



Full wwPDB X-ray Structure Validation Report ⓘ

Sep 30, 2021 – 04:12 PM JST

PDB ID : 6KYV
Title : Crystal Structure of RIG-I and hairpin RNA with G-U wobble base pairs
Authors : Kim, K.-H.; Hwang, J.; Kim, J.H.; Son, K.-P.; Jang, Y.; Kim, M.; Kang, S.-J.;
Lee, J.-O.; Choi, B.-S.
Deposited on : 2019-09-20
Resolution : 3.00 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 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 : 2.23.2
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) : 2.23.2

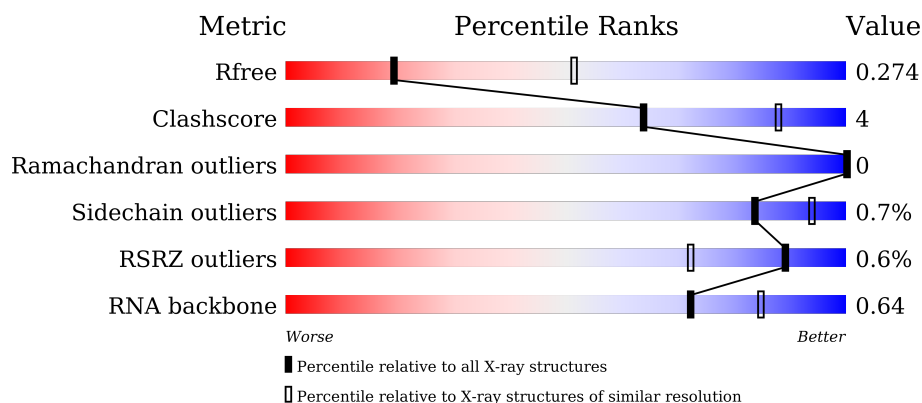
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.00 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	2092 (3.00-3.00)
Clashscore	141614	2416 (3.00-3.00)
Ramachandran outliers	138981	2333 (3.00-3.00)
Sidechain outliers	138945	2336 (3.00-3.00)
RSRZ outliers	127900	1990 (3.00-3.00)
RNA backbone	3102	1173 (3.30-2.70)

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 $\leq 5\%$. 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	22	<div> <div></div> <div>82%</div> <div>5%</div> <div>14%</div> </div>
1	C	22	<div> <div></div> <div>82%</div> <div>9%</div> <div>9%</div> </div>
1	E	22	<div> <div></div> <div>82%</div> <div>9%</div> <div>9%</div> </div>
1	G	22	<div> <div></div> <div>59%</div> <div>32%</div> <div>9%</div> </div>

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Mol	Chain	Length	Quality of chain
1	I	22	<div> <div>5%</div> <div>82%</div> <div>9%</div> <div>9%</div> </div>
1	K	22	<div> <div>9%</div> <div>73%</div> <div>18%</div> <div>9%</div> </div>
2	B	684	<div> <div>87%</div> <div>11%</div> <div>.</div> </div>
2	D	684	<div> <div>87%</div> <div>11%</div> <div>.</div> </div>
2	F	684	<div> <div>%</div> <div>86%</div> <div>12%</div> <div>.</div> </div>
2	H	684	<div> <div>86%</div> <div>13%</div> <div>.</div> </div>
2	J	684	<div> <div>%</div> <div>84%</div> <div>14%</div> <div>.</div> </div>
2	L	684	<div> <div>87%</div> <div>11%</div> <div>.</div> </div>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 34612 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 RNA chain called RNA (5'-R(*GP*GP*UP*AP*GP*AP*CP*GP*CP*UP*UP*CP*GP*GP*CP*GP*UP*UP*UP*GP*CP*C)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	22	Total	C	N	O	P	0	0	0
			465	208	80	156	21			
1	C	22	Total	C	N	O	P	0	0	0
			465	208	80	156	21			
1	E	22	Total	C	N	O	P	0	0	0
			465	208	80	156	21			
1	G	22	Total	C	N	O	P	0	0	0
			465	208	80	156	21			
1	I	22	Total	C	N	O	P	0	0	0
			465	208	80	156	21			
1	K	22	Total	C	N	O	P	0	0	0
			465	208	80	156	21			

- Molecule 2 is a protein called Probable ATP-dependent RNA helicase DDX58.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	672	Total	C	N	O	S	0	0	0
			5288	3374	900	981	33			
2	D	672	Total	C	N	O	S	0	0	0
			5294	3377	902	982	33			
2	F	672	Total	C	N	O	S	0	0	0
			5289	3376	904	976	33			
2	H	672	Total	C	N	O	S	0	0	0
			5329	3403	909	984	33			
2	J	672	Total	C	N	O	S	0	0	0
			5310	3388	908	981	33			
2	L	672	Total	C	N	O	S	0	0	0
			5306	3384	906	983	33			

There are 18 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	239	SER	-	expression tag	UNP O95786
B	240	ASN	-	expression tag	UNP O95786
B	241	ALA	-	expression tag	UNP O95786
D	239	SER	-	expression tag	UNP O95786
D	240	ASN	-	expression tag	UNP O95786
D	241	ALA	-	expression tag	UNP O95786
F	239	SER	-	expression tag	UNP O95786
F	240	ASN	-	expression tag	UNP O95786
F	241	ALA	-	expression tag	UNP O95786
H	239	SER	-	expression tag	UNP O95786
H	240	ASN	-	expression tag	UNP O95786
H	241	ALA	-	expression tag	UNP O95786
J	239	SER	-	expression tag	UNP O95786
J	240	ASN	-	expression tag	UNP O95786
J	241	ALA	-	expression tag	UNP O95786
L	239	SER	-	expression tag	UNP O95786
L	240	ASN	-	expression tag	UNP O95786
L	241	ALA	-	expression tag	UNP O95786


- 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	0	0
			1	1		
3	D	1	Total	Zn	0	0
			1	1		
3	F	1	Total	Zn	0	0
			1	1		
3	H	1	Total	Zn	0	0
			1	1		
3	J	1	Total	Zn	0	0
			1	1		
3	L	1	Total	Zn	0	0
			1	1		

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: RNA (5'-R(*GP*GP*UP*AP*GP*AP*CP*GP*CP*UP*UP*CP*GP*GP*CP*GP*UP*UP*UP*GP*CP*C)-3')

Chain A: 




- Molecule 1: RNA (5'-R(*GP*GP*UP*AP*GP*AP*CP*GP*CP*UP*UP*CP*GP*GP*CP*GP*UP*UP*UP*GP*CP*C)-3')

Chain C: 



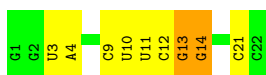
- Molecule 1: RNA (5'-R(*GP*GP*UP*AP*GP*AP*CP*GP*CP*UP*UP*CP*GP*GP*CP*GP*UP*UP*UP*GP*CP*C)-3')

Chain E: 




- Molecule 1: RNA (5'-R(*GP*GP*UP*AP*GP*AP*CP*GP*CP*UP*UP*CP*GP*GP*CP*GP*UP*UP*UP*GP*CP*C)-3')

Chain G: 

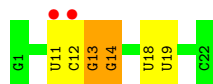


- Molecule 1: RNA (5'-R(*GP*GP*UP*AP*GP*AP*CP*GP*CP*UP*UP*CP*GP*GP*CP*GP*UP*UP*UP*GP*CP*C)-3')

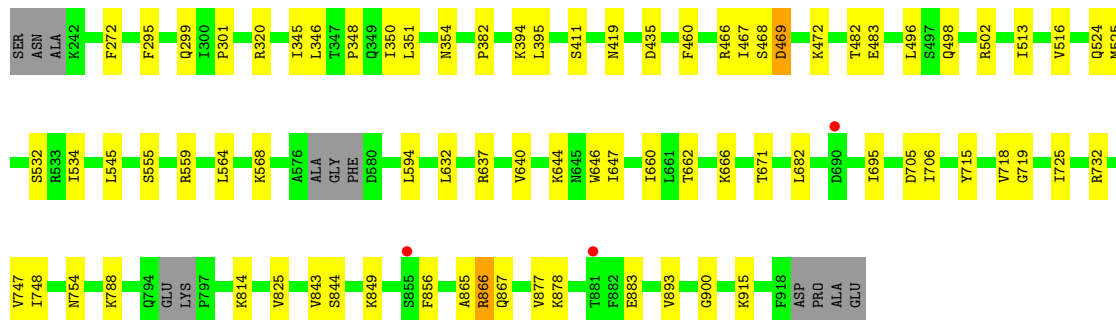
Chain I: 



