



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 12, 2026 – 09:38 PM UTC

PDB ID : 4LK2 / pdb_00004lk2
Title : Crystal structure of RNA splicing effector Prp5
Authors : Zhang, Z.-M.; Li, J.; Yang, F.; Xu, Y.; Zhou, J.
Deposited on : 2013-07-05
Resolution : 2.12 Å(reported)

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

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity	:	4-5-2 with Phenix2.0
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

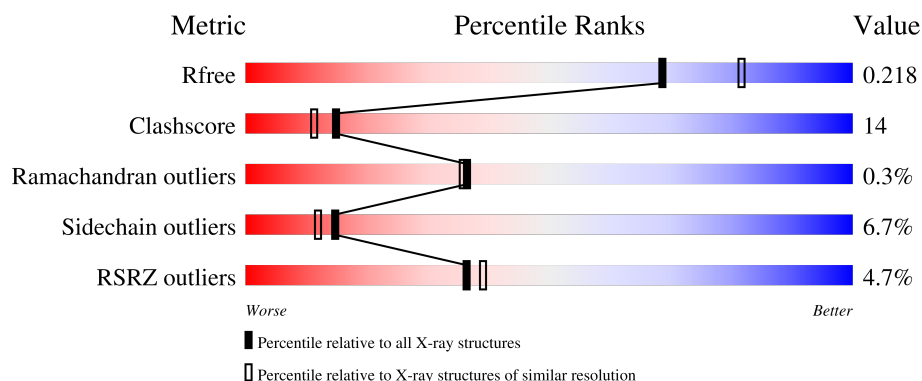
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.12 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}	180053	8290 (2.14-2.10)
Clashscore	190562	8817 (2.14-2.10)
Ramachandran outliers	187476	8738 (2.14-2.10)
Sidechain outliers	187428	8739 (2.14-2.10)
RSRZ outliers	180081	8294 (2.14-2.10)

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	493	4% 65% 24% 10%
1	B	493	4% 64% 22% 12%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 7012 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 Pre-mRNA-processing ATP-dependent RNA helicase PRP5.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	444	Total	C	N	O	S	0	5	0
			3436	2175	591	648	22			
1	B	436	Total	C	N	O	S	0	1	0
			3338	2114	571	632	21			

- Molecule 2 is NICKEL (II) ION (CCD ID: NI) (formula: Ni).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Ni	0	0
			1	1		

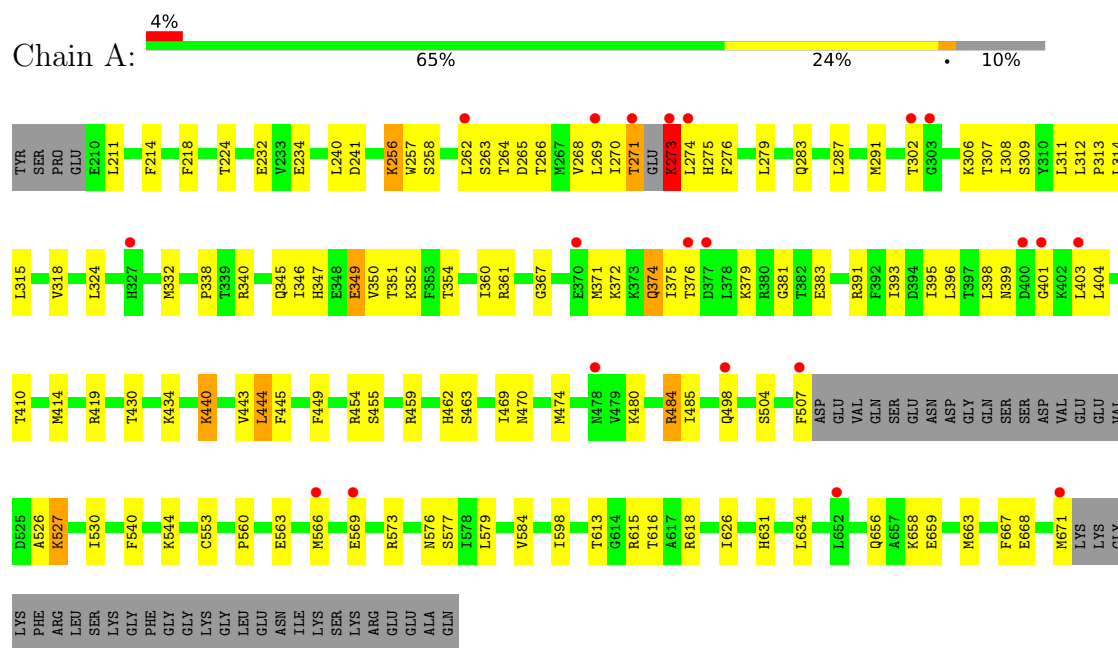
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	109	Total	O	0	0
			109	109		
3	B	128	Total	O	0	0
			128	128		

