wwPDB X-ray Structure Validation Summary Report (i)

Oct 14, 2023 – 06:42 PM EDT

PDB ID	:	8CX9
Title	:	Structure of the SARS-COV2 PLpro (C111S) in complex with a dimeric Ubv that inhibits activity by an unusual allosteric mechanism
Authors	:	Singer, A.U.; Slater, C.L.; Patel, A.; Russel, R.; Mark, B.L.; Sidhu, S.S.
Deposited on	:	2022-05-20
Resolution	:	3.50 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:

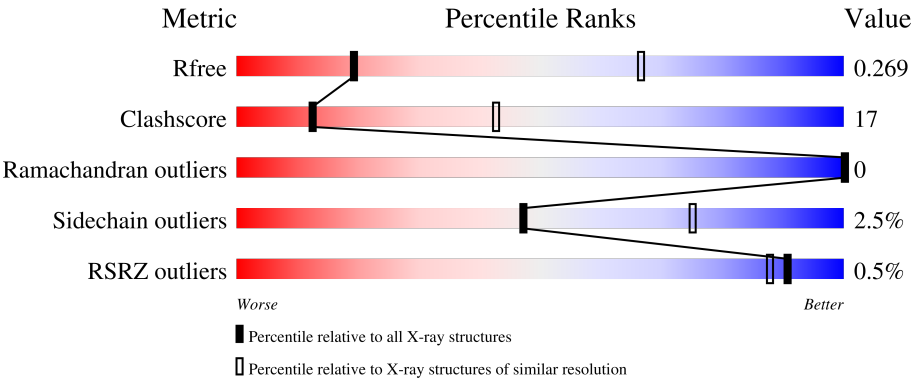
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.36
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson 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 3.50 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	1659 (3.60-3.40)
Clashscore	141614	1036 (3.58-3.42)
Ramachandran outliers	138981	1005 (3.58-3.42)
Sidechain outliers	138945	1006 (3.58-3.42)
RSRZ outliers	127900	1559 (3.60-3.40)

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	320	65% 32%
1	B	320	73% 24%
1	C	320	69% 28%
1	D	320	% 67% 28%
2	E	103	36% 35% 28%

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Mol	Chain	Length	Quality of chain
2	F	103	41% 32% 27%
2	G	103	42% 30% 27%
2	H	103	45% 27% 28%
2	I	103	47% 26% 27%
2	J	103	50% 20% 28%
2	K	103	2% 44% 27% 29%
2	L	103	5% 38% 32% 28%

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
4	BR	A	402	-	-	X	-

2 Entry composition (i)

There are 6 unique types of molecules in this entry. The entry contains 13989 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 Papain-like protease nsp3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	B	314	Total	C	N	O	S	0	0	0
			2430	1542	399	470	19			
1	A	314	Total	C	N	O	S	0	0	0
			2450	1555	402	474	19			
1	C	312	Total	C	N	O	S	0	0	0
			2387	1514	392	462	19			
1	D	310	Total	C	N	O	S	0	0	0
			2365	1503	385	458	19			

There are 20 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	-4	GLY	-	expression tag	UNP P0DTC1
B	-3	PRO	-	expression tag	UNP P0DTC1
B	-2	LEU	-	expression tag	UNP P0DTC1
B	-1	GLY	-	expression tag	UNP P0DTC1
B	0	SER	-	expression tag	UNP P0DTC1
A	-4	GLY	-	expression tag	UNP P0DTC1
A	-3	PRO	-	expression tag	UNP P0DTC1
A	-2	LEU	-	expression tag	UNP P0DTC1
A	-1	GLY	-	expression tag	UNP P0DTC1
A	0	SER	-	expression tag	UNP P0DTC1
C	-4	GLY	-	expression tag	UNP P0DTC1
C	-3	PRO	-	expression tag	UNP P0DTC1
C	-2	LEU	-	expression tag	UNP P0DTC1
C	-1	GLY	-	expression tag	UNP P0DTC1
C	0	SER	-	expression tag	UNP P0DTC1
D	-4	GLY	-	expression tag	UNP P0DTC1
D	-3	PRO	-	expression tag	UNP P0DTC1
D	-2	LEU	-	expression tag	UNP P0DTC1
D	-1	GLY	-	expression tag	UNP P0DTC1
D	0	SER	-	expression tag	UNP P0DTC1

- Molecule 2 is a protein called Ubiquitin variant UbV.CV2.1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	E	74	Total	C	N	O	S	0	0	0
			571	358	97	113	3			
2	I	75	Total	C	N	O	S	0	0	0
			551	349	95	105	2			
2	F	75	Total	C	N	O	S	0	0	0
			580	365	100	112	3			
2	G	75	Total	C	N	O	S	0	0	0
			545	343	93	106	3			
2	K	73	Total	C	N	O	S	0	0	0
			505	316	86	101	2			
2	L	74	Total	C	N	O	S	0	0	0
			517	322	90	102	3			
2	H	74	Total	C	N	O	S	0	0	0
			545	344	94	104	3			
2	J	74	Total	C	N	O	S	0	0	0
			530	333	91	103	3			

There are 320 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
E	-24	HIS	-	expression tag	UNP P62987
E	-23	HIS	-	expression tag	UNP P62987
E	-22	HIS	-	expression tag	UNP P62987
E	-21	HIS	-	expression tag	UNP P62987
E	-20	HIS	-	expression tag	UNP P62987
E	-19	HIS	-	expression tag	UNP P62987
E	-18	VAL	-	expression tag	UNP P62987
E	-17	THR	-	expression tag	UNP P62987
E	-16	SER	-	expression tag	UNP P62987
E	-15	LEU	-	expression tag	UNP P62987
E	-14	TYR	-	expression tag	UNP P62987
E	-13	LYS	-	expression tag	UNP P62987
E	-12	LYS	-	expression tag	UNP P62987
E	-11	ALA	-	expression tag	UNP P62987
E	-10	GLY	-	expression tag	UNP P62987
E	-9	SER	-	expression tag	UNP P62987
E	-8	THR	-	expression tag	UNP P62987
E	-7	ASP	-	expression tag	UNP P62987
E	-6	TYR	-	expression tag	UNP P62987
E	-5	LYS	-	expression tag	UNP P62987
E	-4	ASP	-	expression tag	UNP P62987
E	-3	ASP	-	expression tag	UNP P62987

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Chain	Residue	Modelled	Actual	Comment	Reference
E	-2	ASP	-	expression tag	UNP P62987
E	-1	ASP	-	expression tag	UNP P62987
E	0	LYS	-	expression tag	UNP P62987
E	4	SER	PHE	variant	UNP P62987
E	9	MET	THR	variant	UNP P62987
E	10	ARG	GLY	variant	UNP P62987
E	12	ARG	THR	variant	UNP P62987
E	46	SER	ALA	variant	UNP P62987
E	48	MET	LYS	variant	UNP P62987
E	49	LEU	GLN	variant	UNP P62987
E	64	GLY	GLU	variant	UNP P62987
E	68	THR	HIS	variant	UNP P62987
E	70	GLY	VAL	variant	UNP P62987
E	72	ILE	ARG	variant	UNP P62987
E	75	ALA	GLY	variant	UNP P62987
E	76	ASN	GLY	variant	UNP P62987
E	77	GLY	-	insertion	UNP P62987
E	78	VAL	ILE	variant	UNP P62987
I	-24	HIS	-	expression tag	UNP P62987
I	-23	HIS	-	expression tag	UNP P62987
I	-22	HIS	-	expression tag	UNP P62987
I	-21	HIS	-	expression tag	UNP P62987
I	-20	HIS	-	expression tag	UNP P62987
I	-19	HIS	-	expression tag	UNP P62987
I	-18	VAL	-	expression tag	UNP P62987
I	-17	THR	-	expression tag	UNP P62987
I	-16	SER	-	expression tag	UNP P62987
I	-15	LEU	-	expression tag	UNP P62987
I	-14	TYR	-	expression tag	UNP P62987
I	-13	LYS	-	expression tag	UNP P62987
I	-12	LYS	-	expression tag	UNP P62987
I	-11	ALA	-	expression tag	UNP P62987
I	-10	GLY	-	expression tag	UNP P62987
I	-9	SER	-	expression tag	UNP P62987
I	-8	THR	-	expression tag	UNP P62987
I	-7	ASP	-	expression tag	UNP P62987
I	-6	TYR	-	expression tag	UNP P62987
I	-5	LYS	-	expression tag	UNP P62987
I	-4	ASP	-	expression tag	UNP P62987
I	-3	ASP	-	expression tag	UNP P62987
I	-2	ASP	-	expression tag	UNP P62987
I	-1	ASP	-	expression tag	UNP P62987