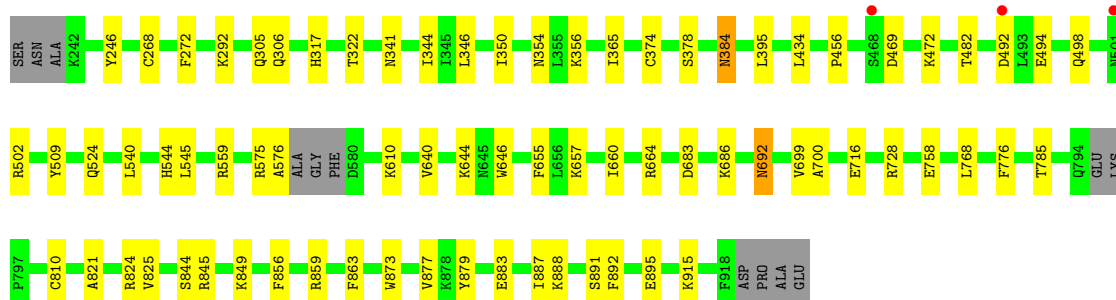
- Molecule 1: RNA (5'-R(*GP*GP*UP*AP*GP*AP*CP*GP*CP*UP*UP*CP*GP*GP*CP*GP*UP*UP*UP*GP*CP*C)-3')



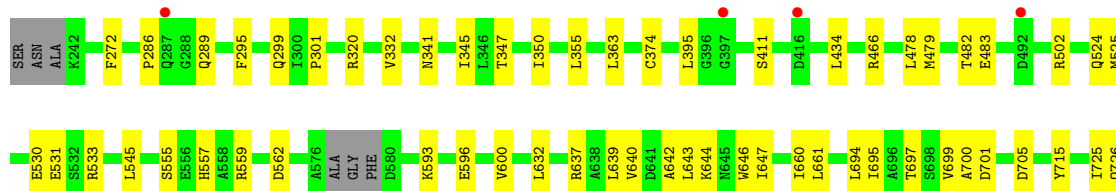
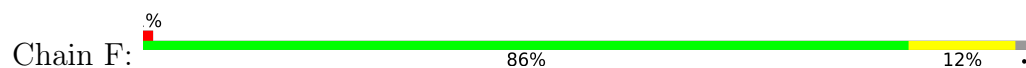
- Molecule 2: Probable ATP-dependent RNA helicase DDX58



- Molecule 2: Probable ATP-dependent RNA helicase DDX58



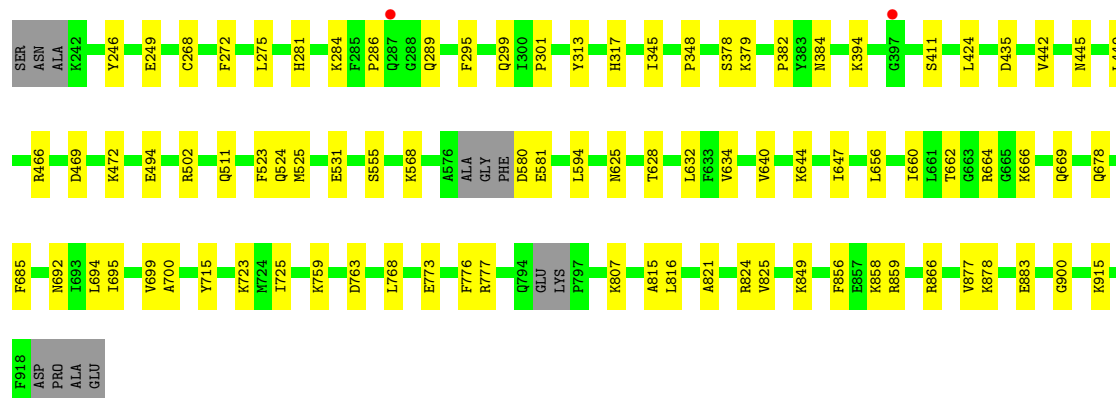
- Molecule 2: Probable ATP-dependent RNA helicase DDX58





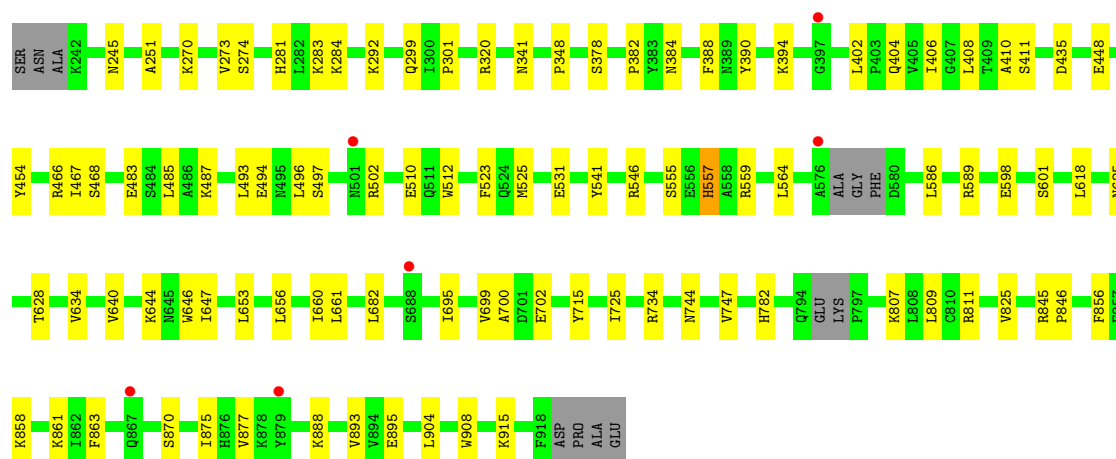
- Molecule 2: Probable ATP-dependent RNA helicase DDX58

Chain H: 86% 13%



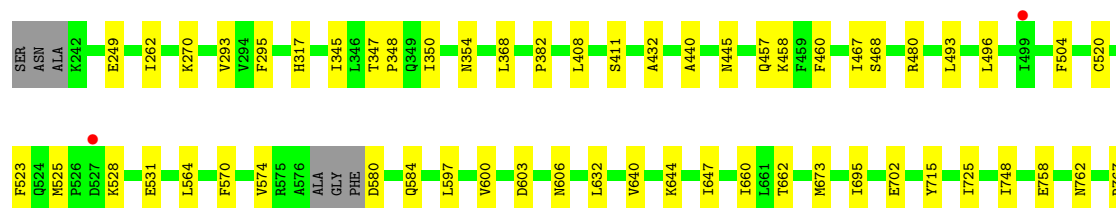
- Molecule 2: Probable ATP-dependent RNA helicase DDX58

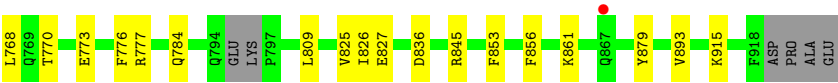
Chain J: 84% 14%



- Molecule 2: Probable ATP-dependent RNA helicase DDX58

Chain L: 87% 11%





4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	112.17Å 175.13Å 308.96Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	36.56 – 3.00 36.56 – 3.00	Depositor EDS
% Data completeness (in resolution range)	95.6 (36.56-3.00) 90.1 (36.56-3.00)	Depositor EDS
R_{merge}	0.17	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.89 (at 3.00Å)	Xtriage
Refinement program	PHENIX 1.13_2998	Depositor
R, R_{free}	0.222 , 0.274 0.222 , 0.274	Depositor DCC
R_{free} test set	5998 reflections (5.13%)	wwPDB-VP
Wilson B-factor (Å ²)	59.0	Xtriage
Anisotropy	0.307	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 34.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	34612	wwPDB-VP
Average B, all atoms (Å ²)	62.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 8.92% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ 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	0.17	0/518	0.75	0/806
1	C	0.18	0/518	0.75	0/806
1	E	0.18	0/518	0.75	0/806
1	G	0.18	0/518	0.75	0/806
1	I	0.18	0/518	0.77	0/806
1	K	0.18	0/518	0.78	0/806
2	B	0.24	0/5393	0.39	0/7292
2	D	0.24	0/5399	0.39	0/7299
2	F	0.24	0/5394	0.39	0/7291
2	H	0.24	0/5435	0.39	0/7340
2	J	0.24	0/5415	0.39	0/7317
2	L	0.24	0/5411	0.39	0/7313
All	All	0.23	0/35555	0.44	0/48688