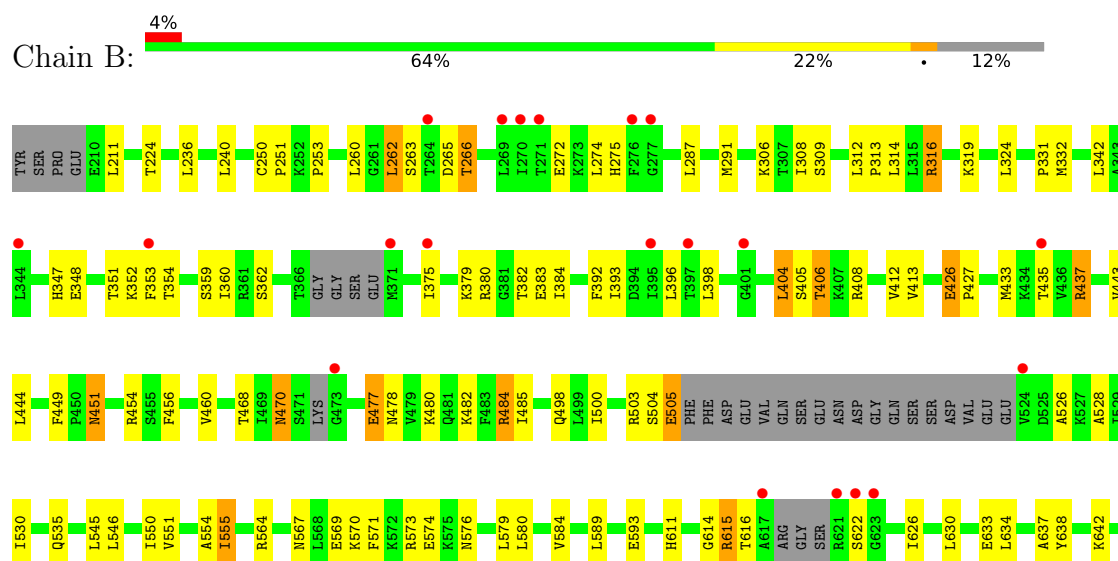
3 Residue-property plots

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: Pre-mRNA-processing ATP-dependent RNA helicase PRP5



• Molecule 1: Pre-mRNA-processing ATP-dependent RNA helicase PRP5



D646	I649	L652	D653	Q656	L660	M663	F667	M671	LYS	LYS	GLY	LYS	PHE	ARG	LEU	SER	LYS	GLY	PHE	GLY	GLY	LYS	GLY	LEU	GLU	ASN	ILE	LYS	SER	LYS	ARG	GLU	ALA	GLN
------	------	------	------	------	------	------	------	------	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----

4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	87.16Å 86.08Å 127.99Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	38.88 – 2.12 38.88 – 2.12	Depositor EDS
% Data completeness (in resolution range)	93.1 (38.88-2.12) 97.3 (38.88-2.12)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.98 (at 2.05Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.6.2_432)	Depositor
R, R_{free}	0.212 , 0.255 0.209 , 0.218	Depositor DCC
R_{free} test set	2946 reflections (4.83%)	wwPDB-VP
Wilson B-factor (Å ²)	33.3	Xtriage
Anisotropy	0.753	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 49.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.033 for k,h,-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	7012	wwPDB-VP
Average B, all atoms (Å ²)	43.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.47% 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: NI

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	0.76	1/3502 (0.0%)	0.89	2/4728 (0.0%)
1	B	0.49	0/3388	0.80	2/4576 (0.0%)
All	All	0.64	1/6890 (0.0%)	0.85	4/9304 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	271	THR	C-N	33.68	1.80	1.33