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Chain	Residue	Modelled	Actual	Comment	Reference
I	0	LYS	-	expression tag	UNP P62987
I	4	SER	PHE	variant	UNP P62987
I	9	MET	THR	variant	UNP P62987
I	10	ARG	GLY	variant	UNP P62987
I	12	ARG	THR	variant	UNP P62987
I	46	SER	ALA	variant	UNP P62987
I	48	MET	LYS	variant	UNP P62987
I	49	LEU	GLN	variant	UNP P62987
I	64	GLY	GLU	variant	UNP P62987
I	68	THR	HIS	variant	UNP P62987
I	70	GLY	VAL	variant	UNP P62987
I	72	ILE	ARG	variant	UNP P62987
I	75	ALA	GLY	variant	UNP P62987
I	76	ASN	GLY	variant	UNP P62987
I	77	GLY	-	insertion	UNP P62987
I	78	VAL	ILE	variant	UNP P62987
F	-24	HIS	-	expression tag	UNP P62987
F	-23	HIS	-	expression tag	UNP P62987
F	-22	HIS	-	expression tag	UNP P62987
F	-21	HIS	-	expression tag	UNP P62987
F	-20	HIS	-	expression tag	UNP P62987
F	-19	HIS	-	expression tag	UNP P62987
F	-18	VAL	-	expression tag	UNP P62987
F	-17	THR	-	expression tag	UNP P62987
F	-16	SER	-	expression tag	UNP P62987
F	-15	LEU	-	expression tag	UNP P62987
F	-14	TYR	-	expression tag	UNP P62987
F	-13	LYS	-	expression tag	UNP P62987
F	-12	LYS	-	expression tag	UNP P62987
F	-11	ALA	-	expression tag	UNP P62987
F	-10	GLY	-	expression tag	UNP P62987
F	-9	SER	-	expression tag	UNP P62987
F	-8	THR	-	expression tag	UNP P62987
F	-7	ASP	-	expression tag	UNP P62987
F	-6	TYR	-	expression tag	UNP P62987
F	-5	LYS	-	expression tag	UNP P62987
F	-4	ASP	-	expression tag	UNP P62987
F	-3	ASP	-	expression tag	UNP P62987
F	-2	ASP	-	expression tag	UNP P62987
F	-1	ASP	-	expression tag	UNP P62987
F	0	LYS	-	expression tag	UNP P62987
F	4	SER	PHE	variant	UNP P62987

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Chain	Residue	Modelled	Actual	Comment	Reference
F	9	MET	THR	variant	UNP P62987
F	10	ARG	GLY	variant	UNP P62987
F	12	ARG	THR	variant	UNP P62987
F	46	SER	ALA	variant	UNP P62987
F	48	MET	LYS	variant	UNP P62987
F	49	LEU	GLN	variant	UNP P62987
F	64	GLY	GLU	variant	UNP P62987
F	68	THR	HIS	variant	UNP P62987
F	70	GLY	VAL	variant	UNP P62987
F	72	ILE	ARG	variant	UNP P62987
F	75	ALA	GLY	variant	UNP P62987
F	76	ASN	GLY	variant	UNP P62987
F	77	GLY	-	insertion	UNP P62987
F	78	VAL	ILE	variant	UNP P62987
G	-24	HIS	-	expression tag	UNP P62987
G	-23	HIS	-	expression tag	UNP P62987
G	-22	HIS	-	expression tag	UNP P62987
G	-21	HIS	-	expression tag	UNP P62987
G	-20	HIS	-	expression tag	UNP P62987
G	-19	HIS	-	expression tag	UNP P62987
G	-18	VAL	-	expression tag	UNP P62987
G	-17	THR	-	expression tag	UNP P62987
G	-16	SER	-	expression tag	UNP P62987
G	-15	LEU	-	expression tag	UNP P62987
G	-14	TYR	-	expression tag	UNP P62987
G	-13	LYS	-	expression tag	UNP P62987
G	-12	LYS	-	expression tag	UNP P62987
G	-11	ALA	-	expression tag	UNP P62987
G	-10	GLY	-	expression tag	UNP P62987
G	-9	SER	-	expression tag	UNP P62987
G	-8	THR	-	expression tag	UNP P62987
G	-7	ASP	-	expression tag	UNP P62987
G	-6	TYR	-	expression tag	UNP P62987
G	-5	LYS	-	expression tag	UNP P62987
G	-4	ASP	-	expression tag	UNP P62987
G	-3	ASP	-	expression tag	UNP P62987
G	-2	ASP	-	expression tag	UNP P62987
G	-1	ASP	-	expression tag	UNP P62987
G	0	LYS	-	expression tag	UNP P62987
G	4	SER	PHE	variant	UNP P62987
G	9	MET	THR	variant	UNP P62987
G	10	ARG	GLY	variant	UNP P62987

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Chain	Residue	Modelled	Actual	Comment	Reference
G	12	ARG	THR	variant	UNP P62987
G	46	SER	ALA	variant	UNP P62987
G	48	MET	LYS	variant	UNP P62987
G	49	LEU	GLN	variant	UNP P62987
G	64	GLY	GLU	variant	UNP P62987
G	68	THR	HIS	variant	UNP P62987
G	70	GLY	VAL	variant	UNP P62987
G	72	ILE	ARG	variant	UNP P62987
G	75	ALA	GLY	variant	UNP P62987
G	76	ASN	GLY	variant	UNP P62987
G	77	GLY	-	insertion	UNP P62987
G	78	VAL	ILE	variant	UNP P62987
K	-24	HIS	-	expression tag	UNP P62987
K	-23	HIS	-	expression tag	UNP P62987
K	-22	HIS	-	expression tag	UNP P62987
K	-21	HIS	-	expression tag	UNP P62987
K	-20	HIS	-	expression tag	UNP P62987
K	-19	HIS	-	expression tag	UNP P62987
K	-18	VAL	-	expression tag	UNP P62987
K	-17	THR	-	expression tag	UNP P62987
K	-16	SER	-	expression tag	UNP P62987
K	-15	LEU	-	expression tag	UNP P62987
K	-14	TYR	-	expression tag	UNP P62987
K	-13	LYS	-	expression tag	UNP P62987
K	-12	LYS	-	expression tag	UNP P62987
K	-11	ALA	-	expression tag	UNP P62987
K	-10	GLY	-	expression tag	UNP P62987
K	-9	SER	-	expression tag	UNP P62987
K	-8	THR	-	expression tag	UNP P62987
K	-7	ASP	-	expression tag	UNP P62987
K	-6	TYR	-	expression tag	UNP P62987
K	-5	LYS	-	expression tag	UNP P62987
K	-4	ASP	-	expression tag	UNP P62987
K	-3	ASP	-	expression tag	UNP P62987
K	-2	ASP	-	expression tag	UNP P62987
K	-1	ASP	-	expression tag	UNP P62987
K	0	LYS	-	expression tag	UNP P62987
K	4	SER	PHE	variant	UNP P62987
K	9	MET	THR	variant	UNP P62987
K	10	ARG	GLY	variant	UNP P62987
K	12	ARG	THR	variant	UNP P62987
K	46	SER	ALA	variant	UNP P62987