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	465	0	238	3	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	C	465	0	238	3	0
1	E	465	0	238	2	0
1	G	465	0	238	5	0
1	I	465	0	238	1	0
1	K	465	0	238	2	0
2	B	5288	0	5232	45	0
2	D	5294	0	5241	44	0
2	F	5289	0	5245	46	0
2	H	5329	0	5316	48	0
2	J	5310	0	5283	54	0
2	L	5306	0	5267	42	0
3	B	1	0	0	0	0
3	D	1	0	0	0	0
3	F	1	0	0	0	0
3	H	1	0	0	0	0
3	J	1	0	0	0	0
3	L	1	0	0	0	0
All	All	34612	0	33012	285	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 4.

All (285) 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)
2:F:866:ARG:HG2	2:F:867:GLN:HG2	1.61	0.80
2:L:600:VAL:O	2:L:606:ASN:ND2	2.23	0.72
2:F:525:MET:HB2	2:F:531:GLU:HB2	1.71	0.71
2:D:657:LYS:H	2:D:692:ASN:HB2	1.55	0.70
2:H:825:VAL:HG21	2:H:915:LYS:HB3	1.78	0.66
1:K:13:G:O2'	1:K:14:G:O4'	2.13	0.66
2:D:859:ARG:NH1	2:D:883:GLU:OE2	2.30	0.65
2:H:466:ARG:NH1	2:H:555:SER:O	2.31	0.64
2:D:664:ARG:O	2:D:849:LYS:NZ	2.30	0.63
2:J:845:ARG:HD3	2:J:846:PRO:HD2	1.80	0.63
1:C:11:U:N3	2:D:498:GLN:O	2.30	0.63
1:E:13:G:O2'	1:E:14:G:O4'	2.14	0.63
2:H:669:GLN:O	2:H:849:LYS:NZ	2.25	0.63
1:C:13:G:O2'	1:C:14:G:O4'	2.16	0.62
1:I:13:G:O2'	1:I:14:G:O4'	2.16	0.62
2:F:559:ARG:HD3	2:F:646:TRP:HD1	1.64	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:644:LYS:HB2	2:D:660:ILE:HD11	1.81	0.61
2:F:483:GLU:OE2	2:F:502:ARG:NH2	2.33	0.61
2:H:647:ILE:HD12	2:H:695:ILE:HD11	1.83	0.61
2:J:564:LEU:HD21	2:J:598:GLU:HG2	1.83	0.61
1:G:13:G:O2'	1:G:14:G:O4'	2.16	0.60
2:H:568:LYS:HE2	2:H:594:LEU:HD11	1.83	0.60
2:J:653:LEU:HB3	2:J:656:LEU:HD12	1.84	0.60
2:L:825:VAL:HG21	2:L:915:LYS:HB3	1.82	0.60
2:B:814:LYS:HE3	2:B:893:VAL:HG11	1.83	0.59
2:F:524:GLN:NE2	2:F:900:GLY:O	2.35	0.59
2:F:478:LEU:HD21	2:F:593:LYS:HG3	1.83	0.59
2:B:705:ASP:OD2	2:B:732:ARG:NH2	2.36	0.59
2:B:395:LEU:HD12	2:B:788:LYS:HD3	1.84	0.58
2:J:466:ARG:NH1	2:J:555:SER:O	2.36	0.58
2:H:644:LYS:HD2	2:H:660:ILE:HD11	1.84	0.58
2:D:469:ASP:HB3	2:D:472:LYS:HB3	1.86	0.58
1:E:19:U:OP2	2:F:637:ARG:NH1	2.37	0.57
2:J:825:VAL:HG21	2:J:915:LYS:HB3	1.85	0.57
2:F:825:VAL:HG21	2:F:915:LYS:HB3	1.87	0.57
2:J:493:LEU:HA	2:J:496:LEU:HD13	1.87	0.57
2:J:483:GLU:HB3	2:J:487:LYS:HE3	1.86	0.57
2:J:394:LYS:NZ	2:J:435:ASP:OD1	2.32	0.57
2:L:644:LYS:HB2	2:L:660:ILE:HD11	1.86	0.56
2:J:483:GLU:OE1	2:J:502:ARG:NH2	2.38	0.56
2:L:758:GLU:O	2:L:762:ASN:ND2	2.39	0.56
2:H:494:GLU:OE2	2:H:502:ARG:NH1	2.28	0.56
2:D:640:VAL:HG12	2:D:660:ILE:HD12	1.88	0.56
2:F:807:LYS:HG3	2:F:897:ILE:HD11	1.88	0.56
2:H:424:LEU:HD11	2:H:768:LEU:HG	1.87	0.56
2:L:525:MET:O	2:L:528:LYS:NZ	2.39	0.56
2:D:559:ARG:HD3	2:D:646:TRP:HD1	1.69	0.56
2:F:859:ARG:NH1	2:F:883:GLU:OE2	2.39	0.56
2:D:825:VAL:HG21	2:D:915:LYS:HB3	1.88	0.55
2:H:524:GLN:NE2	2:H:900:GLY:O	2.40	0.55
2:L:348:PRO:HB2	2:L:382:PRO:HB2	1.88	0.55
2:B:524:GLN:NE2	2:B:900:GLY:O	2.39	0.55
2:H:859:ARG:NH1	2:H:883:GLU:OE2	2.40	0.55
2:B:350:ILE:O	2:B:354:ASN:ND2	2.30	0.55
2:B:867:GLN:H	2:D:576:ALA:HA	1.72	0.55
2:B:718:VAL:HB	2:B:747:VAL:HG13	1.89	0.55
1:A:13:G:O2'	1:A:14:G:O4'	2.16	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:637:ARG:HH22	2:B:671:THR:HB	1.71	0.54
2:H:378:SER:O	2:H:384:ASN:ND2	2.40	0.54
2:L:525:MET:HB2	2:L:531:GLU:HB2	1.90	0.54
2:D:494:GLU:OE2	2:D:502:ARG:NH1	2.41	0.54
2:J:625:ASN:O	2:J:628:THR:OG1	2.23	0.54
2:H:759:LYS:NZ	2:H:763:ASP:OD2	2.36	0.54
2:B:644:LYS:HB2	2:B:660:ILE:HD11	1.90	0.54
2:J:292:LYS:NZ	2:J:341:ASN:O	2.41	0.54
2:J:634:VAL:HG12	2:J:715:TYR:HB3	1.89	0.53
2:D:856:PHE:HD2	2:D:877:VAL:HG21	1.73	0.53
2:L:647:ILE:HD12	2:L:695:ILE:HD11	1.89	0.53
2:F:632:LEU:HD11	2:F:715:TYR:HB2	1.89	0.53
2:D:657:LYS:HB2	2:D:692:ASN:H	1.73	0.53
2:D:378:SER:O	2:D:384:ASN:ND2	2.42	0.53
2:D:610:LYS:NZ	2:D:716:GLU:OE1	2.42	0.53
2:H:525:MET:HB2	2:H:531:GLU:HB2	1.91	0.53
2:J:510:GLU:OE2	2:J:546:ARG:NH2	2.42	0.53
1:G:21:C:OP2	2:H:664:ARG:NH1	2.42	0.53
2:B:666:LYS:O	2:B:849:LYS:NZ	2.40	0.53
1:G:9:C:O2	2:H:511:GLN:NE2	2.38	0.53
2:B:844:SER:OG	2:D:844:SER:OG	2.21	0.52
2:H:662:THR:O	2:H:678:GLN:NE2	2.38	0.52
1:C:11:U:O2'	1:C:13:G:O6	2.24	0.52
2:L:293:VAL:HG22	2:L:368:LEU:HB3	1.90	0.52
2:J:644:LYS:HB2	2:J:660:ILE:HD11	1.92	0.52
2:B:825:VAL:HG21	2:B:915:LYS:HB3	1.92	0.52
2:J:888:LYS:HA	2:J:908:TRP:HE1	1.75	0.52
2:B:467:ILE:HG13	2:B:468:SER:H	1.73	0.52
2:F:699:VAL:HG12	2:F:700:ALA:H	1.74	0.52
2:J:497:SER:HG	2:J:512:TRP:HD1	1.57	0.52
