



# Full wwPDB X-ray Structure Validation Report ⓘ

Oct 25, 2017 – 11:24 PM EDT

PDB ID : 3KQH  
Title : Three Conformational Snapshots of the Hepatitis C Virus NS3 Helicase Reveal a Ratchet Translocation Mechanism  
Authors : Gu, M.; Rice, C.M.  
Deposited on : unknown  
Resolution : 2.40 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](mailto:validation@mail.wwpdb.org)

A user guide is available at

<http://wwpdb.org/validation/2016/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

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The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20030345
Percentile statistics	:	20161228.v01 (using entries in the PDB archive December 28th 2016)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	rb-20030345

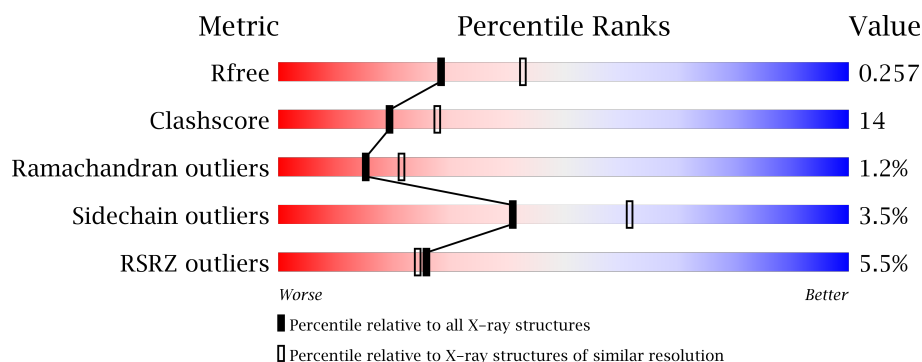
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.40 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))
$R_{free}$	100719	3166 (2.40-2.40)
Clashscore	112137	3674 (2.40-2.40)
Ramachandran outliers	110173	3616 (2.40-2.40)
Sidechain outliers	110143	3617 (2.40-2.40)
RSRZ outliers	101464	3195 (2.40-2.40)

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 on the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria. 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	436	<div> <div>5%</div> <div>70%</div> <div>29%</div> <div>.</div> </div>
1	B	436	<div> <div>6%</div> <div>68%</div> <div>30%</div> <div>.</div> </div>
2	C	6	<div> <div>67%</div> <div>33%</div> </div>
2	D	6	<div> <div>17%</div> <div>33%</div> <div>67%</div> </div>

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 6934 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 Serine protease/NTPase/helicase NS3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	436	Total	C	N	O	S	0	0	0
			3275	2074	555	625	21			
1	B	436	Total	C	N	O	S	0	0	0
			3275	2074	555	625	21			

- Molecule 2 is a DNA chain called 5'-D(\*AP\*AP\*AP\*AP\*AP\*A)-3'.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	6	Total	C	N	O	P	0	0	0
			123	60	30	28	5			
2	D	6	Total	C	N	O	P	0	0	0
			123	60	30	28	5			

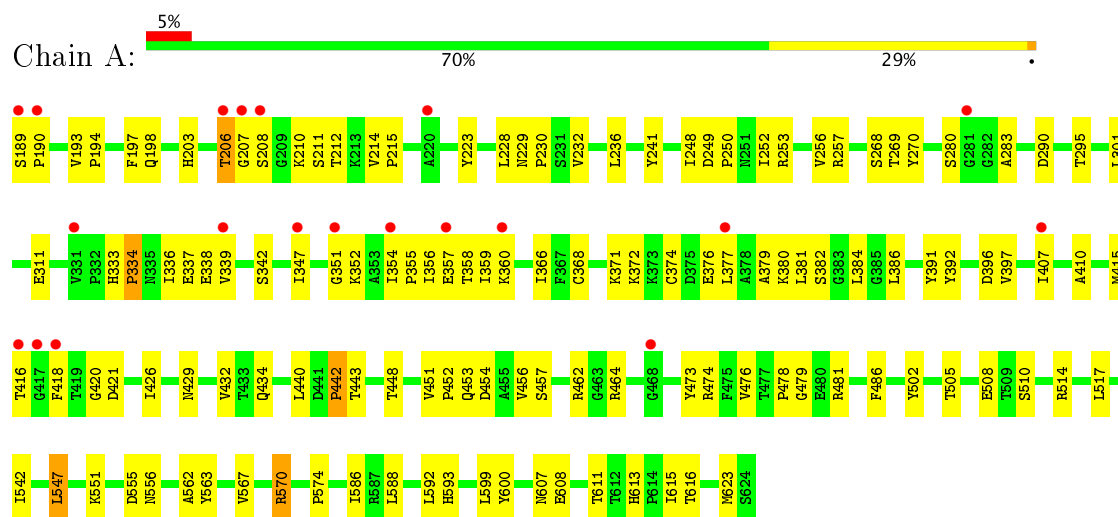
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	72	Total	O	0	0
			72	72		
3	B	60	Total	O	0	0
			60	60		
3	C	5	Total	O	0	0
			5	5		
3	D	1	Total	O	0	0
			1	1		