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Chain	Residue	Modelled	Actual	Comment	Reference
K	48	MET	LYS	variant	UNP P62987
K	49	LEU	GLN	variant	UNP P62987
K	64	GLY	GLU	variant	UNP P62987
K	68	THR	HIS	variant	UNP P62987
K	70	GLY	VAL	variant	UNP P62987
K	72	ILE	ARG	variant	UNP P62987
K	75	ALA	GLY	variant	UNP P62987
K	76	ASN	GLY	variant	UNP P62987
K	77	GLY	-	insertion	UNP P62987
K	78	VAL	ILE	variant	UNP P62987
L	-24	HIS	-	expression tag	UNP P62987
L	-23	HIS	-	expression tag	UNP P62987
L	-22	HIS	-	expression tag	UNP P62987
L	-21	HIS	-	expression tag	UNP P62987
L	-20	HIS	-	expression tag	UNP P62987
L	-19	HIS	-	expression tag	UNP P62987
L	-18	VAL	-	expression tag	UNP P62987
L	-17	THR	-	expression tag	UNP P62987
L	-16	SER	-	expression tag	UNP P62987
L	-15	LEU	-	expression tag	UNP P62987
L	-14	TYR	-	expression tag	UNP P62987
L	-13	LYS	-	expression tag	UNP P62987
L	-12	LYS	-	expression tag	UNP P62987
L	-11	ALA	-	expression tag	UNP P62987
L	-10	GLY	-	expression tag	UNP P62987
L	-9	SER	-	expression tag	UNP P62987
L	-8	THR	-	expression tag	UNP P62987
L	-7	ASP	-	expression tag	UNP P62987
L	-6	TYR	-	expression tag	UNP P62987
L	-5	LYS	-	expression tag	UNP P62987
L	-4	ASP	-	expression tag	UNP P62987
L	-3	ASP	-	expression tag	UNP P62987
L	-2	ASP	-	expression tag	UNP P62987
L	-1	ASP	-	expression tag	UNP P62987
L	0	LYS	-	expression tag	UNP P62987
L	4	SER	PHE	variant	UNP P62987
L	9	MET	THR	variant	UNP P62987
L	10	ARG	GLY	variant	UNP P62987
L	12	ARG	THR	variant	UNP P62987
L	46	SER	ALA	variant	UNP P62987
L	48	MET	LYS	variant	UNP P62987
L	49	LEU	GLN	variant	UNP P62987

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Chain	Residue	Modelled	Actual	Comment	Reference
L	64	GLY	GLU	variant	UNP P62987
L	68	THR	HIS	variant	UNP P62987
L	70	GLY	VAL	variant	UNP P62987
L	72	ILE	ARG	variant	UNP P62987
L	75	ALA	GLY	variant	UNP P62987
L	76	ASN	GLY	variant	UNP P62987
L	77	GLY	-	insertion	UNP P62987
L	78	VAL	ILE	variant	UNP P62987
H	-24	HIS	-	expression tag	UNP P62987
H	-23	HIS	-	expression tag	UNP P62987
H	-22	HIS	-	expression tag	UNP P62987
H	-21	HIS	-	expression tag	UNP P62987
H	-20	HIS	-	expression tag	UNP P62987
H	-19	HIS	-	expression tag	UNP P62987
H	-18	VAL	-	expression tag	UNP P62987
H	-17	THR	-	expression tag	UNP P62987
H	-16	SER	-	expression tag	UNP P62987
H	-15	LEU	-	expression tag	UNP P62987
H	-14	TYR	-	expression tag	UNP P62987
H	-13	LYS	-	expression tag	UNP P62987
H	-12	LYS	-	expression tag	UNP P62987
H	-11	ALA	-	expression tag	UNP P62987
H	-10	GLY	-	expression tag	UNP P62987
H	-9	SER	-	expression tag	UNP P62987
H	-8	THR	-	expression tag	UNP P62987
H	-7	ASP	-	expression tag	UNP P62987
H	-6	TYR	-	expression tag	UNP P62987
H	-5	LYS	-	expression tag	UNP P62987
H	-4	ASP	-	expression tag	UNP P62987
H	-3	ASP	-	expression tag	UNP P62987
H	-2	ASP	-	expression tag	UNP P62987
H	-1	ASP	-	expression tag	UNP P62987
H	0	LYS	-	expression tag	UNP P62987
H	4	SER	PHE	variant	UNP P62987
H	9	MET	THR	variant	UNP P62987
H	10	ARG	GLY	variant	UNP P62987
H	12	ARG	THR	variant	UNP P62987
H	46	SER	ALA	variant	UNP P62987
H	48	MET	LYS	variant	UNP P62987
H	49	LEU	GLN	variant	UNP P62987
H	64	GLY	GLU	variant	UNP P62987
H	68	THR	HIS	variant	UNP P62987

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Chain	Residue	Modelled	Actual	Comment	Reference
H	70	GLY	VAL	variant	UNP P62987
H	72	ILE	ARG	variant	UNP P62987
H	75	ALA	GLY	variant	UNP P62987
H	76	ASN	GLY	variant	UNP P62987
H	77	GLY	-	insertion	UNP P62987
H	78	VAL	ILE	variant	UNP P62987
J	-24	HIS	-	expression tag	UNP P62987
J	-23	HIS	-	expression tag	UNP P62987
J	-22	HIS	-	expression tag	UNP P62987
J	-21	HIS	-	expression tag	UNP P62987
J	-20	HIS	-	expression tag	UNP P62987
J	-19	HIS	-	expression tag	UNP P62987
J	-18	VAL	-	expression tag	UNP P62987
J	-17	THR	-	expression tag	UNP P62987
J	-16	SER	-	expression tag	UNP P62987
J	-15	LEU	-	expression tag	UNP P62987
J	-14	TYR	-	expression tag	UNP P62987
J	-13	LYS	-	expression tag	UNP P62987
J	-12	LYS	-	expression tag	UNP P62987
J	-11	ALA	-	expression tag	UNP P62987
J	-10	GLY	-	expression tag	UNP P62987
J	-9	SER	-	expression tag	UNP P62987
J	-8	THR	-	expression tag	UNP P62987
J	-7	ASP	-	expression tag	UNP P62987
J	-6	TYR	-	expression tag	UNP P62987
J	-5	LYS	-	expression tag	UNP P62987
J	-4	ASP	-	expression tag	UNP P62987
J	-3	ASP	-	expression tag	UNP P62987
J	-2	ASP	-	expression tag	UNP P62987
J	-1	ASP	-	expression tag	UNP P62987
J	0	LYS	-	expression tag	UNP P62987
J	4	SER	PHE	variant	UNP P62987
J	9	MET	THR	variant	UNP P62987
J	10	ARG	GLY	variant	UNP P62987
J	12	ARG	THR	variant	UNP P62987
J	46	SER	ALA	variant	UNP P62987
J	48	MET	LYS	variant	UNP P62987
J	49	LEU	GLN	variant	UNP P62987
J	64	GLY	GLU	variant	UNP P62987
J	68	THR	HIS	variant	UNP P62987
J	70	GLY	VAL	variant	UNP P62987
J	72	ILE	ARG	variant	UNP P62987

