



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 07:07 AM UTC

PDB ID : 2HIV / pdb_00002hiv
Title : ATP-dependent DNA ligase from *S. solfataricus*
Authors : Pascal, J.M.; Ellenberger, T.
Deposited on : 2006-06-29
Resolution : 2.05 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

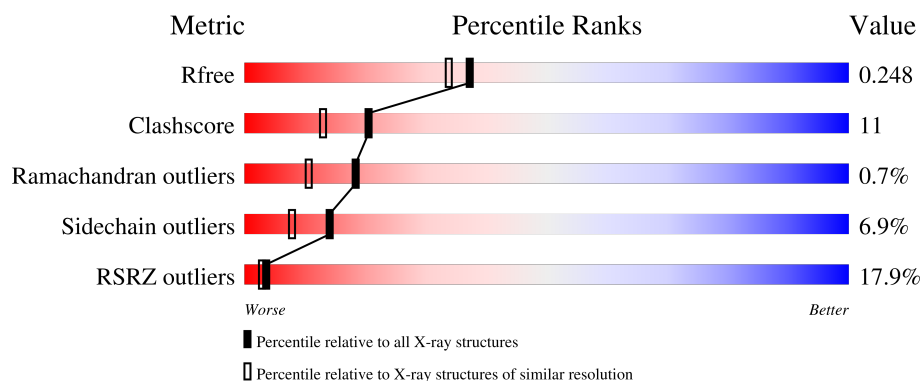
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.05 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	2260 (2.04-2.04)
Clashscore	190562	2333 (2.04-2.04)
Ramachandran outliers	187476	2318 (2.04-2.04)
Sidechain outliers	187428	2318 (2.04-2.04)
RSRZ outliers	180081	2260 (2.04-2.04)

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	621	17% 71% 19% 7%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 4836 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 Thermostable DNA ligase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	577	Total	C	N	O	S	0	6	0
			4611	2954	776	866	15			

There are 20 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-19	MET	-	cloning artifact	UNP Q980T8
A	-18	GLY	-	cloning artifact	UNP Q980T8
A	-17	SER	-	cloning artifact	UNP Q980T8
A	-16	SER	-	cloning artifact	UNP Q980T8
A	-15	HIS	-	expression tag	UNP Q980T8
A	-14	HIS	-	expression tag	UNP Q980T8
A	-13	HIS	-	expression tag	UNP Q980T8
A	-12	HIS	-	expression tag	UNP Q980T8
A	-11	HIS	-	expression tag	UNP Q980T8
A	-10	HIS	-	expression tag	UNP Q980T8
A	-9	SER	-	cloning artifact	UNP Q980T8
A	-8	SER	-	cloning artifact	UNP Q980T8
A	-7	GLY	-	cloning artifact	UNP Q980T8
A	-6	LEU	-	cloning artifact	UNP Q980T8
A	-5	VAL	-	cloning artifact	UNP Q980T8
A	-4	PRO	-	cloning artifact	UNP Q980T8
A	-3	ARG	-	cloning artifact	UNP Q980T8
A	-2	GLY	-	cloning artifact	UNP Q980T8
A	-1	SER	-	cloning artifact	UNP Q980T8
A	0	HIS	-	cloning artifact	UNP Q980T8