2:H:625:ASN:O	2:H:628:THR:OG1	2.24	0.51
2:L:640:VAL:HG12	2:L:660:ILE:HD12	1.92	0.51
2:H:299:GLN:HB3	2:H:301:PRO:HD2	1.91	0.51
2:J:411:SER:HB2	2:J:725:ILE:HD12	1.91	0.51
2:J:661:LEU:HD22	2:J:682:LEU:HD21	1.92	0.51
2:F:411:SER:HB2	2:F:725:ILE:HD12	1.93	0.51
2:H:348:PRO:HB2	2:H:382:PRO:HB2	1.93	0.51
2:D:292:LYS:HD2	2:D:365:ILE:HG13	1.93	0.51
2:J:586:LEU:HD23	2:J:589:ARG:HD2	1.93	0.51
2:F:466:ARG:NH1	2:F:555:SER:O	2.39	0.51
2:L:262:ILE:HB	2:L:408:LEU:HD23	1.94	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:L:350:ILE:O	2:L:354:ASN:ND2	2.34	0.50
2:H:821:ALA:O	2:H:824:ARG:NH1	2.41	0.50
2:H:858:LYS:HA	2:H:877:VAL:HG12	1.93	0.50
2:B:532:SER:OG	2:B:866:ARG:NH2	2.43	0.50
2:B:559:ARG:HD3	2:B:646:TRP:HD1	1.76	0.50
2:J:274:SER:HB3	2:J:408:LEU:HD11	1.93	0.50
2:J:647:ILE:HD12	2:J:695:ILE:HD11	1.92	0.50
2:J:454:TYR:OH	2:J:734:ARG:NH2	2.44	0.50
2:F:856:PHE:HD2	2:F:877:VAL:HG21	1.76	0.50
2:H:249:GLU:OE1	2:H:445:ASN:ND2	2.40	0.50
2:J:281:HIS:HA	2:J:284:LYS:HE2	1.93	0.50
2:L:480:ARG:HG3	2:L:504:PHE:HE2	1.77	0.50
2:H:666:LYS:O	2:H:849:LYS:NZ	2.44	0.50
2:F:286:PRO:HD2	2:F:289:GLN:HB2	1.94	0.49
2:D:524:GLN:NE2	2:D:895:GLU:OE2	2.45	0.49
2:J:245:ASN:N	2:J:448:GLU:OE2	2.44	0.49
2:J:348:PRO:HB2	2:J:382:PRO:HB2	1.94	0.49
2:J:807:LYS:HB2	2:J:895:GLU:HB3	1.94	0.49
2:J:384:ASN:O	2:J:388:PHE:N	2.38	0.49
2:B:647:ILE:HD12	2:B:695:ILE:HD11	1.94	0.49
2:B:295:PHE:HB3	2:B:345:ILE:HG13	1.95	0.49
2:J:251:ALA:HB2	2:J:273:VAL:HG13	1.96	0.48
2:D:540:LEU:O	2:D:544:HIS:ND1	2.46	0.48
2:F:299:GLN:HB3	2:F:301:PRO:HD2	1.95	0.48
2:D:655:PHE:O	2:D:692:ASN:ND2	2.46	0.48
2:D:821:ALA:O	2:D:824:ARG:NH1	2.40	0.48
2:B:496:LEU:HD23	2:B:516:VAL:HG13	1.96	0.48
2:H:469:ASP:HB3	2:H:472:LYS:HB3	1.95	0.47
2:D:246:TYR:OH	2:D:268:CYS:O	2.26	0.47
2:H:411:SER:HB2	2:H:725:ILE:HD12	1.95	0.47
2:L:809:LEU:HB2	2:L:893:VAL:HG13	1.95	0.47
2:H:815:ALA:HB2	2:H:866:ARG:HH11	1.79	0.47
2:J:467:ILE:HG13	2:J:468:SER:H	1.79	0.47
2:H:275:LEU:HB3	2:H:313:TYR:CE1	2.49	0.47
2:J:485:LEU:HB3	2:J:541:TYR:CZ	2.50	0.47
2:J:744:ASN:HB3	2:J:747:VAL:HG23	1.96	0.47
2:F:843:VAL:HG13	2:F:865:ALA:HB2	1.97	0.47
2:J:299:GLN:HB3	2:J:301:PRO:HD2	1.96	0.47
2:B:640:VAL:HG12	2:B:660:ILE:HD12	1.97	0.47
2:D:699:VAL:HG12	2:D:700:ALA:H	1.79	0.47
2:B:394:LYS:NZ	2:B:435:ASP:OD1	2.48	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:637:ARG:HD2	2:B:662:THR:HB	1.96	0.46
2:J:525:MET:HB2	2:J:531:GLU:HB2	1.96	0.46
2:L:411:SER:HB2	2:L:725:ILE:HD12	1.96	0.46
2:L:432:ALA:HB1	2:L:784:GLN:HG3	1.97	0.46
2:B:469:ASP:HB3	2:B:472:LYS:HB3	1.96	0.46
2:D:344:ILE:HG22	2:D:346:LEU:HD13	1.97	0.46
2:B:299:GLN:HB3	2:B:301:PRO:HD2	1.97	0.46
2:B:568:LYS:HE2	2:B:594:LEU:HD11	1.97	0.46
2:F:647:ILE:HD12	2:F:695:ILE:HD11	1.98	0.46
1:A:11:U:N3	2:B:498:GLN:O	2.49	0.46
2:F:701:ASP:HB3	2:F:726:GLN:HE22	1.79	0.46
2:J:404:GLN:HE21	2:J:406:ILE:HD11	1.80	0.46
2:J:863:PHE:HB3	2:J:870:SER:HA	1.98	0.46
2:F:661:LEU:HB2	2:F:694:LEU:HD11	1.98	0.46
2:F:759:LYS:NZ	2:F:763:ASP:OD2	2.47	0.46
2:H:379:LYS:H	2:H:379:LYS:HG3	1.53	0.46
2:D:322:THR:OG1	2:J:320:ARG:NH2	2.49	0.46
2:H:275:LEU:HB3	2:H:313:TYR:HE1	1.80	0.46
2:L:662:THR:H	2:L:673:MET:HE3	1.81	0.46
2:B:411:SER:HB2	2:B:725:ILE:HD12	1.98	0.45
2:D:888:LYS:HZ2	2:D:891:SER:HG	1.62	0.45
2:J:858:LYS:HA	2:J:877:VAL:HG12	1.98	0.45
2:L:725:ILE:HD11	2:L:758:GLU:HG3	1.97	0.45
2:D:305:GLN:OE1	2:D:306:GLN:NE2	2.50	0.45
2:F:742:THR:HG21	2:F:747:VAL:HB	1.98	0.45
2:B:482:THR:HB	2:B:545:LEU:HD21	1.99	0.45
2:H:246:TYR:HE1	2:H:268:CYS:HB2	1.81	0.45
2:L:347:THR:HB	2:L:350:ILE:HD13	1.98	0.45
2:B:637:ARG:NH2	2:B:671:THR:HB	2.32	0.45
2:F:809:LEU:HB2	2:F:893:VAL:HG13	1.99	0.45
2:L:570:PHE:O	2:L:574:VAL:HG23	2.17	0.45
2:D:845:ARG:HB2	2:D:863:PHE:HE1	1.82	0.44
2:H:632:LEU:HD11	2:H:715:TYR:HB2	2.00	0.44
2:B:460:PHE:HB3	2:B:748:ILE:HD12	1.99	0.44
2:B:682:LEU:HD22	2:B:706:ILE:HG21	1.99	0.44
2:H:394:LYS:NZ	2:H:435:ASP:OD1	2.38	0.44
2:J:640:VAL:HG12	2:J:660:ILE:HD12	1.99	0.44
2:J:809:LEU:HB2	2:J:893:VAL:HG13	1.99	0.44
2:B:719:GLY:O	2:B:754:ASN:ND2	2.48	0.44
2:F:320:ARG:NH1	2:F:341:ASN:OD1	2.50	0.44
2:J:378:SER:O	2:J:384:ASN:ND2	2.47	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:J:618:LEU:HD13	2:J:656:LEU:HD11	2.00	0.44
2:D:768:LEU:HD21	2:D:776:PHE:CZ	2.52	0.44
2:H:699:VAL:HG12	2:H:700:ALA:H	1.83	0.44
2:F:355:LEU:HD11	2:F:363:LEU:HD21	2.00	0.44
2:F:395:LEU:HD11	2:F:784:GLN:HB3	1.99	0.44