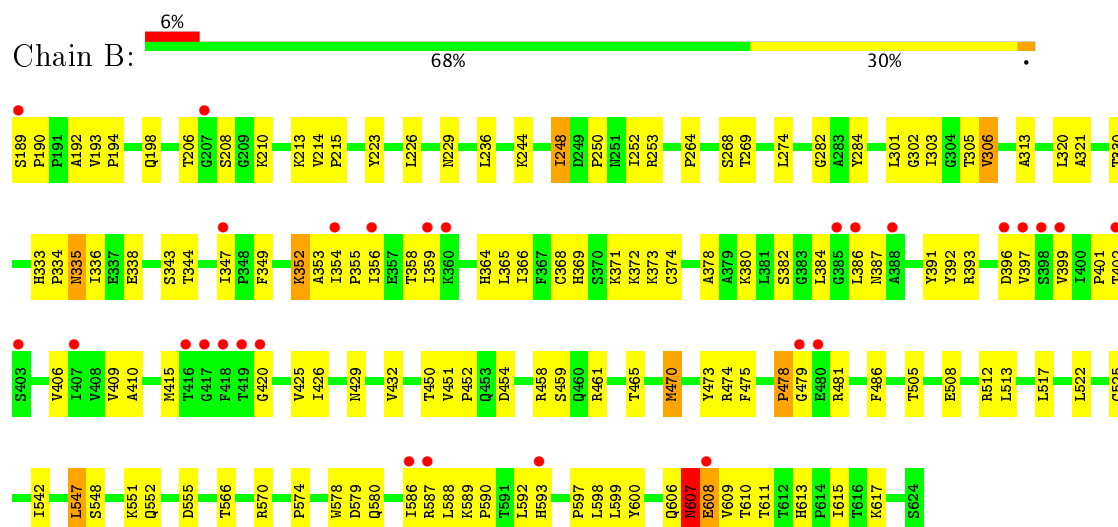
### 3 Residue-property plots

These plots are drawn for all protein, RNA and DNA 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: Serine protease/NTPase/helicase NS3



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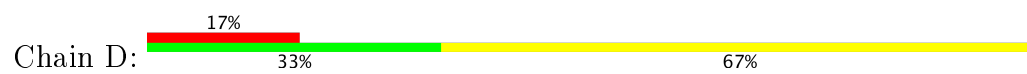


- Molecule 2: 5'-D(\*AP\*AP\*AP\*AP\*AP\*A)-3'





- Molecule 2: 5'-D(\*AP\*AP\*AP\*AP\*AP\*A)-3'



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	45.20Å 115.38Å 197.31Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 2.40 37.49 – 2.39	Depositor EDS
% Data completeness (in resolution range)	93.5 (50.00-2.40) 92.7 (37.49-2.39)	Depositor EDS
$R_{merge}$	0.06	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.16 (at 2.39Å)	Xtriage
Refinement program	CNS	Depositor
R, $R_{free}$	0.226 , 0.258 0.225 , 0.257	Depositor DCC
$R_{free}$ test set	1917 reflections (4.95%)	DCC
Wilson B-factor (Å <sup>2</sup> )	38.1	Xtriage
Anisotropy	0.347	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.33 , 47.0	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.93	EDS
Total number of atoms	6934	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	47.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.14% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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](#)

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	0.40	0/3356	0.57	0/4585
1	B	0.39	0/3356	0.57	0/4585
2	C	0.47	0/140	0.79	0/214
2	D	0.43	0/140	0.72	0/214
All	All	0.40	0/6992	0.58	0/9598

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	563	TYR	Sidechain

### 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	3275	0	3236	92	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	B	3275	0	3236	103	0
2	C	123	0	68	2	0
2	D	123	0	68	5	0
3	A	72	0	0	7	0
3	B	60	0	0	7	0
3	C	5	0	0	1	0
3	D	1	0	0	0	0
All	All	6934	0	6608	193	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 14.