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	271	THR	O-C-N	-22.14	92.94	122.83
1	B	437	ARG	CA-C-N	5.63	125.73	119.87
1	B	437	ARG	C-N-CA	5.63	125.73	119.87
1	A	273	LYS	N-CA-C	5.27	125.75	111.00

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	3436	0	3485	96	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	B	3338	0	3371	92	0
2	A	1	0	0	0	0
3	A	109	0	0	4	0
3	B	128	0	0	5	0
All	All	7012	0	6856	188	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 (188) 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:271:THR:C	1:A:273:LYS:N	1.80	1.37
1:A:485:ILE:H	1:A:656:GLN:HE22	1.03	0.94
1:A:484:ARG:HD3	1:A:626:ILE:HD11	1.51	0.91
1:B:555:ILE:HD11	1:B:564:ARG:HA	1.58	0.86
1:A:560:PRO:HG2	1:A:563:GLU:HG2	1.58	0.85
1:A:256:LYS:HG3	1:A:258:SER:H	1.43	0.83
1:B:449:PHE:CE2	1:B:454:ARG:HG2	2.13	0.83
1:B:449:PHE:CZ	1:B:454:ARG:HG2	2.12	0.83
1:A:332:MET:HE1	1:A:383:GLU:HG3	1.60	0.81
1:B:484[A]:ARG:HH11	1:B:484[A]:ARG:HG2	1.46	0.80
1:B:406:THR:HG21	1:B:435:THR:HG22	1.61	0.79
1:B:526:ALA:H	1:B:576:ASN:ND2	1.80	0.78
1:B:332:MET:HE1	1:B:383:GLU:HG3	1.64	0.78
1:B:569:GLU:HG3	1:B:573:ARG:NH1	1.98	0.77
1:A:485:ILE:H	1:A:656:GLN:NE2	1.80	0.76
1:B:637:ALA:HB2	1:B:663:MET:HE3	1.68	0.76
1:A:218:PHE:HB2	1:A:291[A]:MET:HE1	1.69	0.75
1:B:236:LEU:HD23	1:B:253:PRO:HG2	1.68	0.74
1:B:480:LYS:HE2	1:B:482:LYS:NZ	2.01	0.74
1:A:291[A]:MET:HG2	1:A:313:PRO:HG3	1.68	0.74
1:A:658:LYS:O	1:A:658:LYS:HD2	1.87	0.73
1:A:257:TRP:HH2	1:A:308:ILE:HD13	1.53	0.72
1:B:468:THR:HG23	3:B:766:HOH:O	1.88	0.72
1:A:485:ILE:N	1:A:656:GLN:HE22	1.85	0.71
1:B:451:ASN:HD22	1:B:451:ASN:H	1.36	0.71
1:A:271:THR:O	1:A:273:LYS:N	2.22	0.71
1:B:477:GLU:H	1:B:477:GLU:CD	1.99	0.70
1:B:484[A]:ARG:HH11	1:B:484[A]:ARG:CG	2.05	0.69
1:B:503:ARG:HH12	1:B:576:ASN:HB3	1.59	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:451:ASN:H	1:B:451:ASN:ND2	1.93	0.65
1:A:306:LYS:HB2	1:A:445:PHE:CD1	2.31	0.65
1:B:478:ASN:O	1:B:622:SER:HB2	1.97	0.65
1:B:526:ALA:H	1:B:576:ASN:HD22	1.42	0.65
1:A:613:THR:O	1:A:616:THR:HG22	1.97	0.65
1:A:347:HIS:O	1:A:351:THR:HG23	1.95	0.64
1:B:375:ILE:HG22	1:B:379:LYS:HE3	1.79	0.64
1:A:395:ILE:O	1:A:403:LEU:HD21	1.98	0.64
1:B:316:ARG:HD2	3:B:770:HOH:O	1.97	0.63
1:B:503:ARG:HD3	1:B:528:ALA:HB2	1.80	0.63