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Chain	Residue	Modelled	Actual	Comment	Reference
J	75	ALA	GLY	variant	UNP P62987
J	76	ASN	GLY	variant	UNP P62987
J	77	GLY	-	insertion	UNP P62987
J	78	VAL	ILE	variant	UNP P62987

- Molecule 3 is ZINC ION (three-letter code: ZN) (formula: Zn) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	B	1	Total Zn 1 1	0	0
3	A	1	Total Zn 1 1	0	0
3	C	1	Total Zn 1 1	0	0
3	D	1	Total Zn 1 1	0	0

- Molecule 4 is BROMIDE ION (three-letter code: BR) (formula: Br).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	B	1	Total Br 1 1	0	0
4	A	2	Total Br 2 2	0	0
4	D	1	Total Br 1 1	0	0
4	E	1	Total Br 1 1	0	0
4	F	1	Total Br 1 1	0	0

- Molecule 5 is SODIUM ION (three-letter code: NA) (formula: Na).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	1	Total Na 1 1	0	0
5	C	1	Total Na 1 1	0	0

- Molecule 6 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

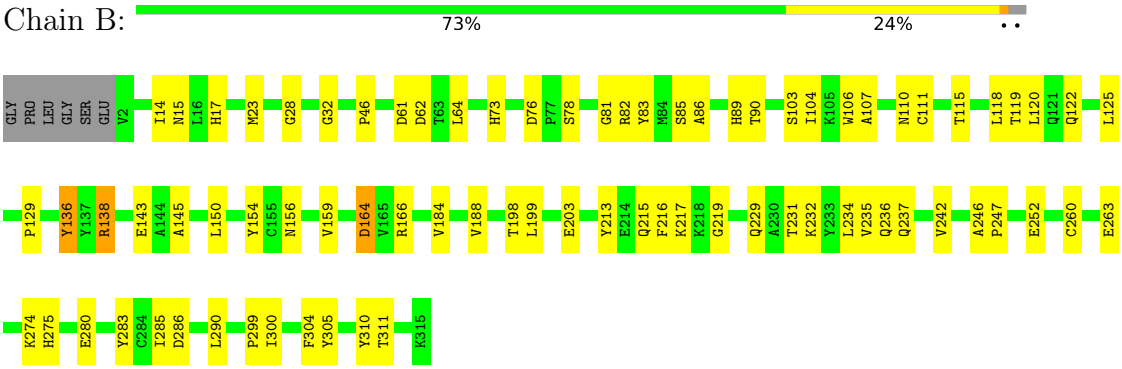
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	D	1	Total Cl 1 1	0	0

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.

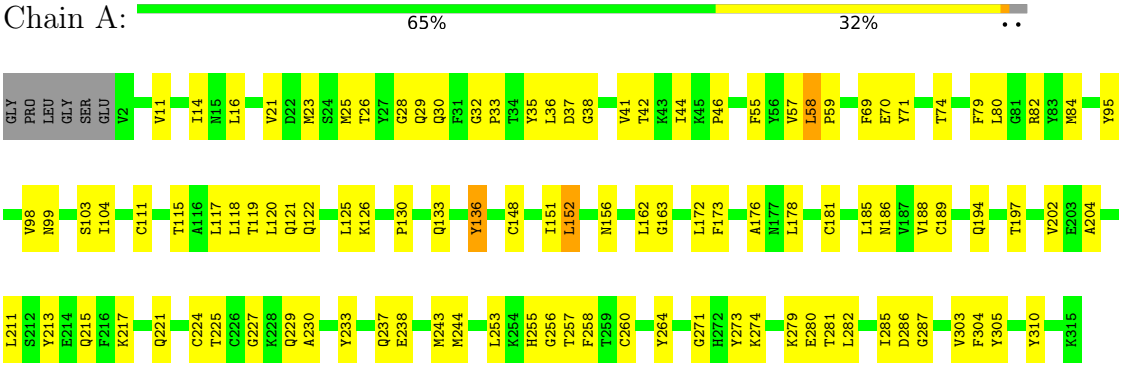
- Molecule 1: Papain-like protease nsp3

Chain B:



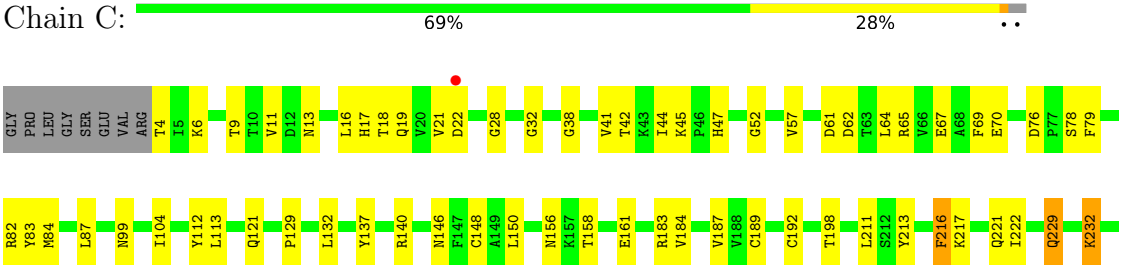
- Molecule 1: Papain-like protease nsp3

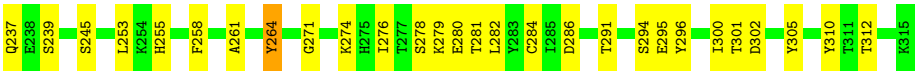
Chain A:



- Molecule 1: Papain-like protease nsp3

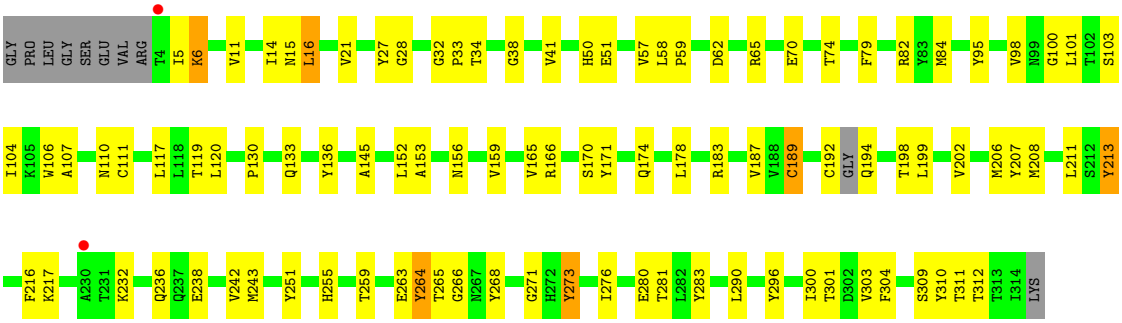
Chain C:





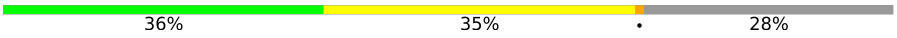
• Molecule 1: Papain-like protease nsp3

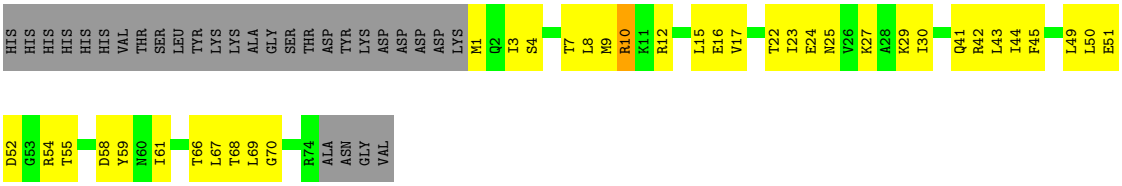
Chain D:



• Molecule 2: Ubiquitin variant UbV.CV2.1

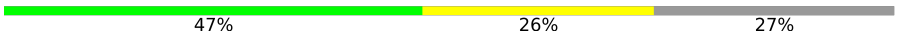
Chain E:





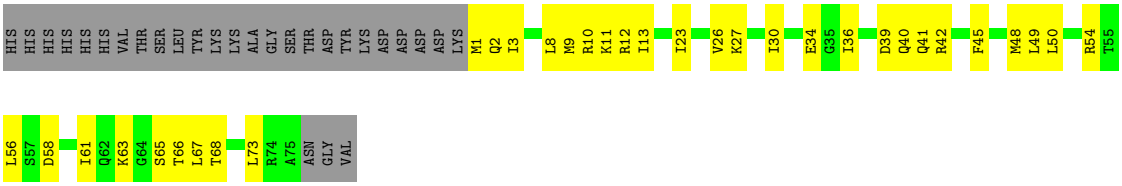
• Molecule 2: Ubiquitin variant UbV.CV2.1

Chain I:



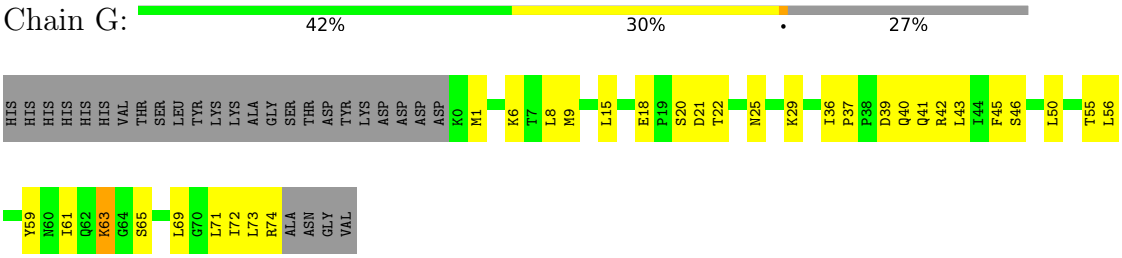
• Molecule 2: Ubiquitin variant UbV.CV2.1

Chain F:



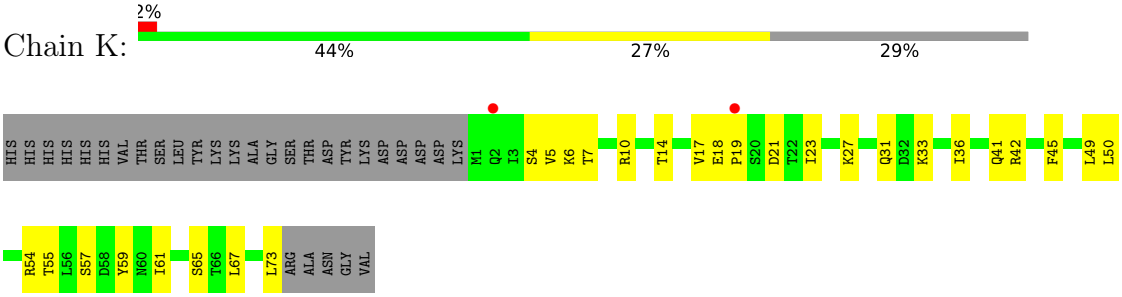
• Molecule 2: Ubiquitin variant UbV.CV2.1

Chain G:



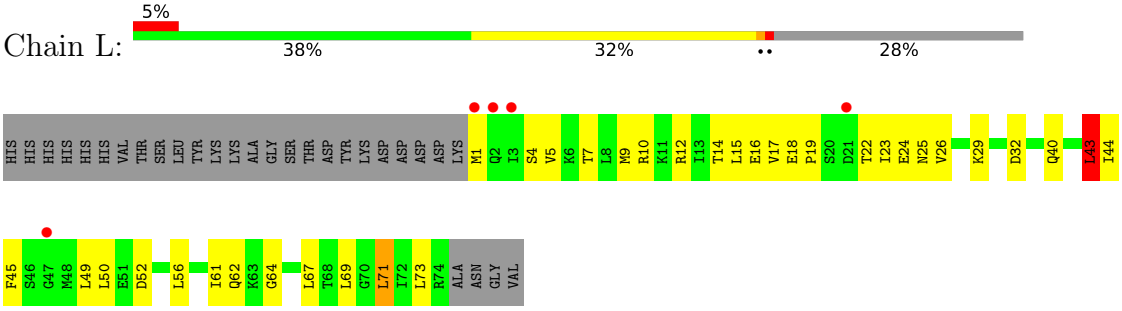
• Molecule 2: Ubiquitin variant UbV.CV2.1

Chain K:



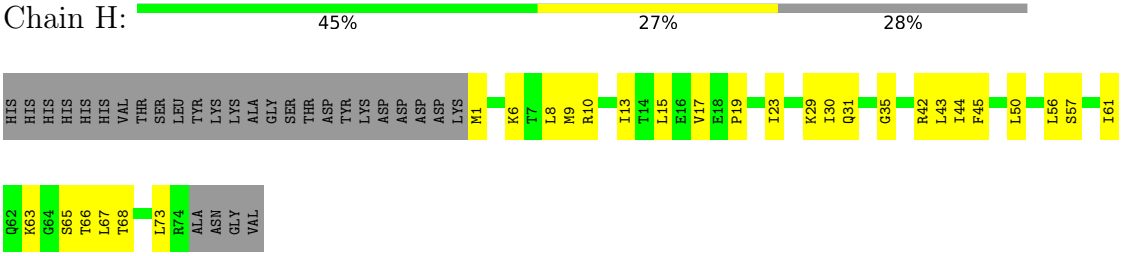
• Molecule 2: Ubiquitin variant UbV.CV2.1

Chain L:



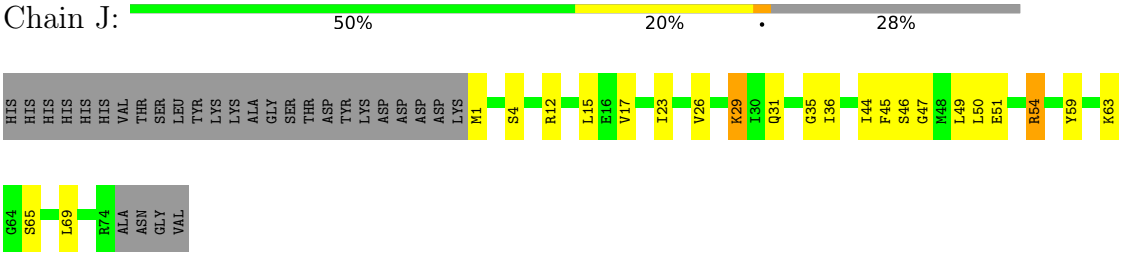
• Molecule 2: Ubiquitin variant UbV.CV2.1

Chain H:



• Molecule 2: Ubiquitin variant UbV.CV2.1

Chain J:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α, β, γ	54.58Å 174.86Å 121.56Å 90.00° 95.74° 90.00°	Depositor
Resolution (Å)	47.79 – 3.50 47.79 – 3.50	Depositor EDS
% Data completeness (in resolution range)	81.5 (47.79-3.50) 81.6 (47.79-3.50)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.31 (at 3.48Å)	Xtriage
Refinement program	PHENIX 1.19.2_4158	Depositor
R, R_{free}	0.225 , 0.270 0.224 , 0.269	Depositor DCC
R_{free} test set	1998 reflections (8.56%)	wwPDB-VP
Wilson B-factor (Å^2)	61.5	Xtriage
Anisotropy	0.562	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.30 , 45.8	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	0.88	EDS
Total number of atoms	13989	wwPDB-VP
Average B, all atoms (Å^2)	63.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.30% 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, NA, CL, BR

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	0.29	0/2509	0.55	1/3416 (0.0%)
1	B	0.30	0/2488	0.55	0/3390
1	C	0.29	0/2445	0.59	0/3338
1	D	0.32	0/2422	0.59	0/3310
2	E	0.30	0/575	0.66	0/775
2	F	0.35	0/584	0.68	0/785
2	G	0.31	0/549	0.68	1/742 (0.1%)
2	H	0.37	0/549	0.70	2/741 (0.3%)
2	I	0.30	0/555	0.70	0/750
2	J	0.31	0/534	0.70	0/724
2	K	0.27	0/509	0.61	0/692
2	L	0.30	0/520	0.78	2/707 (0.3%)
All	All	0.31	0/14239	0.61	6/19370 (0.0%)

There are no bond length outliers.