All (193) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:611:THR:HB	3:A:49:HOH:O	1.55	1.04
1:B:366:ILE:HG23	1:B:426:ILE:HB	1.59	0.83
1:B:607:ASN:ND2	1:B:608:GLU:H	1.77	0.82
1:A:397:VAL:HG11	1:A:418:PHE:HB2	1.63	0.81
1:B:555:ASP:HB2	3:B:45:HOH:O	1.84	0.78
1:B:366:ILE:HD12	1:B:426:ILE:HD12	1.67	0.76
1:B:415:MET:HB2	3:B:43:HOH:O	1.87	0.74
1:A:208:SER:HB2	1:B:470:MET:SD	2.28	0.73
1:A:359:ILE:HD12	1:A:407:ILE:HD13	1.72	0.71
1:A:194:PRO:HG3	1:A:198:GLN:HB2	1.71	0.71
1:B:320:LEU:HD13	1:B:522:LEU:HD21	1.73	0.71
1:A:379:ALA:HA	3:A:76:HOH:O	1.92	0.70
1:A:611:THR:HG23	1:A:616:THR:HG21	1.72	0.70
1:B:505:THR:OG1	1:B:508:GLU:HG3	1.92	0.69
1:A:256:VAL:HG23	1:A:257:ARG:H	1.58	0.67
1:A:611:THR:HG22	1:A:611:THR:O	1.95	0.67
1:A:357:GLU:HA	1:A:360:LYS:HB2	1.76	0.66
1:A:547:LEU:O	1:A:551:LYS:HG3	1.96	0.65
1:B:338:GLU:HG2	1:B:473:TYR:HD2	1.61	0.65
1:A:207:GLY:HA2	1:A:210:LYS:HE3	1.78	0.64
1:B:210:LYS:HD2	1:B:321:ALA:HB1	1.80	0.63
1:A:236:LEU:HD23	1:A:252:ILE:HG21	1.83	0.60
1:A:253:ARG:HB2	1:A:268:SER:HB2	1.84	0.59
1:B:274:LEU:HD21	1:B:302:GLY:HA2	1.83	0.59
1:A:229:ASN:O	1:A:269:THR:HA	2.03	0.58
1:B:349:PHE:HB3	3:B:39:HOH:O	2.03	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:371:LYS:HG3	1:B:392:TYR:CD2	2.37	0.58
1:A:451:VAL:HG23	1:A:452:PRO:HD2	1.83	0.58
1:B:486:PHE:O	1:B:525:CYS:HB2	2.04	0.58
1:B:589:LYS:HB3	1:B:590:PRO:HD3	1.85	0.57
1:A:351:GLY:C	1:A:352:LYS:HD2	2.25	0.57
1:A:366:ILE:HG21	1:A:377:LEU:HD21	1.86	0.57
1:A:371:LYS:HG3	1:A:392:TYR:CD2	2.40	0.57
1:B:401:PRO:HG2	1:B:406:VAL:HG11	1.87	0.56
1:B:548:SER:O	1:B:552:GLN:HG3	2.05	0.56
1:A:391:TYR:HB3	1:A:410:ALA:HB2	1.88	0.56
1:B:356:ILE:HD11	1:B:386:LEU:HD11	1.86	0.56
1:B:248:ILE:HD11	1:B:264:PRO:HD2	1.87	0.55
1:B:282:GLY:HA2	1:B:313:ALA:O	2.06	0.55
1:B:305:THR:OG1	1:B:512:ARG:HD3	2.06	0.55
1:B:574:PRO:HG2	1:B:607:ASN:OD1	2.05	0.55
1:B:586:ILE:HG13	1:B:587:ARG:N	2.22	0.55
1:B:192:ALA:O	1:B:194:PRO:HD3	2.06	0.55
1:B:333:HIS:HB3	1:B:336:ILE:HB	1.89	0.55
1:B:391:TYR:HB3	1:B:410:ALA:HB2	1.87	0.55
1:B:425:VAL:HG23	1:B:465:THR:HB	1.88	0.55
1:B:393:ARG:HG2	1:B:393:ARG:HH11	1.71	0.54
1:A:429:ASN:N	1:A:429:ASN:HD22	2.06	0.54