1:A:584:VAL:HG13	3:A:908:HOH:O	1.97	0.63
1:B:593:GLU:CD	1:B:593:GLU:H	2.06	0.63
1:A:354:THR:HG22	1:A:360:ILE:HG22	1.80	0.63
1:A:264:THR:O	1:A:268:VAL:HG23	1.98	0.63
1:B:535:GLN:HG3	3:B:751:HOH:O	2.00	0.61
1:B:630:LEU:HB2	1:B:633:GLU:HG3	1.82	0.61
1:B:412:VAL:HG11	1:B:433:MET:HE1	1.83	0.61
1:A:312:LEU:HB2	1:A:313:PRO:HD3	1.83	0.60
1:A:324:LEU:HD21	1:A:332:MET:HE2	1.82	0.60
1:B:637:ALA:CB	1:B:663:MET:HE3	2.31	0.59
1:B:265:ASP:OD1	1:B:266:THR:N	2.35	0.58
1:A:332:MET:HE3	1:A:383:GLU:HA	1.85	0.58
1:A:374:GLN:OE1	1:A:391:ARG:HD2	2.04	0.56
1:A:526:ALA:O	1:A:576:ASN:HA	2.05	0.56
1:B:316:ARG:HA	1:B:319:LYS:HE3	1.87	0.56
1:A:265:ASP:OD1	1:A:266:THR:N	2.34	0.56
1:A:484:ARG:CD	1:A:626:ILE:HD11	2.29	0.56
1:B:593:GLU:CD	1:B:593:GLU:N	2.64	0.56
1:B:312:LEU:HB2	1:B:313:PRO:HD3	1.86	0.56
1:B:348:GLU:O	1:B:352:LYS:HG2	2.06	0.56
1:A:340:ARG:HD2	1:A:367:GLY:O	2.06	0.56
1:A:504:SER:HA	1:A:507:PHE:CE2	2.41	0.56
1:A:455:SER:O	1:A:459[B]:ARG:HG3	2.06	0.56
1:B:504:SER:C	1:B:505:GLU:HG3	2.30	0.56
1:B:646:ASP:OD2	1:B:646:ASP:N	2.38	0.55
1:A:351:THR:HA	1:A:354:THR:OG1	2.07	0.55
1:A:314:LEU:O	1:A:318:VAL:HG22	2.07	0.55
1:A:372:LYS:O	1:A:376:THR:HG23	2.06	0.54
1:B:555:ILE:HD11	1:B:564:ARG:CA	2.36	0.54
1:B:359:SER:C	1:B:360:ILE:HD12	2.33	0.54
1:B:426:GLU:N	1:B:427:PRO:HD2	2.23	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:634:LEU:HB3	1:A:667:PHE:HB2	1.90	0.54
1:B:456:PHE:O	1:B:460:VAL:HG12	2.07	0.54
1:A:569:GLU:O	1:A:573:ARG:HB2	2.07	0.54
1:A:269:LEU:O	1:A:273:LYS:CB	2.56	0.54
1:B:324:LEU:HD21	1:B:332:MET:HE2	1.90	0.53
1:A:256:LYS:HG2	1:A:258:SER:OG	2.08	0.53
1:A:270:ILE:C	1:A:274:LEU:H	2.17	0.53
1:A:218:PHE:CB	1:A:291[A]:MET:HE1	2.37	0.53
1:B:555:ILE:CD1	1:B:564:ARG:HA	2.37	0.52
1:A:553:CYS:O	1:A:579:LEU:HD12	2.09	0.52
1:A:414:MET:HB2	1:A:444:LEU:HD12	1.91	0.52
1:A:484:ARG:HH22	1:A:498:GLN:NE2	2.08	0.52
1:A:659:GLU:HG2	1:A:663:MET:HE3	1.90	0.52
1:B:545:LEU:O	1:B:550:ILE:HB	2.10	0.52
1:B:480:LYS:HE2	1:B:482:LYS:HZ1	1.74	0.51
1:B:503:ARG:NH1	1:B:576:ASN:HB3	2.25	0.51
1:B:347:HIS:O	1:B:351:THR:HG23	2.10	0.51
1:B:554:ALA:O	1:B:555:ILE:HD12	2.10	0.51
1:A:527:LYS:HE2	1:A:577:SER:OG	2.12	0.50
1:A:283:GLN:NE2	1:A:308:ILE:HD11	2.26	0.50
1:B:412:VAL:HG11	1:B:433:MET:CE	2.41	0.50