2:F:644:LYS:HD2	2:F:660:ILE:HD11	1.98	0.44
2:B:348:PRO:HB2	2:B:382:PRO:HB2	1.98	0.44
2:D:559:ARG:HD3	2:D:646:TRP:CD1	2.52	0.44
2:B:843:VAL:HG13	2:B:865:ALA:HB2	2.00	0.44
2:B:866:ARG:HG2	2:B:867:GLN:HG2	2.00	0.44
2:L:493:LEU:HA	2:L:496:LEU:HD13	1.99	0.44
2:L:632:LEU:HD11	2:L:715:TYR:HB2	2.00	0.44
2:F:347:THR:HB	2:F:350:ILE:HD13	2.00	0.43
2:F:559:ARG:HD3	2:F:646:TRP:CD1	2.51	0.43
2:J:270:LYS:HZ3	2:J:410:ALA:HB2	1.83	0.43
2:J:699:VAL:HG12	2:J:700:ALA:H	1.83	0.43
2:J:858:LYS:HG2	2:J:875:ILE:HD12	2.00	0.43
2:B:320:ARG:HD2	2:F:332:VAL:HG11	1.99	0.43
2:B:564:LEU:HG	2:B:568:LYS:HE3	2.00	0.43
2:D:456:PRO:HD2	2:D:728:ARG:HD3	2.00	0.43
2:D:888:LYS:NZ	2:D:891:SER:OG	2.39	0.43
2:F:295:PHE:HB3	2:F:345:ILE:HG13	2.01	0.43
2:F:374:CYS:HB2	2:F:434:LEU:HD11	1.99	0.43
2:F:557:HIS:HB3	2:F:639:LEU:HD21	2.00	0.43
2:H:856:PHE:HD2	2:H:877:VAL:HG21	1.84	0.43
2:L:773:GLU:OE2	2:L:777:ARG:NH1	2.50	0.43
2:F:640:VAL:HG12	2:F:660:ILE:HD12	2.00	0.43
2:H:640:VAL:HG12	2:H:660:ILE:HD12	2.00	0.43
2:L:564:LEU:HD11	2:L:597:LEU:HD13	1.99	0.43
2:D:810:CYS:HB2	2:D:873:TRP:HZ2	1.84	0.43
2:F:559:ARG:NH2	2:F:562:ASP:OD1	2.39	0.43
2:H:773:GLU:OE2	2:H:777:ARG:NH1	2.48	0.43
2:H:286:PRO:HD2	2:H:289:GLN:HB2	2.00	0.43
2:H:656:LEU:HA	2:H:692:ASN:HD21	1.83	0.43
2:L:295:PHE:HB3	2:L:345:ILE:HG13	2.01	0.43
2:L:493:LEU:HD12	2:L:496:LEU:HD22	2.01	0.43
2:L:520:CYS:HA	2:L:523:PHE:HD1	1.84	0.43
2:B:466:ARG:NH1	2:B:555:SER:O	2.52	0.43
2:H:295:PHE:HB3	2:H:345:ILE:HG13	2.00	0.43
2:B:856:PHE:HD2	2:B:877:VAL:HG21	1.84	0.42
2:L:249:GLU:OE1	2:L:445:ASN:ND2	2.51	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:634:VAL:HG12	2:H:715:TYR:HB3	2.01	0.42
2:D:350:ILE:O	2:D:354:ASN:ND2	2.38	0.42
2:J:598:GLU:HA	2:J:601:SER:HB2	2.01	0.42
2:L:603:ASP:HB3	2:L:606:ASN:ND2	2.34	0.42
2:B:346:LEU:HD23	2:B:351:LEU:HB2	2.01	0.42
2:F:888:LYS:HA	2:F:908:TRP:HE1	1.84	0.42
1:K:18:U:H2'	1:K:19:U:C6	2.54	0.42
2:D:887:ILE:HB	2:D:892:PHE:HE2	1.85	0.42
2:L:262:ILE:HG12	2:L:440:ALA:HB3	2.01	0.42
2:B:483:GLU:OE1	2:B:502:ARG:NH2	2.48	0.42
2:F:639:LEU:HA	2:F:642:ALA:HB3	2.01	0.42
1:G:10:U:O2'	1:G:11:U:H2'	2.19	0.42
2:J:390:TYR:CE2	2:J:402:LEU:HB3	2.55	0.42
2:J:856:PHE:HD2	2:J:877:VAL:HG21	1.84	0.42
2:L:580:ASP:O	2:L:584:GLN:HG2	2.20	0.42
2:L:457:GLN:NE2	2:L:458:LYS:O	2.53	0.42
2:L:460:PHE:HB3	2:L:748:ILE:HD12	2.02	0.42
2:D:482:THR:HB	2:D:545:LEU:HD21	2.02	0.42
2:L:826:ILE:HG22	2:L:827:GLU:HG3	2.00	0.42
2:B:632:LEU:HD11	2:B:715:TYR:HB2	2.01	0.42
2:F:482:THR:HB	2:F:545:LEU:HD21	2.01	0.42
2:J:494:GLU:OE1	2:J:494:GLU:N	2.52	0.42
2:D:728:ARG:NH2	2:D:758:GLU:OE1	2.50	0.41
2:H:281:HIS:O	2:H:284:LYS:HG2	2.20	0.41
2:J:270:LYS:NZ	2:J:702:GLU:OE2	2.43	0.41
2:L:767:ARG:O	2:L:770:THR:OG1	2.36	0.41
2:F:479:MET:HG3	2:F:545:LEU:HD22	2.02	0.41
2:F:530:GLU:OE2	2:F:533:ARG:NE	2.53	0.41
2:F:644:LYS:HB2	2:F:660:ILE:HD11	2.02	0.41
2:H:246:TYR:CE1	2:H:268:CYS:HB2	2.56	0.41
2:D:292:LYS:NZ	2:D:341:ASN:O	2.53	0.41
2:D:502:ARG:HD2	2:D:509:TYR:CD1	2.56	0.41
2:L:270:LYS:NZ	2:L:702:GLU:OE2	2.36	0.41
2:D:356:LYS:HE2	2:D:356:LYS:HB3	1.84	0.41
2:H:442:VAL:HG21	2:H:449:LEU:HD22	2.03	0.41
2:H:580:ASP:OD1	2:H:581:GLU:N	2.51	0.41
2:H:807:LYS:HE2	2:H:816:LEU:HD13	2.03	0.41
2:L:853:PHE:CZ	2:L:856:PHE:HB2	2.56	0.41
2:B:513:ILE:HD12	2:B:513:ILE:HA	1.91	0.41
2:F:812:LYS:HB3	2:F:869:CYS:SG	2.61	0.41
1:G:3:U:H2'	1:G:4:A:C8	2.56	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:685:PHE:CE1	2:H:694:LEU:HB2	2.56	0.41
2:J:559:ARG:HD3	2:J:646:TRP:HD1	1.86	0.41
2:L:467:ILE:HG13	2:L:468:SER:H	1.86	0.41
2:L:768:LEU:HD21	2:L:776:PHE:CZ	2.56	0.41
2:L:845:ARG:O	2:L:861:LYS:N	2.49	0.41
2:B:525:MET:SD	2:B:534:ILE:HD12	2.60	0.41
2:B:878:LYS:HA	2:B:883:GLU:HA	2.02	0.41
2:D:395:LEU:HD13	2:D:785:THR:HA	2.03	0.41
2:F:596:GLU:O	2:F:600:VAL:HG23	2.21	0.41
2:F:640:VAL:HG21	2:F:697:THR:HG23	2.01	0.41
2:L:520:CYS:HA	2:L:523:PHE:CD1	2.56	0.41
1:A:13:G:O2'	1:A:14:G:OP2	2.39	0.40
2:J:811:ARG:CZ	2:J:904:LEU:HG	2.51	0.40
2:J:861:LYS:HE3	2:J:863:PHE:CZ	2.56	0.40
2:D:374:CYS:HB2	2:D:434:LEU:HD11	2.03	0.40
2:J:557:HIS:O	2:J:715:TYR:OH	2.26	0.40
2:H:859:ARG:HH12	2:H:878:LYS:HB2	1.87	0.40
2:L:347:THR:HG22	2:L:348:PRO:HD2	2.03	0.40
2:B:867:GLN:HE22	2:D:575:ARG:HG3	1.86	0.40
2:F:632:LEU:HD21	2:F:643:LEU:HD13	2.04	0.40
2:J:283:LYS:HD3	2:J:283:LYS:HA	1.85	0.40
2:D:683:ASP:O	2:D:686:LYS:HD3	2.21	0.40
2:H:723:LYS:HE3	2:H:723:LYS:HB2	1.89	0.40