1:B:607:ASN:ND2	1:B:608:GLU:N	2.51	0.54
1:B:599:LEU:HD22	1:B:615:ILE:HG22	1.90	0.54
1:A:339:VAL:O	1:A:474:ARG:HA	2.07	0.53
1:A:334:PRO:HA	3:B:73:HOH:O	2.08	0.53
1:A:189:SER:N	1:A:190:PRO:HD2	2.24	0.53
1:B:193:VAL:HG21	1:B:223:TYR:CE2	2.44	0.52
1:B:335:ASN:ND2	1:B:335:ASN:H	2.07	0.52
1:B:229:ASN:O	1:B:269:THR:HA	2.10	0.52
1:B:335:ASN:HD22	1:B:335:ASN:H	1.58	0.52
1:B:189:SER:HB3	1:B:190:PRO:HD3	1.91	0.52
1:B:613:HIS:HE1	1:B:615:ILE:HD13	1.74	0.52
1:B:607:ASN:HD22	1:B:608:GLU:H	1.56	0.51
1:B:429:ASN:OD1	1:B:475:PHE:HB2	2.10	0.51
1:A:228:LEU:HA	1:A:268:SER:O	2.10	0.51
1:A:502:TYR:OH	2:D:6:DA:H1'	2.11	0.51
1:B:336:ILE:HD11	1:B:465:THR:O	2.11	0.51
1:A:358:THR:HG22	1:A:359:ILE:HG23	1.92	0.51
1:B:380:LYS:HB2	1:B:380:LYS:NZ	2.26	0.51
1:A:452:PRO:HG2	1:A:481:ARG:HH12	1.76	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:397:VAL:CG1	1:A:418:PHE:HB2	2.39	0.50
1:B:335:ASN:HD22	1:B:335:ASN:N	2.08	0.50
1:A:280:SER:HB2	1:A:283:ALA:HB2	1.93	0.50
1:A:366:ILE:HD12	1:A:426:ILE:HD12	1.93	0.50
1:A:333:HIS:HB3	1:A:336:ILE:HB	1.93	0.50
1:A:542:ILE:HD13	1:A:547:LEU:HG	1.94	0.50
1:A:611:THR:HG23	1:A:616:THR:CG2	2.41	0.50
1:A:434:GLN:NE2	1:A:556:ASN:HD22	2.10	0.49
1:B:248:ILE:O	1:B:250:PRO:HD3	2.11	0.49
1:A:453:GLN:HB2	1:A:457:SER:HB3	1.94	0.49
1:A:197:PHE:O	1:A:198:GLN:HG2	2.12	0.49
1:A:232:VAL:HG23	2:D:5:DA:H5"	1.95	0.49
1:B:578:TRP:CZ2	1:B:589:LYS:HG3	2.48	0.49
1:A:382:SER:HA	1:A:386:LEU:O	2.12	0.48
1:A:574:PRO:HG2	1:A:607:ASN:OD1	2.13	0.48
1:B:478:PRO:HG2	1:B:479:GLY:H	1.77	0.48
1:A:451:VAL:CG2	1:A:452:PRO:HD2	2.43	0.48
1:A:505:THR:OG1	1:A:508:GLU:HG3	2.12	0.48
1:A:567:VAL:HG21	1:A:599:LEU:HD11	1.93	0.48
1:B:451:VAL:HG23	1:B:452:PRO:HD2	1.93	0.48
1:B:542:ILE:HD13	1:B:547:LEU:HG	1.94	0.48
1:B:599:LEU:HD22	1:B:615:ILE:CG2	2.44	0.48
1:A:593:HIS:HB3	3:A:10:HOH:O	2.13	0.48
1:B:372:LYS:HG3	1:B:373:LYS:N	2.27	0.48
1:A:368:CYS:HB2	1:A:374:CYS:SG	2.54	0.48
1:A:380:LYS:HE2	1:A:384:LEU:HD11	1.94	0.48
1:B:365:LEU:C	1:B:366:ILE:HD13	2.34	0.48
1:A:440:LEU:HD23	1:A:623:MET:HE2	1.95	0.48
1:A:442:PRO:HG2	1:A:443:THR:HG23	1.96	0.48
1:B:349:PHE:CG	1:B:354:ILE:HD11	2.49	0.48
1:B:358:THR:HG23	1:B:474:ARG:CZ	2.44	0.48