1:B:480:LYS:HE2	1:B:482:LYS:HZ2	1.75	0.49
1:A:306:LYS:HB2	1:A:445:PHE:CE1	2.47	0.49
1:B:404:LEU:HD13	1:B:405:SER:N	2.28	0.49
1:B:260:LEU:HB3	1:B:262:LEU:HD13	1.94	0.49
1:A:257:TRP:CH2	1:A:308:ILE:HD13	2.39	0.49
1:A:265:ASP:CG	1:A:266:THR:H	2.20	0.49
1:A:419:ARG:HD2	3:A:902:HOH:O	2.13	0.48
1:B:634:LEU:HB3	1:B:667:PHE:HB2	1.96	0.48
1:A:241:ASP:O	1:A:470:ASN:HB2	2.13	0.48
1:B:404:LEU:HD13	1:B:405:SER:H	1.79	0.48
1:A:371:MET:O	1:A:375:ILE:HG13	2.14	0.48
1:B:331:PRO:HD3	1:B:404:LEU:HD11	1.94	0.48
1:A:240:LEU:O	1:A:241:ASP:HB2	2.14	0.48
1:A:265:ASP:CG	1:A:266:THR:N	2.72	0.48
1:B:211:LEU:HD23	1:B:437:ARG:HB2	1.95	0.47
1:B:555:ILE:CD1	1:B:567:ASN:HB2	2.45	0.47
1:B:500:ILE:HD11	1:B:530:ILE:HD11	1.96	0.47
1:A:287:LEU:O	1:A:291[B]:MET:HB2	2.14	0.47
1:B:314:LEU:C	1:B:314:LEU:HD23	2.40	0.47
1:A:263:SER:HB3	1:A:265:ASP:OD1	2.15	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:371:MET:O	1:A:374:GLN:HG2	2.15	0.47
1:A:462:HIS:O	1:A:463:SER:C	2.58	0.47
1:B:352:LYS:N	1:B:352:LYS:HD3	2.30	0.46
1:A:396:LEU:O	1:A:401:GLY:HA2	2.15	0.46
1:A:273:LYS:O	1:A:276:PHE:CE2	2.68	0.46
1:A:631:HIS:HA	1:A:663:MET:HE2	1.98	0.46
1:A:469:ILE:O	1:A:470:ASN:C	2.59	0.46
1:B:526:ALA:N	1:B:576:ASN:HD22	2.12	0.46
1:A:257:TRP:HH2	1:A:308:ILE:CD1	2.25	0.45
1:A:540:PHE:CE2	1:A:544:LYS:HG3	2.51	0.45
1:B:287:LEU:O	1:B:291:MET:HG3	2.16	0.45
1:B:449:PHE:CE1	1:B:454:ARG:HG2	2.48	0.45
1:B:484[A]:ARG:HG3	1:B:626:ILE:HD11	1.99	0.45
1:B:451:ASN:ND2	1:B:451:ASN:N	2.64	0.45
1:A:398:LEU:HD23	1:A:399:ASN:HB2	1.99	0.45
1:B:331:PRO:HA	1:B:382:THR:O	2.17	0.45
1:B:660:LEU:HA	1:B:663:MET:HE2	1.97	0.45
1:A:218:PHE:HB2	1:A:291[A]:MET:CE	2.41	0.45
1:A:338:PRO:HG3	1:A:419:ARG:HB2	1.98	0.45
1:A:440:LYS:N	1:A:440:LYS:HD3	2.32	0.45
1:B:211:LEU:CD2	1:B:437:ARG:HB2	2.47	0.45
1:B:584:VAL:HG13	3:B:735:HOH:O	2.17	0.45
1:A:615:ARG:HD3	3:A:888:HOH:O	2.17	0.44
1:A:403:LEU:H	1:A:403:LEU:HD23	1.82	0.44
1:B:362:SER:HA	1:B:384:ILE:O	2.17	0.44
1:B:413:VAL:HG22	1:B:443:VAL:CG2	2.46	0.44
1:A:306:LYS:O	1:A:309:SER:HB2	2.17	0.44
1:B:449:PHE:CD2	1:B:454:ARG:HG2	2.51	0.44
1:A:307:THR:HG22	1:A:311:LEU:HD11	2.00	0.44
1:A:398:LEU:HD22	1:A:403:LEU:HD22	1.99	0.44
1:A:668:GLU:O	1:A:671:MET:HG2	2.17	0.44