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	
2	B	666/684 (97%)	644 (97%)	22 (3%)	0	100	100
2	D	666/684 (97%)	643 (96%)	23 (4%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	F	666/684 (97%)	646 (97%)	20 (3%)	0	100	100
2	H	666/684 (97%)	644 (97%)	22 (3%)	0	100	100
2	J	666/684 (97%)	647 (97%)	19 (3%)	0	100	100
2	L	666/684 (97%)	645 (97%)	21 (3%)	0	100	100
All	All	3996/4104 (97%)	3869 (97%)	127 (3%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains ⓘ

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	
2	B	575/611 (94%)	571 (99%)	4 (1%)	84	94
2	D	576/611 (94%)	570 (99%)	6 (1%)	76	91
2	F	574/611 (94%)	571 (100%)	3 (0%)	88	96
2	H	584/611 (96%)	580 (99%)	4 (1%)	84	94
2	J	580/611 (95%)	577 (100%)	3 (0%)	88	96
2	L	579/611 (95%)	576 (100%)	3 (0%)	88	96
All	All	3468/3666 (95%)	3445 (99%)	23 (1%)	84	94

All (23) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	B	272	PHE
2	B	419	ASN
2	B	469	ASP
2	B	866	ARG
2	D	272	PHE
2	D	317	HIS
2	D	384	ASN
2	D	492	ASP
2	D	692	ASN

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Mol	Chain	Res	Type
2	D	879	TYR
2	F	272	PHE
2	F	705	ASP
2	F	741	LEU
2	H	272	PHE
2	H	317	HIS
2	H	523	PHE
2	H	776	PHE
2	J	523	PHE
2	J	557	HIS
2	J	782	HIS
2	L	317	HIS
2	L	836	ASP
2	L	879	TYR

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

Mol	Chain	Res	Type
2	D	305	GLN
2	D	306	GLN
2	F	782	HIS
2	H	524	GLN
2	L	606	ASN

5.3.3 RNA ⓘ

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
1	A	21/22 (95%)	4 (19%)	0
1	C	21/22 (95%)	3 (14%)	0
1	E	21/22 (95%)	3 (14%)	0
1	G	21/22 (95%)	3 (14%)	0
1	I	21/22 (95%)	4 (19%)	0
1	K	21/22 (95%)	4 (19%)	0
All	All	126/132 (95%)	21 (16%)	0

All (21) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
1	A	11	U
1	A	12	C

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Mol	Chain	Res	Type
1	A	13	G
1	A	14	G
1	C	12	C
1	C	13	G
1	C	14	G
1	E	12	C
1	E	13	G
1	E	14	G
1	G	12	C
1	G	13	G
1	G	14	G
1	I	12	C
1	I	13	G
1	I	14	G
1	I	20	G
1	K	11	U
1	K	12	C
1	K	13	G
1	K	14	G

There are no RNA pucker outliers to report.

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 6 ligands modelled in this entry, 6 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 [i](#)

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, 95th 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	22/22 (100%)	-0.09	0 100 100	44, 51, 95, 96	0
1	C	22/22 (100%)	-0.16	0 100 100	41, 52, 93, 96	0
1	E	22/22 (100%)	-0.06	0 100 100	37, 51, 108, 111	0
1	G	22/22 (100%)	-0.20	0 100 100	41, 51, 95, 101	0
1	I	22/22 (100%)	0.10	1 (4%) 33 12	47, 57, 110, 116	0
1	K	22/22 (100%)	0.19	2 (9%) 9 3	48, 59, 105, 113	0
2	B	672/684 (98%)	-0.33	3 (0%) 92 79	29, 56, 81, 101	0
2	D	672/684 (98%)	-0.23	3 (0%) 92 79	37, 61, 90, 103	0
2	F	672/684 (98%)	-0.24	5 (0%) 87 69	37, 59, 89, 100	0
2	H	672/684 (98%)	-0.26	2 (0%) 94 84	41, 61, 84, 102	0
2	J	672/684 (98%)	-0.22	6 (0%) 84 63	36, 64, 94, 112	0
2	L	672/684 (98%)	-0.21	3 (0%) 92 79	36, 64, 92, 110	0
All	All	4164/4236 (98%)	-0.24	25 (0%) 89 72	29, 60, 90, 116	0

All (25) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	J	576	ALA	5.5
2	H	287	GLN	3.7
2	F	397	GLY	3.4
1	K	11	U	3.0
2	L	867	GLN	2.9
2	D	492	ASP	2.9
2	F	492	ASP	2.9
2	J	501	ASN	2.8
2	L	499	ILE	2.6
2	L	527	ASP	2.5
2	J	688	SER	2.5

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Mol	Chain	Res	Type	RSRZ
1	K	12	C	2.5
2	J	867	GLN	2.4
2	B	855	SER	2.4
2	F	867	GLN	2.3
1	I	11	U	2.3
2	F	287	GLN	2.3
2	D	501	ASN	2.3
2	J	397	GLY	2.2
2	J	879	TYR	2.2
2	F	416	ASP	2.1
2	B	881	THR	2.1
2	H	397	GLY	2.1
2	B	690	ASP	2.0
2	D	468	SER	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.