1:A:337:GLU:OE2	1:B:213:LYS:HG2	2.14	0.48
1:B:330:THR:HG21	1:B:458:ARG:HB3	1.96	0.48
1:A:462:ARG:HG3	1:A:473:TYR:CG	2.49	0.47
1:A:381:LEU:O	1:A:386:LEU:HB2	2.14	0.47
1:B:566:THR:O	1:B:570:ARG:HD3	2.15	0.47
1:A:358:THR:HG23	1:A:474:ARG:NE	2.30	0.47
1:A:397:VAL:HG21	1:A:416:THR:O	2.15	0.47
1:B:372:LYS:HE3	3:C:37:HOH:O	2.15	0.47
1:B:210:LYS:HA	1:B:214:VAL:CG2	2.45	0.47
1:B:359:ILE:HA	1:B:364:HIS:CD2	2.49	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:248:ILE:HG23	1:B:248:ILE:O	2.15	0.46
1:B:303:ILE:O	1:B:306:VAL:HG13	2.14	0.46
1:B:391:TYR:OH	1:B:397:VAL:HG22	2.15	0.46
1:B:236:LEU:HD23	1:B:252:ILE:HG21	1.96	0.46
1:A:510:SER:O	1:A:514:ARG:HB2	2.15	0.46
1:B:391:TYR:O	1:B:410:ALA:CB	2.64	0.46
1:B:461:ARG:HH11	1:B:461:ARG:HG3	1.81	0.46
1:B:333:HIS:ND1	1:B:334:PRO:HD2	2.31	0.46
1:B:508:GLU:HB3	3:B:7:HOH:O	2.15	0.46
1:A:193:VAL:HG21	1:A:223:TYR:CE2	2.51	0.45
1:B:344:THR:O	1:B:355:PRO:HG3	2.16	0.45
1:B:391:TYR:CE2	1:B:415:MET:HG2	2.52	0.45
1:A:429:ASN:N	1:A:429:ASN:ND2	2.64	0.45
1:B:588:LEU:O	1:B:592:LEU:HG	2.16	0.45
1:A:478:PRO:HG2	1:A:479:GLY:H	1.81	0.45
1:B:371:LYS:HG3	1:B:392:TYR:CG	2.51	0.45
1:B:356:ILE:CD1	1:B:386:LEU:HD11	2.46	0.45
1:B:429:ASN:HD22	1:B:429:ASN:N	2.14	0.45
1:A:352:LYS:HD3	1:A:476:VAL:HG13	1.99	0.45
1:B:451:VAL:CG2	1:B:452:PRO:HD2	2.47	0.45
1:B:368:CYS:HB2	1:B:374:CYS:SG	2.57	0.45
1:B:380:LYS:O	1:B:384:LEU:HB2	2.17	0.44
1:B:597:PRO:HA	1:B:610:THR:HG23	2.00	0.44
1:A:230:PRO:HD3	1:A:270:TYR:HE2	1.83	0.44
1:B:432:VAL:HG21	2:C:2:DA:H5"	2.00	0.44
1:A:376:GLU:O	1:A:379:ALA:HB3	2.17	0.44
1:A:206:THR:HA	3:A:50:HOH:O	2.18	0.44
1:A:542:ILE:HD11	1:A:562:ALA:HB3	2.00	0.44
1:B:391:TYR:O	1:B:410:ALA:HB1	2.18	0.44
1:B:194:PRO:HG2	1:B:198:GLN:HB3	1.99	0.44
1:A:454:ASP:OD1	1:A:456:VAL:N	2.51	0.43
1:A:355:PRO:O	1:A:358:THR:HB	2.18	0.43
1:A:241:TYR:CE2	1:B:470:MET:HB3	2.54	0.43
1:B:611:THR:O	1:B:617:LYS:HE2	2.18	0.43
1:A:371:LYS:HG3	1:A:392:TYR:CG	2.54	0.43
1:A:613:HIS:CE1	1:A:615:ILE:HG12	2.54	0.43
1:A:338:GLU:HG2	1:A:473:TYR:HD2	1.84	0.43
1:B:547:LEU:O	1:B:551:LYS:HG3	2.19	0.43
1:B:598:LEU:HB2	1:B:609:VAL:HG13	2.01	0.43
1:A:248:ILE:O	1:A:250:PRO:HD3	2.19	0.42