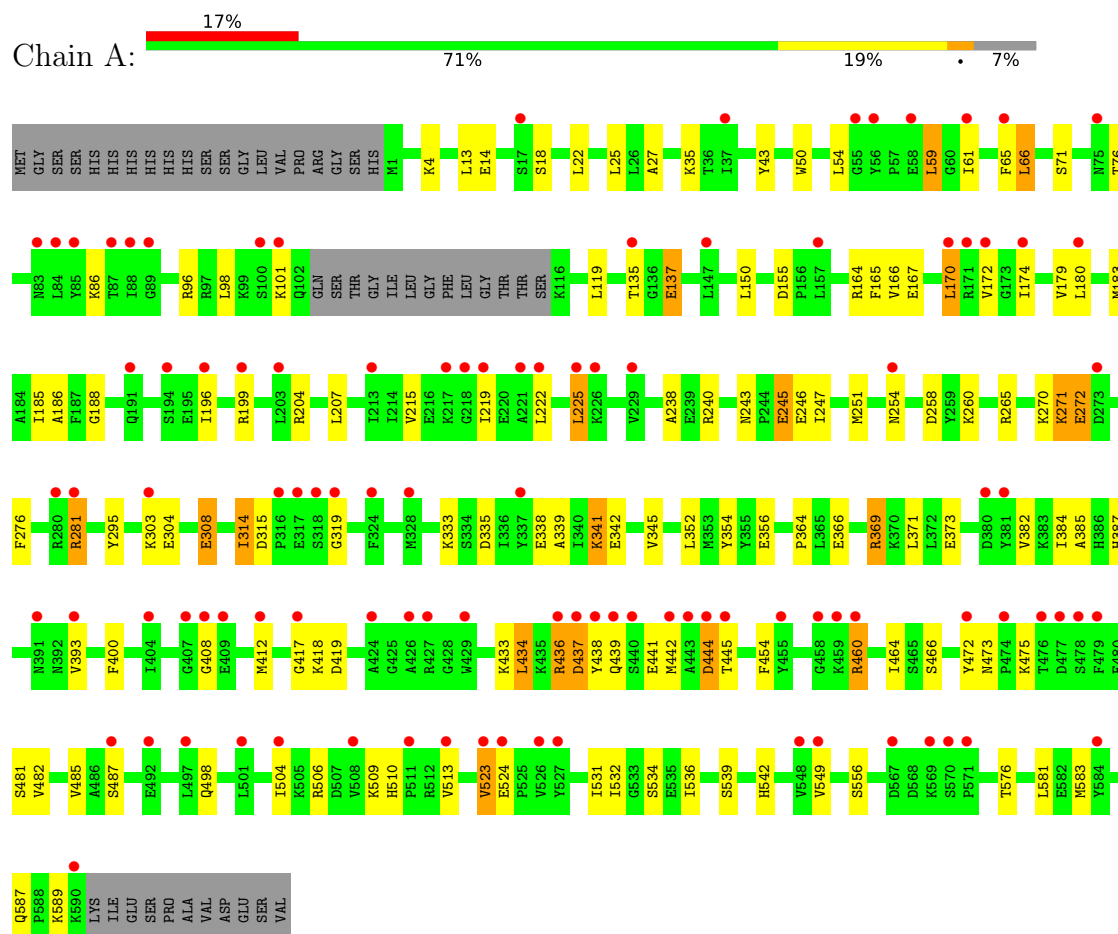
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	225	Total	O	0	0
			225	225		

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: Thermostable DNA ligase



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	119.21Å 169.77Å 77.31Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	42.64 – 2.05 42.64 – 2.05	Depositor EDS
% Data completeness (in resolution range)	97.2 (42.64-2.05) 97.1 (42.64-2.05)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.59 (at 2.05Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.209 , 0.254 0.203 , 0.248	Depositor DCC
R_{free} test set	2423 reflections (4.89%)	wwPDB-VP
Wilson B-factor (Å ²)	35.4	Xtriage
Anisotropy	0.252	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 45.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	4836	wwPDB-VP
Average B, all atoms (Å ²)	49.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.07% 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](#)

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	0.75	2/4708 (0.0%)	0.96	2/6341 (0.0%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	369	ARG	CD-NE	-5.78	1.38	1.46
1	A	308	GLU	CB-CG	-5.66	1.35	1.52

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	369	ARG	NE-CZ-NH2	-6.95	112.94	119.20
1	A	170	LEU	N-CA-C	-6.95	98.62	109.25

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	4611	0	4714	104	2
2	A	225	0	0	11	1
All	All	4836	0	4714	104	2

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 11.