The worst 5 of 6 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	L	43	LEU	CA-CB-CG	6.71	130.74	115.30
1	A	58	LEU	CA-CB-CG	5.63	128.26	115.30
2	H	8	LEU	CA-CB-CG	5.34	127.58	115.30
2	H	43	LEU	CA-CB-CG	5.34	127.57	115.30
2	G	71	LEU	CA-CB-CG	5.18	127.22	115.30

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	2450	0	2324	81	0
1	B	2430	0	2295	55	1
1	C	2387	0	2220	77	0
1	D	2365	0	2185	84	0
2	E	571	0	592	48	0
2	F	580	0	615	29	0
2	G	545	0	548	32	0
2	H	545	0	556	21	1
2	I	551	0	566	29	0
2	J	530	0	519	24	0
2	K	505	0	470	27	0
2	L	517	0	490	33	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
3	C	1	0	0	0	0
3	D	1	0	0	0	0
4	A	2	0	0	3	0
4	B	1	0	0	1	0
4	D	1	0	0	0	0
4	E	1	0	0	0	0
4	F	1	0	0	1	0
5	A	1	0	0	0	0
5	C	1	0	0	0	0
6	D	1	0	0	1	0
All	All	13989	0	13380	461	1

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 17.

The worst 5 of 461 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:189:CYS:SG	1:D:192:CYS:HB2	1.99	1.02
1:D:283:TYR:HD2	1:D:290:LEU:HD11	1.36	0.91
2:J:23:ILE:HD11	2:J:54:ARG:H	1.36	0.87

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:23:MET:HA	1:A:46:PRO:HG2	1.59	0.85
2:I:15:LEU:HD22	2:I:29:LYS:HZ3	1.44	0.82

All (1) 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)
1:B:215:GLN:OE1	2:H:57:SER:OG[1_554]	2.07	0.13

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	312/320 (98%)	295 (95%)	17 (5%)	0	100	100
1	B	312/320 (98%)	293 (94%)	19 (6%)	0	100	100
1	C	310/320 (97%)	284 (92%)	26 (8%)	0	100	100
1	D	306/320 (96%)	284 (93%)	22 (7%)	0	100	100
2	E	72/103 (70%)	65 (90%)	7 (10%)	0	100	100
2	F	73/103 (71%)	65 (89%)	8 (11%)	0	100	100
2	G	73/103 (71%)	65 (89%)	8 (11%)	0	100	100
2	H	72/103 (70%)	64 (89%)	8 (11%)	0	100	100
2	I	73/103 (71%)	64 (88%)	9 (12%)	0	100	100
2	J	72/103 (70%)	68 (94%)	4 (6%)	0	100	100
2	K	71/103 (69%)	67 (94%)	4 (6%)	0	100	100
2	L	72/103 (70%)	59 (82%)	13 (18%)	0	100	100
All	All	1818/2104 (86%)	1673 (92%)	145 (8%)	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	263/278 (95%)	260 (99%)	3 (1%)	73 88
1	B	259/278 (93%)	253 (98%)	6 (2%)	50 77
1	C	250/278 (90%)	243 (97%)	7 (3%)	43 72
1	D	246/278 (88%)	238 (97%)	8 (3%)	38 68
2	E	65/93 (70%)	64 (98%)	1 (2%)	65 84
2	F	66/93 (71%)	66 (100%)	0	100 100
2	G	58/93 (62%)	55 (95%)	3 (5%)	23 56
2	H	58/93 (62%)	57 (98%)	1 (2%)	60 82
2	I	59/93 (63%)	59 (100%)	0	100 100
2	J	54/93 (58%)	50 (93%)	4 (7%)	13 44
2	K	48/93 (52%)	47 (98%)	1 (2%)	53 79
2	L	50/93 (54%)	47 (94%)	3 (6%)	19 52
All	All	1476/1856 (80%)	1439 (98%)	37 (2%)	47 75

5 of 37 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	K	33	LYS
2	J	54	ARG
2	L	32	ASP
2	H	15	LEU
1	C	216	PHE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	174	GLN
1	C	17	HIS
1	C	308	ASN
2	F	41	GLN

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Mol	Chain	Res	Type
2	G	41	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 13 ligands modelled in this entry, 13 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	314/320 (98%)	-0.29	0 100 100	26, 49, 89, 107	0
1	B	314/320 (98%)	-0.39	0 100 100	28, 46, 74, 102	0
1	C	312/320 (97%)	-0.25	1 (0%) 94 91	24, 53, 95, 108	0
1	D	310/320 (96%)	-0.09	2 (0%) 89 86	37, 70, 102, 150	0
2	E	74/103 (71%)	-0.35	0 100 100	35, 50, 77, 93	0
2	F	75/103 (72%)	-0.28	0 100 100	35, 50, 70, 98	0
2	G	75/103 (72%)	0.08	0 100 100	35, 68, 91, 101	0
2	H	74/103 (71%)	-0.10	0 100 100	44, 67, 99, 107	0
2	I	75/103 (72%)	-0.13	0 100 100	43, 82, 101, 108	0
2	J	74/103 (71%)	0.03	0 100 100	59, 92, 102, 119	0
2	K	73/103 (70%)	0.22	2 (2%) 54 48	67, 96, 119, 134	0
2	L	74/103 (71%)	0.57	5 (6%) 17 16	87, 109, 132, 167	0
All	All	1844/2104 (87%)	-0.17	10 (0%) 91 88	24, 59, 105, 167	0

The worst 5 of 10 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	L	2	GLN	2.9
1	D	4	THR	2.8
2	L	1	MET	2.8
2	L	21	ASP	2.7
2	L	47	GLY	2.6

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

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)