1:B:653:ASP:OD1	1:B:653:ASP:N	2.36	0.44
1:B:306:LYS:O	1:B:309:SER:HB2	2.18	0.44
1:B:638:TYR:CZ	1:B:642:LYS:HD2	2.53	0.44
1:B:649:ILE:HA	1:B:652:LEU:CD1	2.48	0.44
1:A:449:PHE:CD2	1:A:454:ARG:NH2	2.86	0.43
1:B:354:THR:HG21	1:B:360:ILE:HG22	2.00	0.43
1:A:270:ILE:HG22	1:A:270:ILE:O	2.17	0.43
1:A:307:THR:O	1:A:311:LEU:HG	2.19	0.43
1:B:211:LEU:HD12	1:B:211:LEU:N	2.33	0.43
1:A:275:HIS:CD2	1:A:283:GLN:HE22	2.37	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:346:ILE:O	1:A:350:VAL:HG23	2.18	0.43
1:A:361[A]:ARG:HH21	1:A:381:GLY:C	2.27	0.43
1:A:504:SER:HA	1:A:507:PHE:CD2	2.54	0.43
1:B:484[B]:ARG:HG3	1:B:626:ILE:HD11	2.01	0.43
1:A:459[B]:ARG:HB3	1:A:459[B]:ARG:CZ	2.49	0.43
1:B:570:LYS:O	1:B:574:GLU:HG3	2.18	0.43
1:A:430:THR:O	1:A:434:LYS:HG3	2.19	0.42
1:A:658:LYS:HD3	1:A:658:LYS:HA	1.83	0.42
1:A:324:LEU:CD2	1:A:332:MET:HE2	2.48	0.42
1:A:354:THR:CG2	1:A:360:ILE:HG22	2.48	0.42
1:B:480:LYS:HB2	1:B:480:LYS:HE3	1.74	0.42
1:B:614:GLY:O	1:B:615:ARG:NH1	2.53	0.42
1:A:631:HIS:HA	1:A:663:MET:CE	2.49	0.42
1:B:412:VAL:HG21	1:B:433:MET:HE1	2.01	0.42
1:B:571:PHE:CD1	1:B:579:LEU:HB2	2.54	0.42
1:B:611:HIS:HD2	3:B:761:HOH:O	2.02	0.42
1:A:398:LEU:HD23	1:A:399:ASN:CB	2.49	0.42
1:A:474:MET:HE3	1:A:474:MET:HB2	1.80	0.42
1:B:332:MET:HA	1:B:408:ARG:O	2.20	0.42
1:A:345:GLN:O	1:A:349:GLU:OE1	2.37	0.42
1:B:375:ILE:O	1:B:379:LYS:HG3	2.20	0.41
1:B:380:ARG:HD3	1:B:380:ARG:HA	1.82	0.41
1:B:449:PHE:CZ	1:B:454:ARG:HA	2.55	0.41
1:A:283:GLN:CD	1:A:308:ILE:HD11	2.45	0.41
1:A:302:THR:HG21	3:A:897:HOH:O	2.19	0.41
1:A:374:GLN:HE21	1:A:374:GLN:HB3	1.59	0.41
1:B:392:PHE:CE2	1:B:396:LEU:HD11	2.56	0.41
1:A:257:TRP:CD1	1:A:279:LEU:HD11	2.56	0.40
1:B:260:LEU:CB	1:B:262:LEU:HD13	2.52	0.40
1:B:470:ASN:HD22	1:B:470:ASN:HA	1.70	0.40
1:A:214:PHE:CE1	1:A:410:THR:HB	2.57	0.40
1:A:393:ILE:HA	1:A:393:ILE:HD13	1.87	0.40
1:B:250:CYS:HA	1:B:251:PRO:HD3	1.95	0.40
1:B:308:ILE:HD12	1:B:353:PHE:HZ	1.86	0.40
1:B:546:LEU:HA	1:B:550:ILE:O	2.22	0.40
1:A:530:ILE:HA	1:A:598:ILE:O	2.22	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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	445/493 (90%)	431 (97%)	13 (3%)	1 (0%)	43	44
1	B	427/493 (87%)	415 (97%)	10 (2%)	2 (0%)	24	22
All	All	872/986 (88%)	846 (97%)	23 (3%)	3 (0%)	36	36