All (104) 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:243:ASN:HD21	1:A:245:GLU:HG2	1.06	1.18
1:A:196:ILE:HD11	1:A:219:ILE:HG23	1.37	1.03
1:A:536:ILE:H	1:A:587:GLN:HE22	1.05	1.02
1:A:179:VAL:HG12	1:A:183:MET:HE2	1.48	0.95
1:A:196:ILE:CD1	1:A:219:ILE:HG23	1.99	0.91
1:A:183:MET:HE1	1:A:207:LEU:HD22	1.53	0.91
1:A:243:ASN:ND2	1:A:245:GLU:HG2	1.89	0.88
1:A:473:ASN:OD1	1:A:475:LYS:HE3	1.74	0.86
1:A:183:MET:HE1	1:A:207:LEU:CD2	2.07	0.83
1:A:180:LEU:HA	1:A:183:MET:HE3	1.61	0.80
1:A:170:LEU:O	2:A:777:HOH:O	2.00	0.80
1:A:281[B]:ARG:HH11	1:A:281[B]:ARG:HG2	1.46	0.79
1:A:454:PHE:HB2	1:A:466:SER:HB3	1.65	0.78
1:A:265:ARG:NE	1:A:308:GLU:OE1	2.15	0.77
1:A:460:ARG:O	1:A:460:ARG:HG3	1.85	0.76
1:A:243:ASN:HD22	1:A:246:GLU:H	1.35	0.75
1:A:442:MET:HA	1:A:534:SER:HA	1.69	0.74
1:A:444:ASP:HB2	1:A:583:MET:CE	2.17	0.74
1:A:196:ILE:HG23	2:A:644:HOH:O	1.88	0.73
1:A:510:HIS:HD2	1:A:524[B]:GLU:OE1	1.71	0.72
1:A:258:ASP:OD2	1:A:369:ARG:NH2	2.23	0.69
1:A:281[B]:ARG:HH11	1:A:281[B]:ARG:CG	2.06	0.69
1:A:314:ILE:HG13	1:A:319:GLY:O	1.93	0.69
1:A:13:LEU:HD21	1:A:172:VAL:HG13	1.76	0.68
1:A:13:LEU:CD2	1:A:172:VAL:HG13	2.24	0.67
1:A:183:MET:CE	1:A:207:LEU:HD22	2.25	0.67
1:A:271:LYS:HE3	1:A:276:PHE:CZ	2.30	0.67
1:A:510:HIS:CD2	1:A:524[B]:GLU:OE1	2.48	0.65
1:A:188:GLY:HA2	1:A:219:ILE:HG13	1.79	0.64
1:A:199:ARG:HD3	2:A:803:HOH:O	1.97	0.64
1:A:444:ASP:HB2	1:A:583:MET:HE2	1.80	0.62
1:A:418:LYS:HD3	2:A:617:HOH:O	2.00	0.62
1:A:14:GLU:OE2	1:A:164:ARG:NH1	2.32	0.62
1:A:542:HIS:HE1	2:A:755:HOH:O	1.83	0.62
1:A:98:LEU:O	1:A:101:LYS:HG2	2.00	0.62
1:A:536:ILE:H	1:A:587:GLN:NE2	1.88	0.61
1:A:539:SER:HB2	1:A:556:SER:HB2	1.81	0.61
1:A:464:ILE:O	1:A:498[A]:GLN:NE2	2.33	0.60
1:A:172:VAL:HG12	1:A:174:ILE:HG23	1.83	0.59
1:A:43:TYR:OH	2:A:619:HOH:O	2.16	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:369:ARG:HD2	2:A:606:HOH:O	2.02	0.59
1:A:258:ASP:CG	1:A:369:ARG:HH22	2.10	0.58
1:A:510:HIS:O	1:A:513:VAL:HG22	2.03	0.58
1:A:238:ALA:CB	1:A:412:MET:HE1	2.36	0.56
1:A:385:ALA:O	1:A:387:HIS:HD2	1.89	0.55
1:A:35:LYS:HB2	1:A:215:VAL:CG1	2.36	0.55
1:A:66:LEU:CD1	1:A:167:GLU:HG3	2.37	0.55
1:A:437:ASP:C	1:A:439:GLN:H	2.13	0.55
1:A:335:ASP:HB3	1:A:338:GLU:HB2	1.89	0.55
1:A:387:HIS:HE1	2:A:661:HOH:O	1.90	0.55