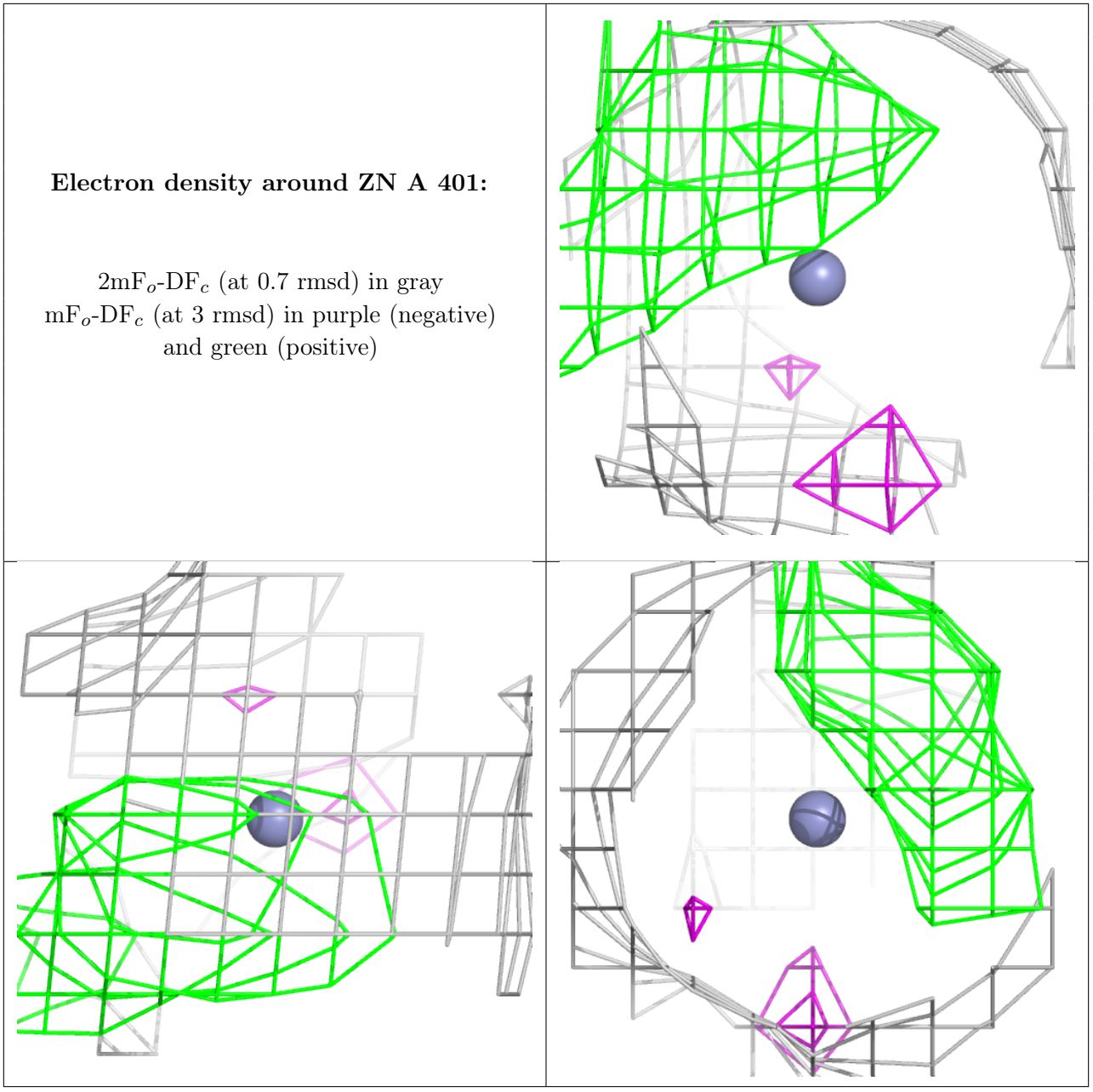
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95^{th} percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' 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
3	ZN	A	401	1/1	0.85	0.13	57,57,57,57	0
5	NA	A	404	1/1	0.86	0.45	26,26,26,26	0
4	BR	D	402	1/1	0.87	0.17	104,104,104,104	0
5	NA	C	402	1/1	0.92	0.39	36,36,36,36	0
4	BR	A	403	1/1	0.93	0.08	86,86,86,86	0
4	BR	A	402	1/1	0.94	0.09	87,87,87,87	0
6	CL	D	403	1/1	0.94	0.15	53,53,53,53	0
4	BR	F	101	1/1	0.95	0.07	82,82,82,82	0
4	BR	B	402	1/1	0.95	0.06	77,77,77,77	0
3	ZN	B	401	1/1	0.96	0.08	66,66,66,66	0
3	ZN	D	401	1/1	0.96	0.12	185,185,185,185	0
4	BR	E	101	1/1	0.97	0.07	62,62,62,62	0
3	ZN	C	401	1/1	0.99	0.10	49,49,49,49	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.

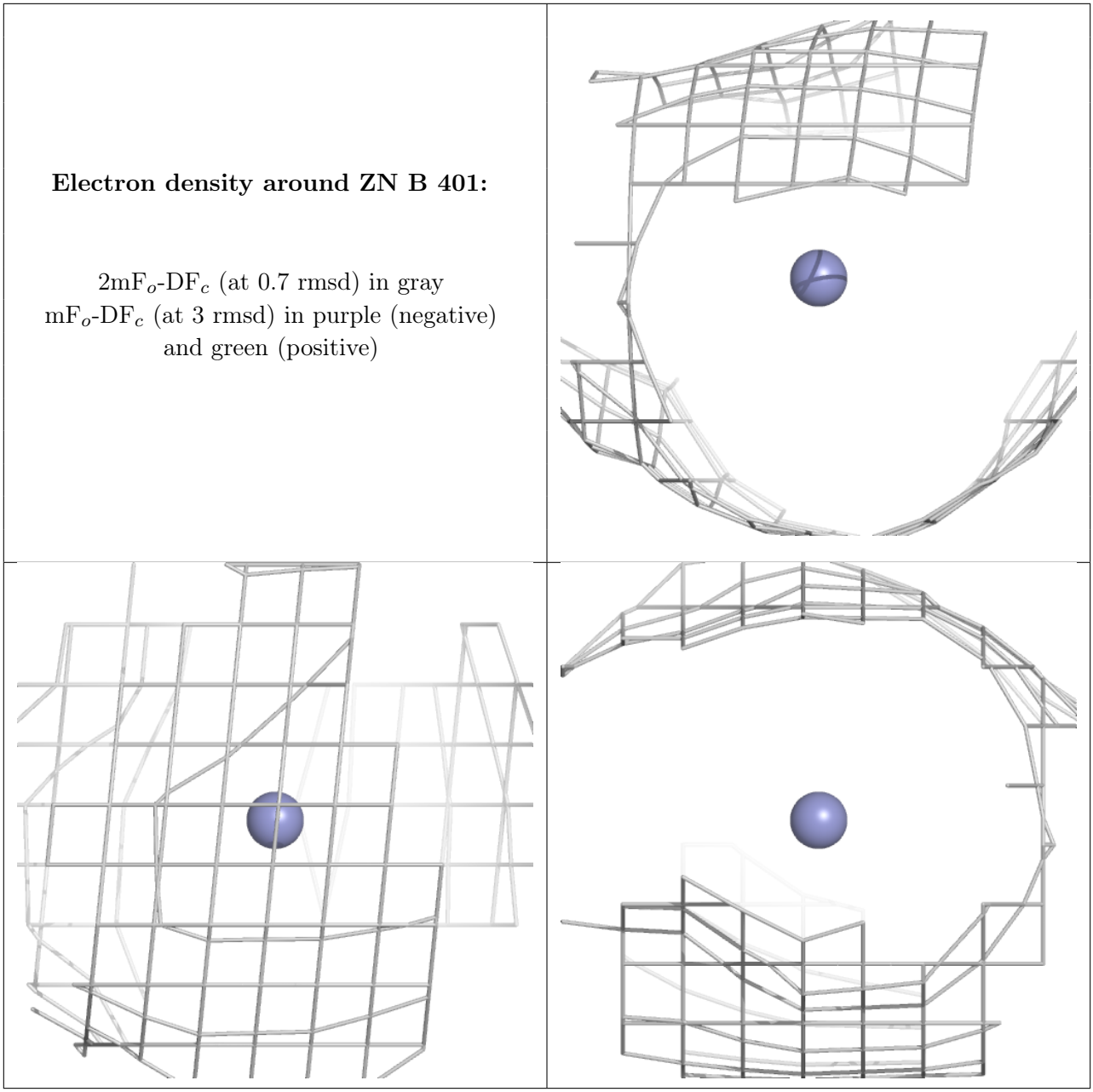
Electron density around ZN A 401:

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



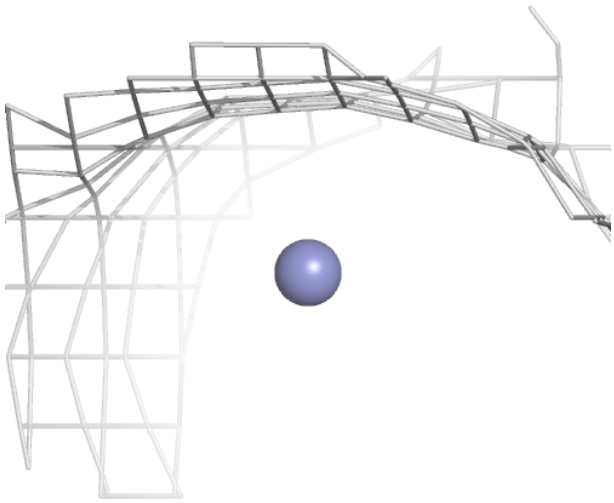
Electron density around ZN B 401:

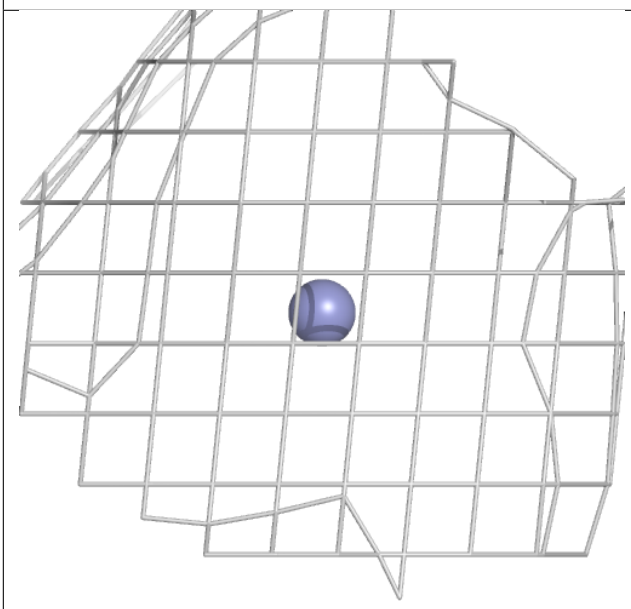
$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

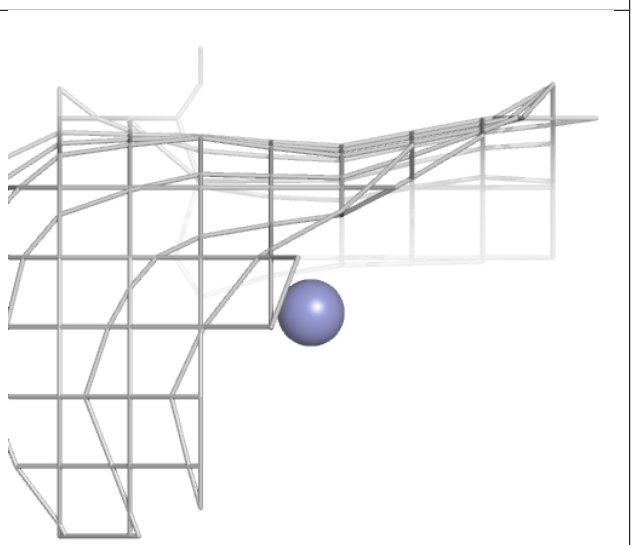


Electron density around ZN D 401:

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

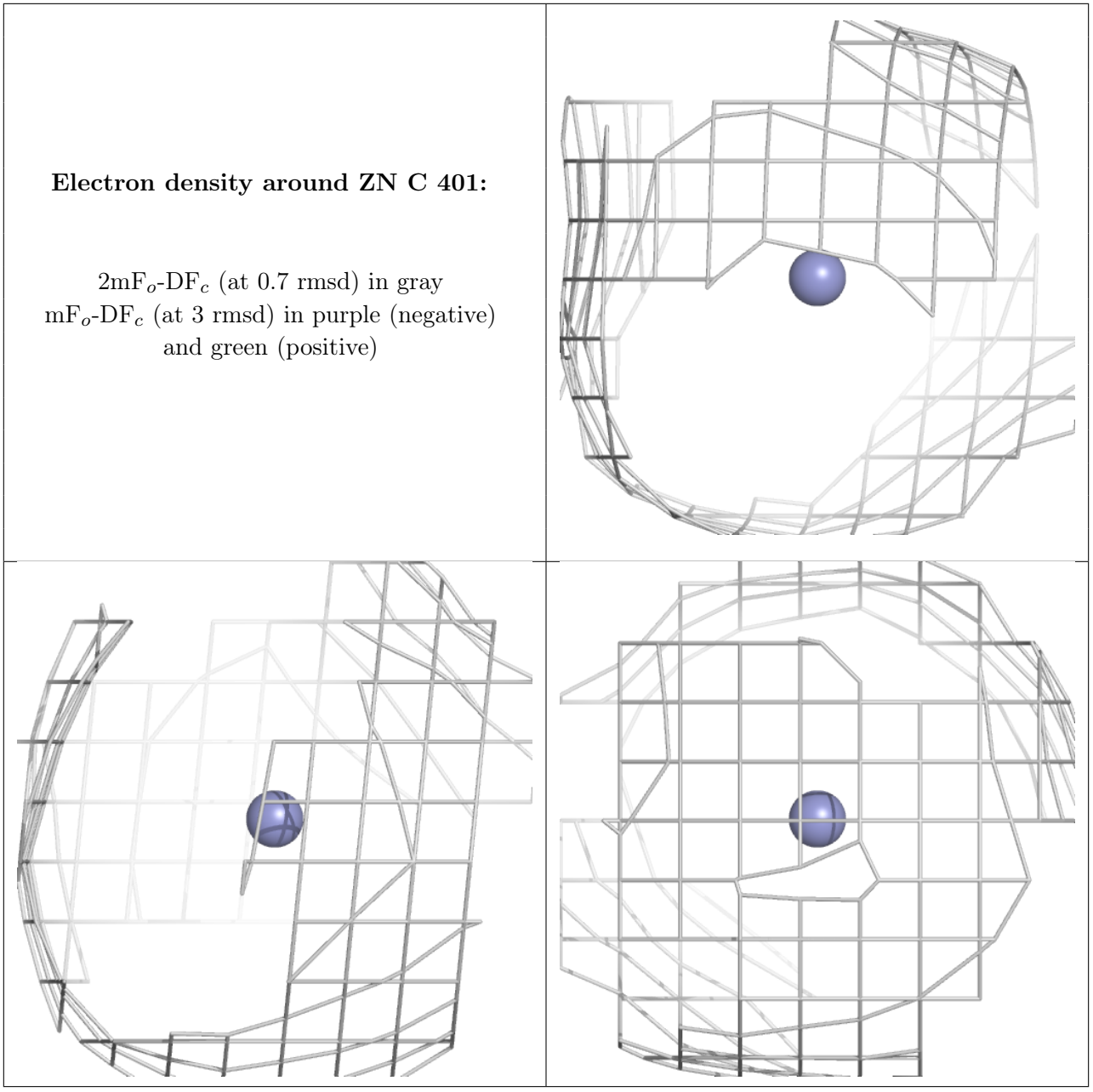






Electron density around ZN C 401:

$2mF_o$-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.