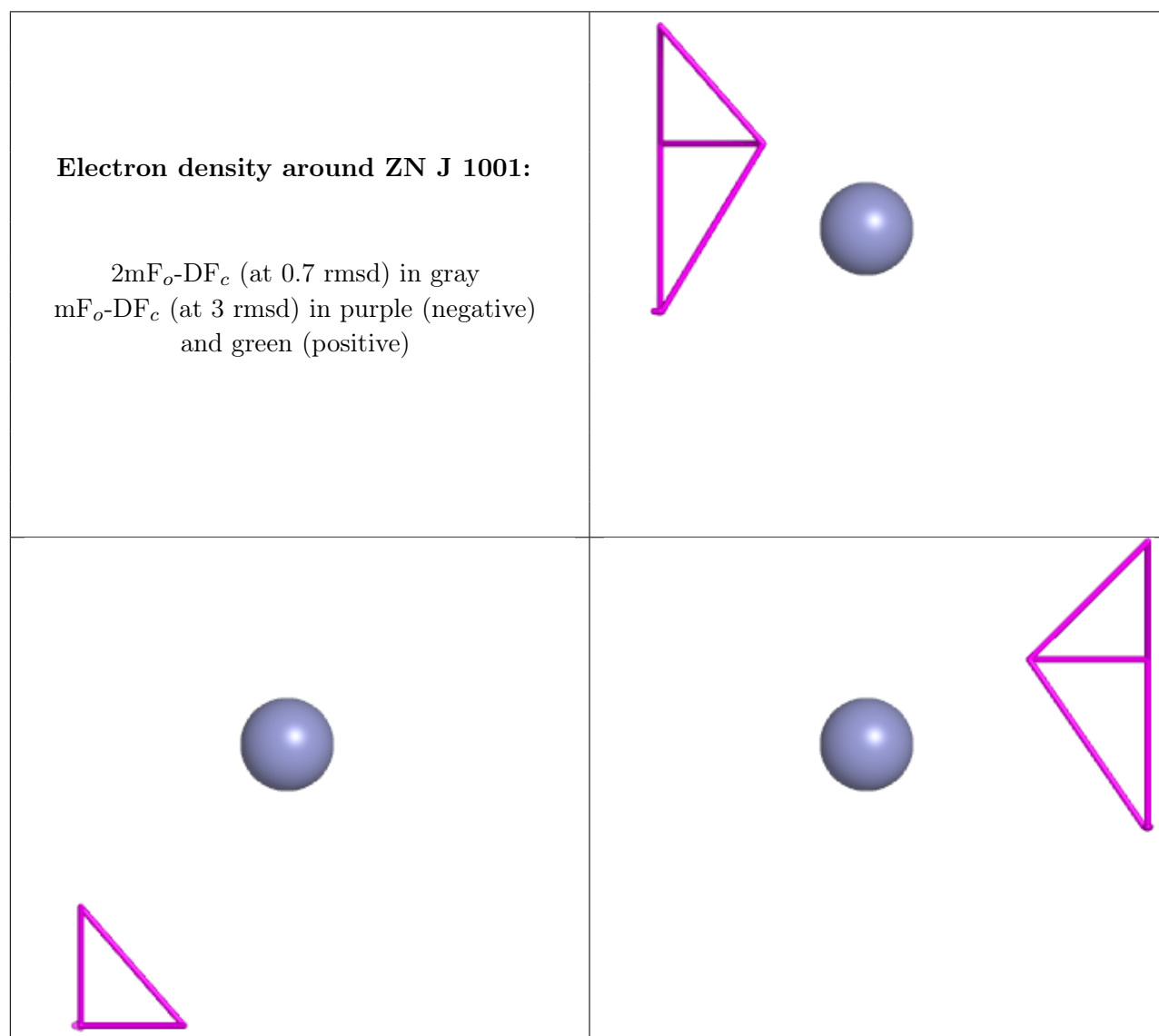
6.4 Ligands [i](#)

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, 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(\AA^2)	Q<0.9
3	ZN	J	1001	1/1	0.94	0.11	100,100,100,100	0
3	ZN	L	1001	1/1	0.95	0.14	89,89,89,89	0
3	ZN	H	1001	1/1	0.97	0.10	65,65,65,65	0
3	ZN	D	1001	1/1	0.97	0.15	76,76,76,76	0
3	ZN	F	1001	1/1	0.97	0.15	77,77,77,77	0
3	ZN	B	1001	1/1	0.99	0.15	62,62,62,62	0

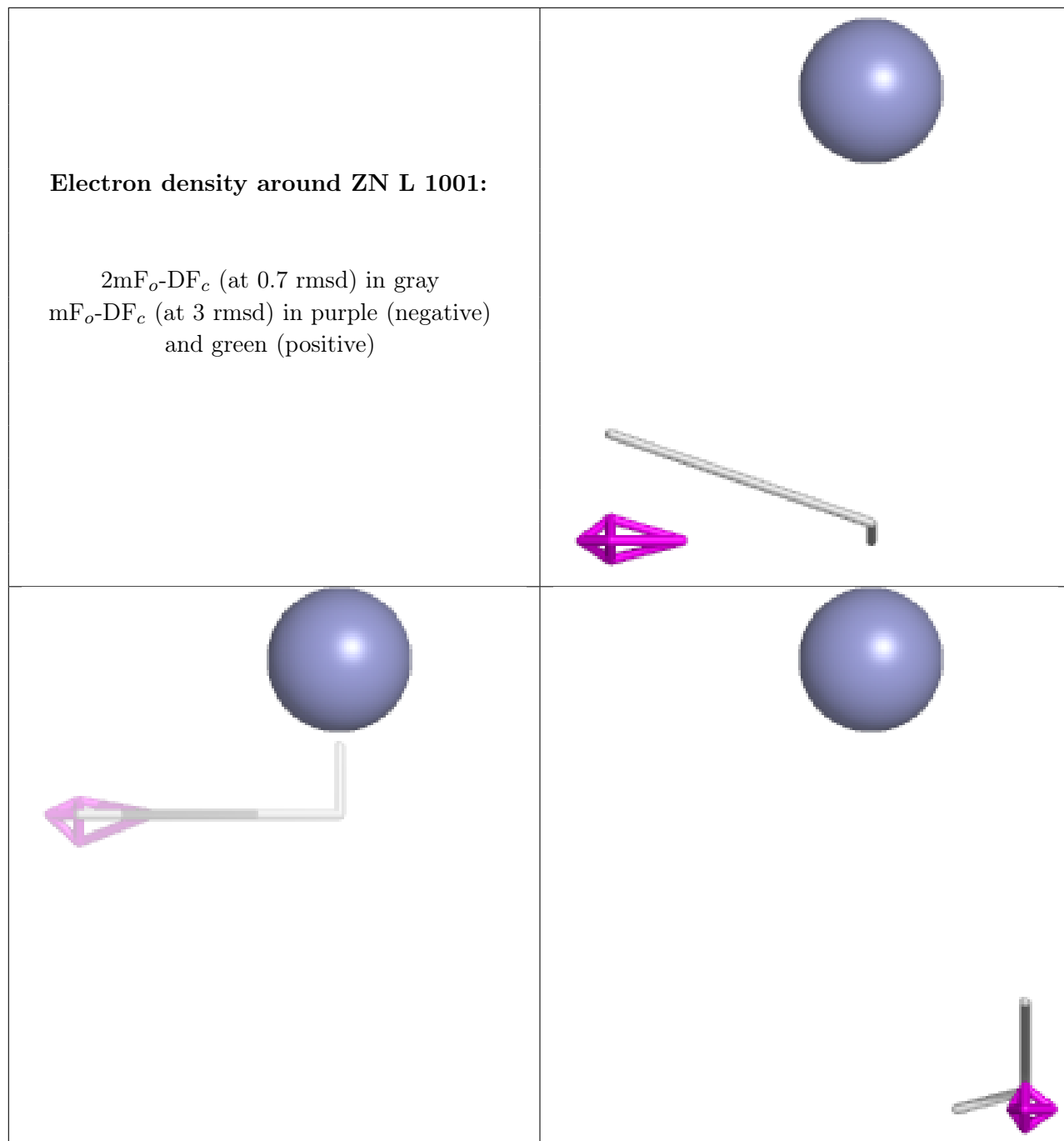
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers

as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.