1:A:281[B]:ARG:HG2	1:A:281[B]:ARG:NH1	2.20	0.54
1:A:137:GLU:O	2:A:813:HOH:O	2.18	0.54
1:A:338:GLU:HA	1:A:341:LYS:HG2	1.89	0.54
1:A:260:LYS:HD3	1:A:412:MET:SD	2.48	0.54
1:A:444:ASP:HB2	1:A:583:MET:HE1	1.90	0.53
1:A:240:ARG:HH21	1:A:437:ASP:CB	2.22	0.53
1:A:61:ILE:HG12	1:A:65:PHE:HB2	1.92	0.52
1:A:504:ILE:HG13	1:A:523:VAL:HG22	1.90	0.52
1:A:270:LYS:O	1:A:304:GLU:HA	2.09	0.51
1:A:400:PHE:CD1	1:A:434:LEU:HD13	2.44	0.51
1:A:369:ARG:O	1:A:373:GLU:HG3	2.10	0.51
1:A:408:GLY:HA2	2:A:756:HOH:O	2.10	0.50
1:A:509:LYS:HG3	1:A:513:VAL:CG2	2.42	0.50
1:A:165:PHE:CD1	1:A:170:LEU:HD11	2.48	0.49
1:A:295:TYR:HB3	1:A:382:VAL:HG23	1.93	0.49
1:A:498[A]:GLN:H	1:A:498[A]:GLN:CD	2.21	0.49
1:A:243:ASN:ND2	1:A:246:GLU:H	2.08	0.49
1:A:393:VAL:CG2	2:A:614:HOH:O	2.61	0.49
1:A:50:TRP:HB2	1:A:54:LEU:HD12	1.93	0.48
1:A:315:ASP:O	1:A:319:GLY:N	2.47	0.47
1:A:481:SER:OG	1:A:542:HIS:HD2	1.96	0.47
1:A:472:TYR:O	1:A:576:THR:HA	2.15	0.47
1:A:436:ARG:C	1:A:438:TYR:H	2.23	0.46
1:A:366:GLU:CD	1:A:417:GLY:H	2.23	0.46
1:A:333:LYS:HB2	1:A:339:ALA:HB2	1.97	0.46
1:A:240:ARG:NH2	1:A:437:ASP:CB	2.79	0.45
1:A:86:LYS:H	1:A:86:LYS:HG2	1.56	0.45
1:A:119:LEU:HD21	1:A:150:LEU:HD23	1.98	0.45
1:A:509:LYS:CG	1:A:513:VAL:HG23	2.46	0.45
1:A:341:LYS:HG3	1:A:342:GLU:N	2.32	0.45
1:A:295:TYR:HB3	1:A:382:VAL:CG2	2.47	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:412:MET:SD	1:A:433:LYS:HB3	2.56	0.45
1:A:66:LEU:HD11	1:A:167:GLU:HG3	1.98	0.45
1:A:271:LYS:HE3	1:A:276:PHE:CE1	2.52	0.44
1:A:509:LYS:CG	1:A:513:VAL:CG2	2.94	0.44
1:A:247:ILE:HD12	1:A:434:LEU:HG	1.99	0.44
1:A:364:PRO:HB3	1:A:417:GLY:O	2.16	0.44
1:A:485:VAL:HG21	1:A:531:ILE:HD12	2.00	0.44
1:A:304:GLU:HG3	1:A:356:GLU:HA	2.01	0.43
1:A:444:ASP:O	1:A:532:ILE:HG13	2.19	0.43
1:A:251:MET:HA	1:A:251:MET:HE2	2.01	0.43
1:A:304:GLU:OE1	1:A:354:TYR:HE2	2.01	0.43
1:A:506:ARG:HD3	1:A:510:HIS:HB2	2.00	0.43
1:A:196:ILE:CD1	1:A:219:ILE:CG2	2.85	0.42
1:A:384:ILE:O	1:A:385:ALA:C	2.60	0.42
1:A:315:ASP:O	1:A:319:GLY:HA2	2.19	0.42
1:A:204:ARG:CZ	1:A:225:LEU:HD21	2.50	0.41
1:A:196:ILE:HD12	1:A:222:LEU:HD12	2.02	0.41
1:A:238:ALA:HB1	1:A:412:MET:HE1	2.02	0.41
1:A:437:ASP:C	1:A:439:GLN:N	2.78	0.41
1:A:59:LEU:HD13	1:A:166:VAL:HA	2.02	0.40
1:A:183:MET:CE	1:A:207:LEU:CD2	2.88	0.40
1:A:27:ALA:HA	1:A:186:ALA:HB2	2.04	0.40

