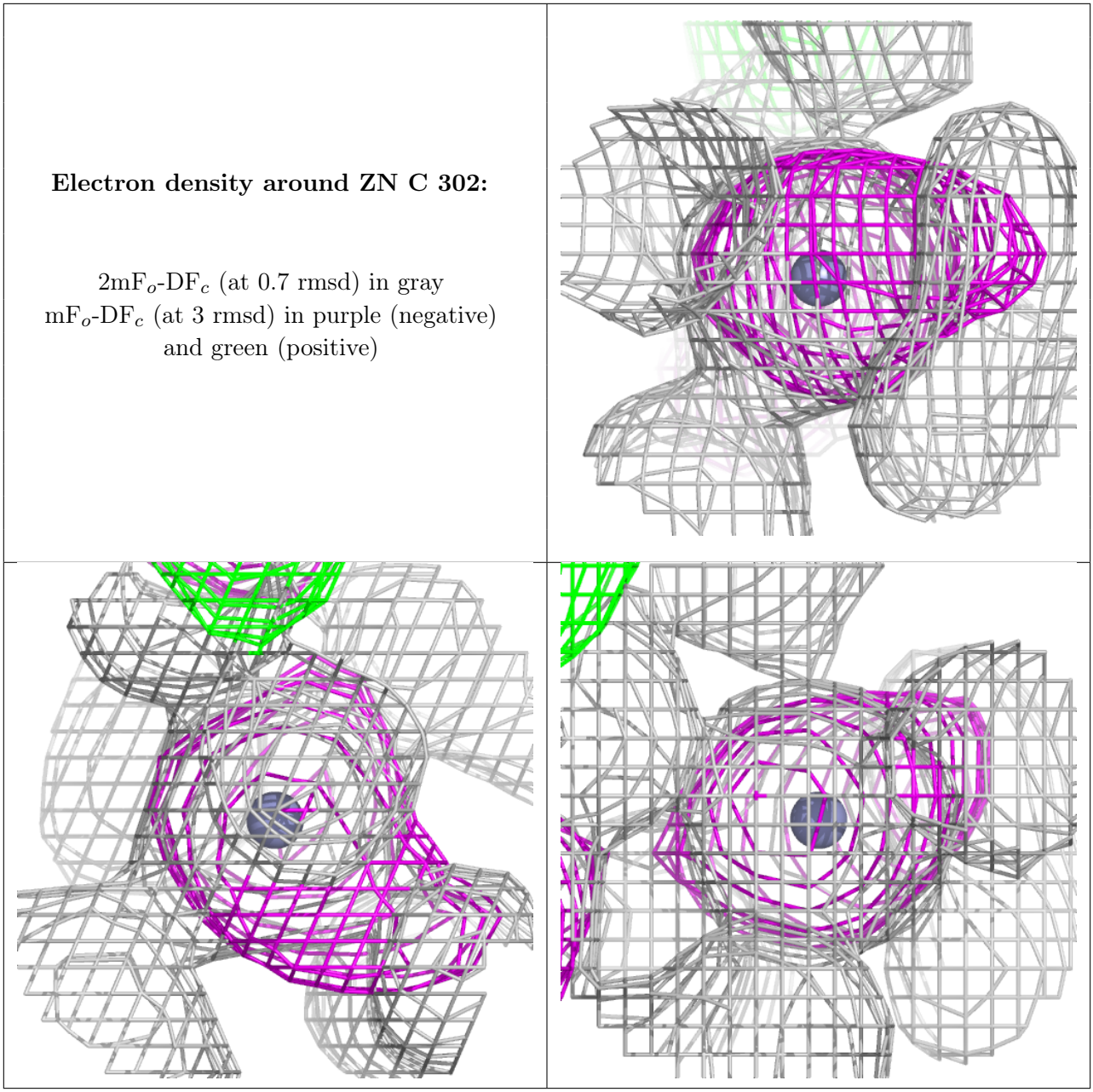
Electron density around ZN C 301:

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



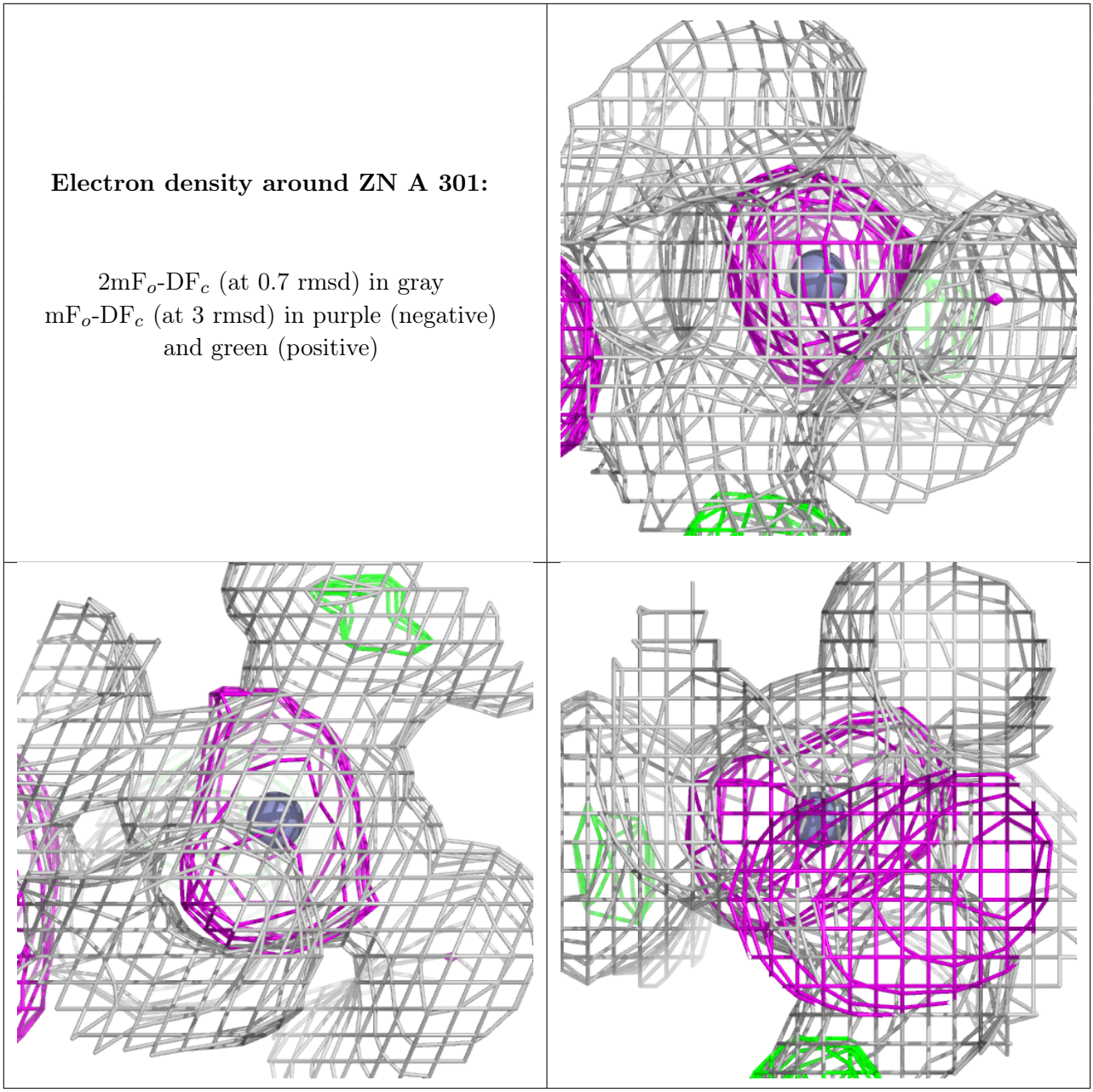
Electron density around ZN C 302:

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



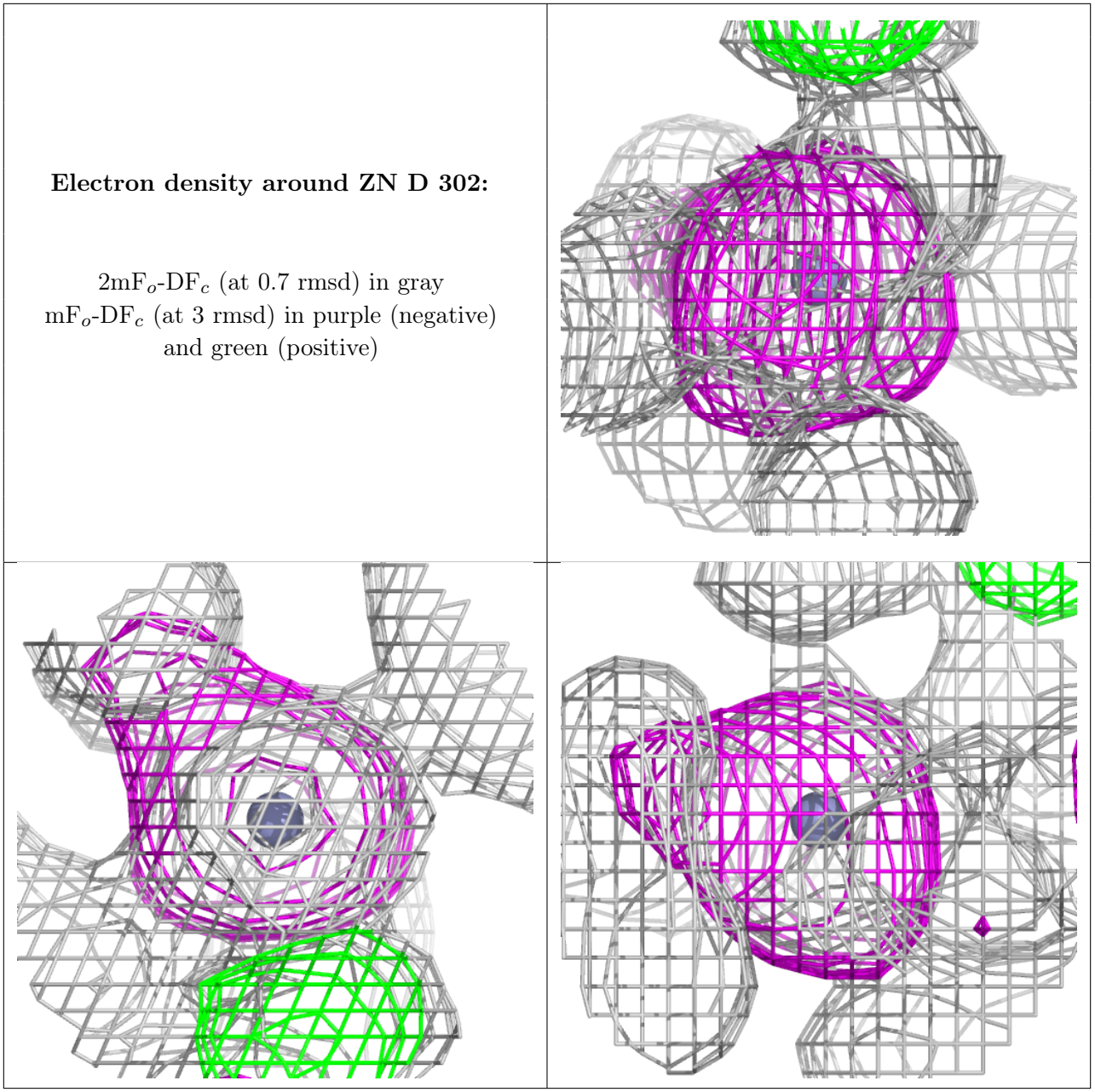
Electron density around ZN A 301:

$2mF_o$-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 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)



6.5 Other polymers (i)

There are no such residues in this entry.