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	272	GLU
1	A	273	LYS
1	B	274	LEU

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	380/437 (87%)	358 (94%)	22 (6%)	18	16
1	B	369/437 (84%)	338 (92%)	31 (8%)	10	7
All	All	749/874 (86%)	696 (93%)	53 (7%)	15	10

All (53) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	211	LEU
1	A	224	THR
1	A	232	GLU

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Mol	Chain	Res	Type
1	A	234[A]	GLU
1	A	234[B]	GLU
1	A	256	LYS
1	A	262	LEU
1	A	315	LEU
1	A	349	GLU
1	A	352[A]	LYS
1	A	352[B]	LYS
1	A	374	GLN
1	A	379	LYS
1	A	404	LEU
1	A	440	LYS
1	A	443	VAL
1	A	444	LEU
1	A	480	LYS
1	A	484	ARG
1	A	527	LYS
1	A	566	MET
1	A	618	ARG
1	B	224	THR
1	B	240	LEU
1	B	262	LEU
1	B	263	SER
1	B	266	THR
1	B	275	HIS
1	B	316	ARG
1	B	342	LEU
1	B	393	ILE
1	B	398	LEU
1	B	404	LEU
1	B	406	THR
1	B	426	GLU
1	B	444	LEU
1	B	451	ASN
1	B	470	ASN
1	B	477	GLU
1	B	484[A]	ARG
1	B	484[B]	ARG
1	B	485	ILE
1	B	498	GLN
1	B	505	GLU
1	B	551	VAL

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Mol	Chain	Res	Type
1	B	555	ILE
1	B	580	LEU
1	B	589	LEU
1	B	615	ARG
1	B	616	THR
1	B	646	ASP
1	B	653	ASP
1	B	656	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (14) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	275	HIS
1	A	478	ASN
1	A	481	GLN
1	A	498	GLN
1	A	576	ASN
1	A	656	GLN
1	B	374	GLN
1	B	451	ASN
1	B	470	ASN
1	B	487	HIS
1	B	547	ASN
1	B	567	ASN
1	B	576	ASN
1	B	656	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 oligosaccharides in this entry.

5.6 Ligand geometry

Of 1 ligands modelled in this entry, 1 is 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, 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	444/493 (90%)	0.43	21 (4%) 36 39	20, 40, 64, 73	5 (1%)
1	B	436/493 (88%)	0.51	20 (4%) 37 40	24, 43, 68, 82	1 (0%)
All	All	880/986 (89%)	0.47	41 (4%) 36 39	20, 42, 65, 82	6 (0%)

All (41) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	271	THR	3.9
1	A	274	LEU	3.8
1	B	617	ALA	3.5
1	B	623	GLY	3.5
1	B	269	LEU	3.2
1	B	401	GLY	3.1
1	B	371	MET	3.0
1	A	376	THR	3.0
1	A	566	MET	2.9
1	B	276	PHE	2.9
1	A	377	ASP	2.7
1	A	403	LEU	2.7
1	B	395	ILE	2.7
1	B	277	GLY	2.7
1	B	473	GLY	2.7
1	A	507	PHE	2.6
1	B	264	THR	2.6
1	A	273	LYS	2.6
1	A	269	LEU	2.6
1	B	270	ILE	2.6
1	B	435	THR	2.5
1	A	400	ASP	2.5
1	A	498	GLN	2.5
1	A	569	GLU	2.5

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Mol	Chain	Res	Type	RSRZ
1	B	397	THR	2.5
1	B	344	LEU	2.4
1	A	370	GLU	2.4
1	B	353	PHE	2.4
1	B	621	ARG	2.4
1	A	478	ASN	2.3
1	A	671	MET	2.3
1	A	262	LEU	2.3
1	A	302	THR	2.3
1	B	271	THR	2.3
1	A	303	GLY	2.2
1	A	401	GLY	2.2
1	B	524	VAL	2.1
1	A	652	LEU	2.1
1	B	622	SER	2.1
1	A	327	HIS	2.1
1	B	375	ILE	2.1

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 oligosaccharides 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($\text{\AA}^2$)	Q<0.9
2	NI	A	701	1/1	0.99	0.04	32,32,32,32	0

6.5 Other polymers [i](#)

There are no such residues in this entry.