1:B:354:ILE:N	1:B:354:ILE:HD12	2.35	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:396:ASP:O	1:B:399:VAL:HG22	2.19	0.42
1:A:214:VAL:N	1:A:215:PRO:HD2	2.34	0.42
1:A:211:SER:HB3	1:A:290:ASP:OD1	2.19	0.42
1:A:347:ILE:HG21	1:A:381:LEU:HD21	2.01	0.42
1:A:434:GLN:HG3	2:D:2:DA:H62	1.84	0.42
1:A:588:LEU:O	1:A:592:LEU:HG	2.18	0.42
1:B:214:VAL:HB	1:B:215:PRO:CD	2.49	0.42
1:B:302:GLY:O	1:B:306:VAL:HG12	2.19	0.42
1:A:382:SER:HB3	3:A:76:HOH:O	2.19	0.42
1:B:226:LEU:HD22	1:B:284:TYR:CE2	2.53	0.42
1:A:212:THR:O	1:A:215:PRO:HG2	2.19	0.42
1:B:374:CYS:HA	1:B:409:VAL:CG1	2.49	0.42
1:B:454:ASP:HB2	1:B:481:ARG:O	2.19	0.42
1:B:429:ASN:N	1:B:429:ASN:ND2	2.67	0.42
1:B:369:HIS:HD2	1:B:450:THR:CG2	2.32	0.42
1:B:579:ASP:CG	1:B:580:GLN:H	2.23	0.42
1:A:432:VAL:HG21	2:D:2:DA:H5'	2.01	0.42
1:A:354:ILE:HD13	1:A:426:ILE:HD13	2.02	0.42
1:A:342:SER:HB2	3:A:123:HOH:O	2.18	0.42
1:A:380:LYS:O	1:A:384:LEU:HD13	2.19	0.42
1:B:593:HIS:HB2	3:B:27:HOH:O	2.19	0.41
1:A:415:MET:CE	1:A:464:ARG:HH12	2.33	0.41
1:A:256:VAL:HG23	1:A:257:ARG:N	2.29	0.41
1:A:354:ILE:HD11	1:A:426:ILE:HG21	2.02	0.41
1:B:450:THR:HG23	2:C:1:DA:C5	2.56	0.41
1:A:570:ARG:CG	1:A:570:ARG:HH11	2.33	0.41
1:B:330:THR:HG22	1:B:459:SER:OG	2.21	0.41
1:B:352:LYS:HB3	1:B:353:ALA:H	1.65	0.41
1:A:356:ILE:HG23	1:A:357:GLU:N	2.36	0.41
1:B:513:LEU:O	1:B:517:LEU:HD13	2.21	0.41
1:B:253:ARG:HB2	1:B:268:SER:HB2	2.03	0.41
1:B:378:ALA:O	1:B:382:SER:HB2	2.21	0.41
1:A:203:HIS:HB2	3:A:63:HOH:O	2.20	0.41
1:B:338:GLU:HA	1:B:473:TYR:O	2.21	0.41
2:D:1:DA:HO5'	2:D:1:DA:H8	1.66	0.41
1:A:358:THR:HG23	1:A:474:ARG:CZ	2.51	0.40
1:B:402:THR:N	3:B:122:HOH:O	2.53	0.40
1:A:280:SER:HB2	1:A:283:ALA:CB	2.52	0.40
1:A:613:HIS:HE1	1:A:615:ILE:HG12	1.84	0.40
1:B:349:PHE:HB2	1:B:354:ILE:HD11	2.04	0.40
1:A:197:PHE:CD2	1:A:311:GLU:HB2	2.57	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:295:THR:HB	1:A:486:PHE:HB3	2.04	0.40
1:A:586:ILE:HD12	1:A:586:ILE:HA	1.91	0.40
1:B:347:ILE:O	1:B:353:ALA:HA	2.21	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	
1	A	434/436 (100%)	398 (92%)	32 (7%)	4 (1%)	20	29
1	B	434/436 (100%)	396 (91%)	32 (7%)	6 (1%)	13	18
All	All	868/872 (100%)	794 (92%)	64 (7%)	10 (1%)	15	21