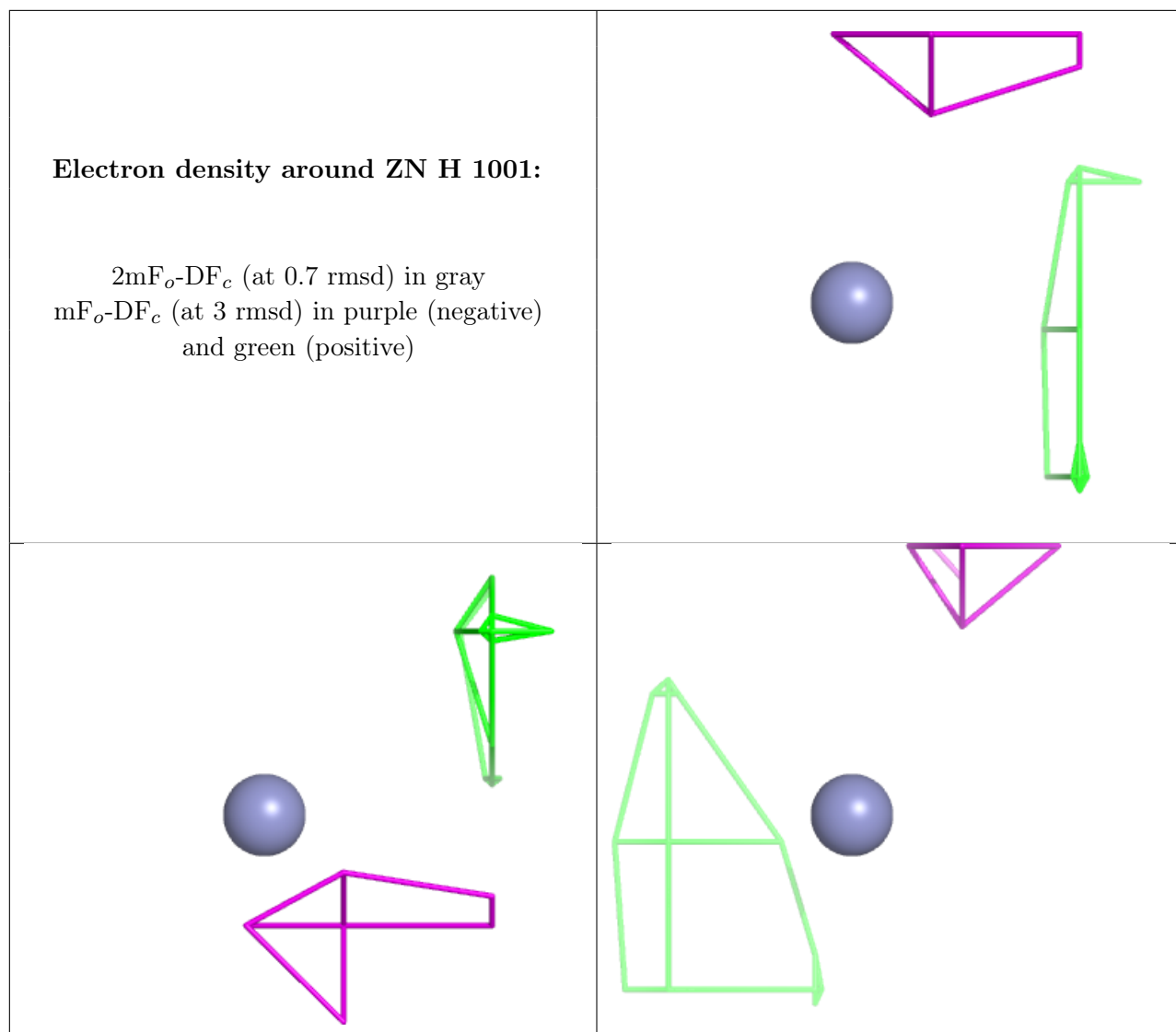
Electron density around ZN L 1001:

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



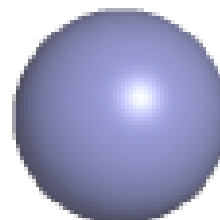
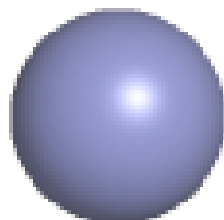
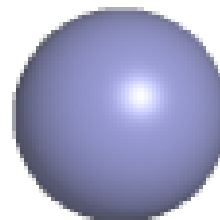
Electron density around ZN H 1001:

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



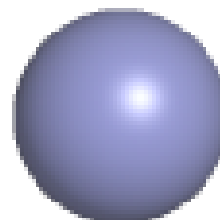
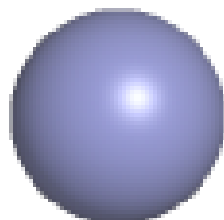
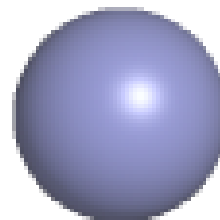
Electron density around ZN D 1001:

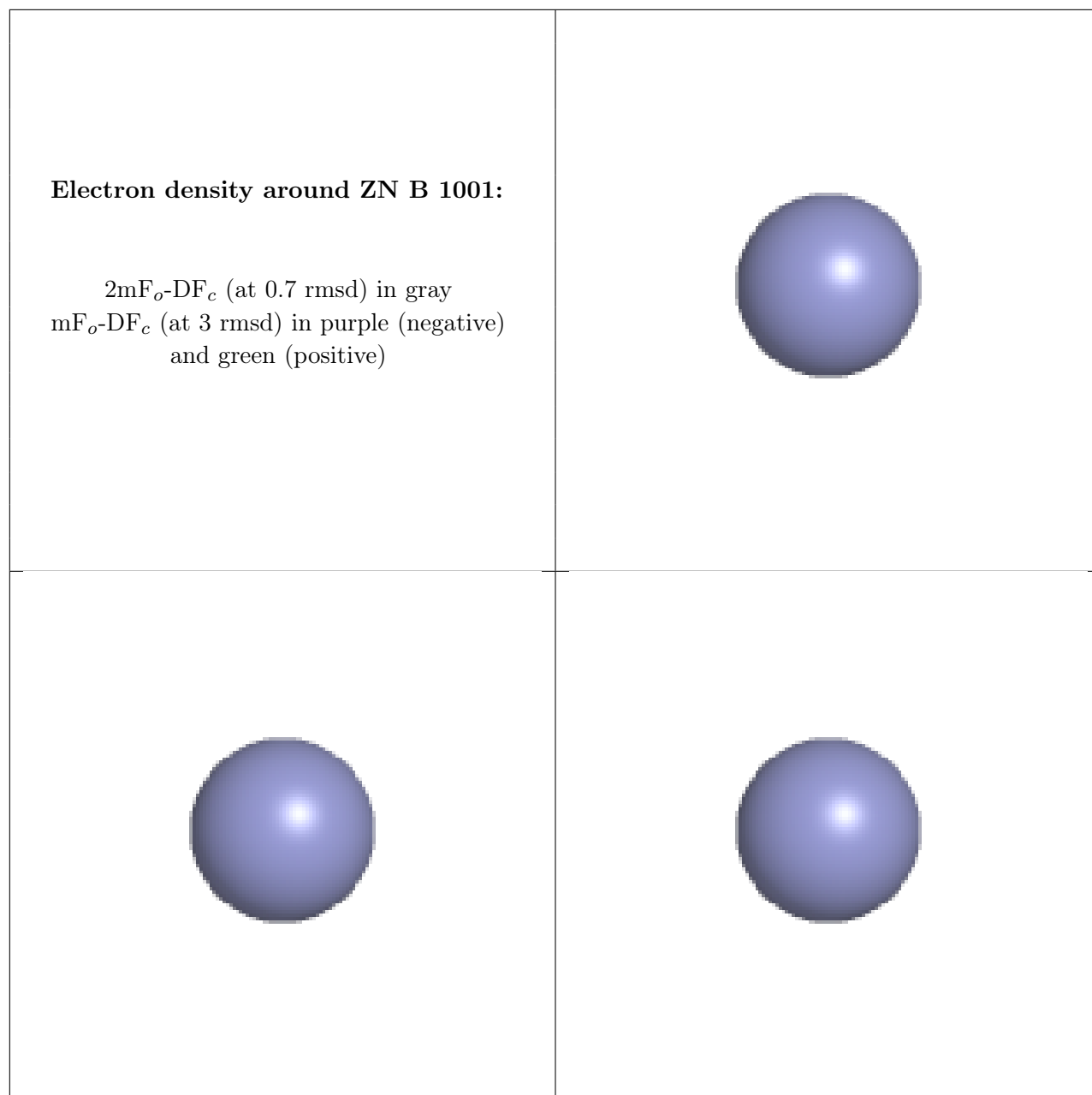
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around ZN F 1001:

$2mF_o-DF_c$ (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.