All (2) 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:A:506:ARG:NH1	2:A:811:HOH:O[3_555]	1.79	0.41
1:A:18:SER:OG	1:A:419:ASP:OD2[6_555]	2.08	0.12

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	579/621 (93%)	553 (96%)	22 (4%)	4 (1%)	18 10

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	272	GLU
1	A	436	ARG
1	A	437	ASP
1	A	441	GLU

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	497/536 (93%)	462 (93%)	35 (7%)	14 7

All (35) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	4	LYS
1	A	22	LEU
1	A	25	LEU
1	A	59	LEU
1	A	66	LEU
1	A	71	SER
1	A	76	THR
1	A	96	ARG
1	A	135	THR
1	A	137	GLU
1	A	155	ASP
1	A	185	ILE
1	A	225	LEU
1	A	245	GLU
1	A	254	ASN
1	A	271	LYS
1	A	272	GLU

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Mol	Chain	Res	Type
1	A	281[A]	ARG
1	A	281[B]	ARG
1	A	303	LYS
1	A	314	ILE
1	A	341	LYS
1	A	345	VAL
1	A	352	LEU
1	A	371	LEU
1	A	434	LEU
1	A	444	ASP
1	A	445	THR
1	A	460	ARG
1	A	482	VAL
1	A	487	SER
1	A	523	VAL
1	A	549	VAL
1	A	581	LEU
1	A	589	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	46	GLN
1	A	243	ASN
1	A	329	HIS
1	A	346	ASN
1	A	387	HIS
1	A	542	HIS
1	A	587	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 [i](#)

There are no oligosaccharides 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 [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	577/621 (92%)	1.34	103 (17%) 3 3	24, 47, 61, 77	6 (1%)

All (103) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	438	TYR	7.2
1	A	172	VAL	5.4
1	A	444	ASP	4.9
1	A	229	VAL	4.9
1	A	442	MET	4.6
1	A	316	PRO	4.5
1	A	170	LEU	4.1
1	A	526	VAL	4.1
1	A	281[A]	ARG	4.0
1	A	89	GLY	4.0
1	A	337	TYR	4.0
1	A	88	ILE	3.8
1	A	319	GLY	3.8
1	A	437	ASP	3.6
1	A	196	ILE	3.6
1	A	171	ARG	3.5
1	A	443	ALA	3.5
1	A	324	PHE	3.3
1	A	440	SER	3.3
1	A	472	TYR	3.3
1	A	479	PHE	3.3
1	A	590	LYS	3.2
1	A	497	LEU	3.2
1	A	222	LEU	3.2
1	A	460	ARG	3.1
1	A	513	VAL	3.1
1	A	225	LEU	3.1

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Mol	Chain	Res	Type	RSRZ
1	A	458	GLY	3.0
1	A	61	ILE	2.9
1	A	476	THR	2.9
1	A	474	PRO	2.9
1	A	85	TYR	2.9
1	A	501	LEU	2.8
1	A	571	PRO	2.8
1	A	199	ARG	2.8
1	A	213	ILE	2.8
1	A	87	THR	2.6
1	A	569	LYS	2.6
1	A	408	GLY	2.6
1	A	317	GLU	2.6
1	A	318	SER	2.6
1	A	221	ALA	2.6
1	A	436	ARG	2.6
1	A	147	LEU	2.6
1	A	393	VAL	2.5
1	A	427[A]	ARG	2.5
1	A	508	VAL	2.5
1	A	477	ASP	2.5
1	A	478	SER	2.5
1	A	426	ALA	2.5
1	A	37	ILE	2.5
1	A	570	SER	2.4
1	A	412	MET	2.4
1	A	459	LYS	2.4
1	A	404	ILE	2.4
1	A	524[A]	GLU	2.4
1	A	523	VAL	2.4
1	A	58	GLU	2.4
1	A	219	ILE	2.4
1	A	303	LYS	2.4
1	A	511	PRO	2.4
1	A	445	THR	2.4
1	A	455	TYR	2.3
1	A	203	LEU	2.3
1	A	254	ASN	2.3
1	A	391	ASN	2.3
1	A	527	TYR	2.3
1	A	584	TYR	2.3
1	A	549	VAL	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	381	TYR	2.2
1	A	409	GLU	2.2
1	A	100	SER	2.2
1	A	83	ASN	2.2
1	A	157	LEU	2.2
1	A	194	SER	2.2
1	A	273	ASP	2.2
1	A	548	VAL	2.2
1	A	180	LEU	2.2
1	A	280	ARG	2.2
1	A	101	LYS	2.2
1	A	84	LEU	2.1
1	A	487	SER	2.1
1	A	226	LYS	2.1
1	A	56	TYR	2.1
1	A	328	MET	2.1
1	A	135	THR	2.1
1	A	191	GLN	2.1
1	A	439	GLN	2.1
1	A	75	ASN	2.1
1	A	380	ASP	2.1
1	A	174	ILE	2.1
1	A	504	ILE	2.1
1	A	17	SER	2.1
1	A	492	GLU	2.1
1	A	55	GLY	2.1
1	A	218	GLY	2.1
1	A	407	GLY	2.1
1	A	429	TRP	2.1
1	A	424	ALA	2.1
1	A	217	LYS	2.0
1	A	65	PHE	2.0
1	A	567	ASP	2.0
1	A	417	GLY	2.0

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 oligosaccharides 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.