All (10) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	607	ASN
1	A	420	GLY
1	B	206	THR
1	B	208	SER
1	B	248	ILE
1	B	420	GLY
1	A	421	ASP
1	A	206	THR
1	B	478	PRO
1	A	334	PRO

### 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	
1	A	356/356 (100%)	344 (97%)	12 (3%)	42	63
1	B	356/356 (100%)	343 (96%)	13 (4%)	39	59
All	All	712/712 (100%)	687 (96%)	25 (4%)	41	61

All (25) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	249	ASP
1	A	301	LEU
1	A	372	LYS
1	A	396	ASP
1	A	442	PRO
1	A	448	THR
1	A	517	LEU
1	A	547	LEU
1	A	555	ASP
1	A	570	ARG
1	A	600	TYR
1	A	608	GLU
1	B	244	LYS
1	B	301	LEU
1	B	306	VAL
1	B	335	ASN
1	B	343	SER
1	B	352	LYS
1	B	387	ASN
1	B	470	MET
1	B	547	LEU
1	B	600	TYR
1	B	606	GLN
1	B	607	ASN
1	B	608	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such

sidechains are listed below:

Mol	Chain	Res	Type
1	A	201	HIS
1	A	387	ASN
1	A	434	GLN
1	A	460	GLN
1	A	518	ASN
1	A	580	GLN
1	A	606	GLN
1	A	607	ASN
1	B	201	HIS
1	B	335	ASN
1	B	387	ASN
1	B	460	GLN
1	B	518	ASN
1	B	552	GLN
1	B	606	GLN
1	B	607	ASN

### 5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

## 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 carbohydrates in this entry.

## 5.6 Ligand geometry [i](#)

There are no ligands in this entry.

## 5.7 Other polymers [i](#)

There are no such residues in this entry.

## 5.8 Polymer linkage issues ⓘ

There are no chain breaks in this entry.



## 6 Fit of model and data ⓘ

### 6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	436/436 (100%)	0.13	20 (4%) 33 31	13, 46, 76, 84	0
1	B	436/436 (100%)	0.25	28 (6%) 20 18	21, 48, 76, 86	0
2	C	6/6 (100%)	0.11	0 100 100	36, 45, 50, 50	0
2	D	6/6 (100%)	0.22	1 (16%) 2 2	40, 46, 50, 51	0
All	All	884/884 (100%)	0.19	49 (5%) 26 24	13, 47, 76, 86	0

All (49) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	419	THR	5.3
1	B	418	PHE	5.0
1	A	331	VAL	4.8
1	B	207	GLY	4.6
1	B	354	ILE	3.9
1	B	189	SER	3.9
1	B	360	LYS	3.7
1	B	388	ALA	3.6
1	B	385	GLY	3.6
1	A	207	GLY	3.6
1	A	360	LYS	3.5
1	B	479	GLY	3.5
1	B	416	THR	3.4
1	B	417	GLY	3.3
1	B	398	SER	3.3
1	B	347	ILE	3.2
1	A	407	ILE	3.1
1	B	407	ILE	3.0
1	A	418	PHE	3.0
1	B	399	VAL	2.9
1	A	347	ILE	2.8

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Mol	Chain	Res	Type	RSRZ
1	A	189	SER	2.8
1	B	586	ILE	2.8
1	A	206	THR	2.8
1	A	190	PRO	2.8
1	A	417	GLY	2.8
1	B	359	ILE	2.8
1	A	339	VAL	2.7
2	D	1	DA	2.6
1	A	208	SER	2.6
1	B	402	THR	2.5
1	A	351	GLY	2.5
1	B	593	HIS	2.4
1	A	281	GLY	2.4
1	B	420	GLY	2.3
1	B	356	ILE	2.3
1	A	377	LEU	2.3
1	B	396	ASP	2.3
1	B	587	ARG	2.2
1	B	608	GLU	2.2
1	A	357	GLU	2.2
1	A	468	GLY	2.2
1	B	397	VAL	2.2
1	A	220	ALA	2.1
1	A	354	ILE	2.1
1	B	480	GLU	2.1
1	A	416	THR	2.1
1	B	386	LEU	2.1
1	B	403	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 carbohydrates in this entry.

## 6.4 Ligands [i](#)

There are no ligands in this entry.

## 6.5 Other polymers [i](#)

There are no such residues in this entry.