



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 02:09 PM UTC

PDB ID : 9LL6 / pdb_00009ll6
Title : Crystal structure of Ornithine decarboxylase H216F mutant without PLP
Authors : Kim, D.S.; Park, J.H.; Adiko, N.N.; Seo, H.T.
Deposited on : 2025-01-17
Resolution : 2.64 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

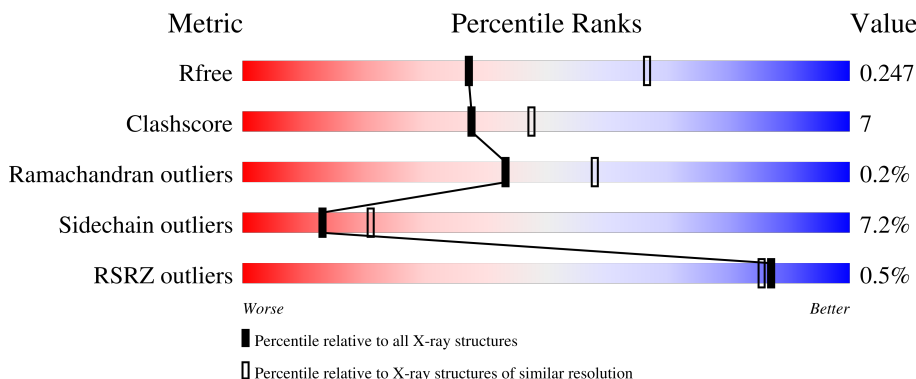
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.64 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	2053 (2.66-2.62)
Clashscore	190562	2097 (2.66-2.62)
Ramachandran outliers	187476	2066 (2.66-2.62)
Sidechain outliers	187428	2066 (2.66-2.62)
RSRZ outliers	180081	2052 (2.66-2.62)

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	694	 79% 18% .
1	B	694	 80% 18% .

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 10713 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 Ornithine decarboxylase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	694	Total	C	N	O	S	0	0	0
			5349	3427	915	999	8			
1	B	694	Total	C	N	O	S	0	0	0
			5349	3427	915	999	8			

There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	17	ASP	ASN	conflict	UNP A0AAN4BJQ6
A	66	TYR	HIS	conflict	UNP A0AAN4BJQ6
A	102	GLU	GLY	conflict	UNP A0AAN4BJQ6
A	216	PHE	HIS	engineered mutation	UNP A0AAN4BJQ6
A	455	VAL	ALA	conflict	UNP A0AAN4BJQ6
A	620	ASP	GLU	conflict	UNP A0AAN4BJQ6
A	654	GLU	PRO	conflict	UNP A0AAN4BJQ6
A	695	GLY	-	expression tag	UNP A0AAN4BJQ6
B	17	ASP	ASN	conflict	UNP A0AAN4BJQ6
B	66	TYR	HIS	conflict	UNP A0AAN4BJQ6
B	102	GLU	GLY	conflict	UNP A0AAN4BJQ6
B	216	PHE	HIS	engineered mutation	UNP A0AAN4BJQ6
B	455	VAL	ALA	conflict	UNP A0AAN4BJQ6
B	620	ASP	GLU	conflict	UNP A0AAN4BJQ6
B	654	GLU	PRO	conflict	UNP A0AAN4BJQ6
B	695	GLY	-	expression tag	UNP A0AAN4BJQ6

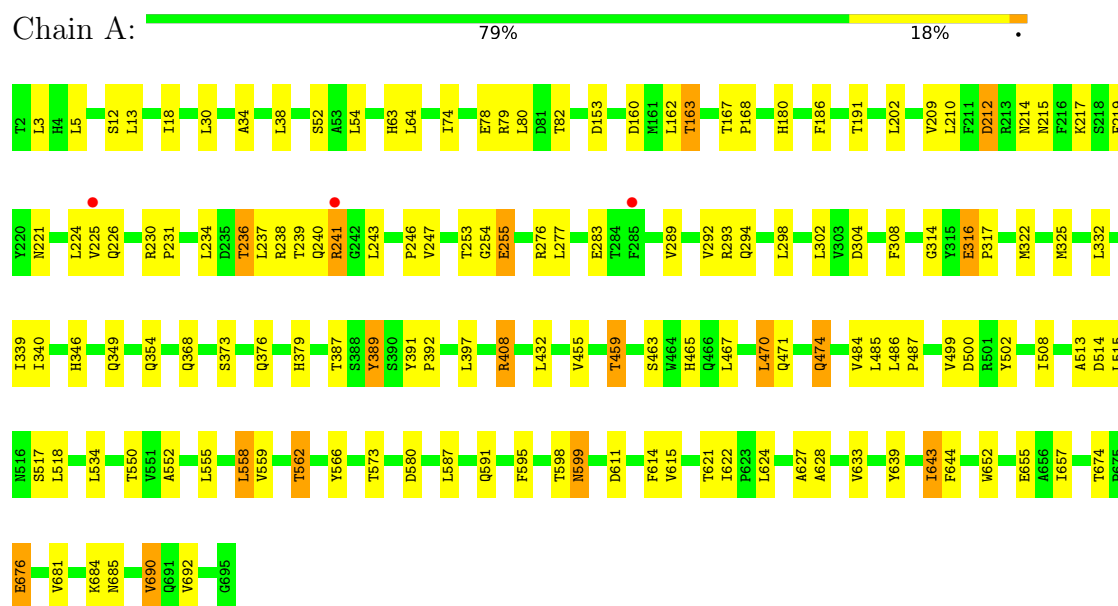
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	8	Total	O	0	0
			8	8		
2	B	7	Total	O	0	0
			7	7		

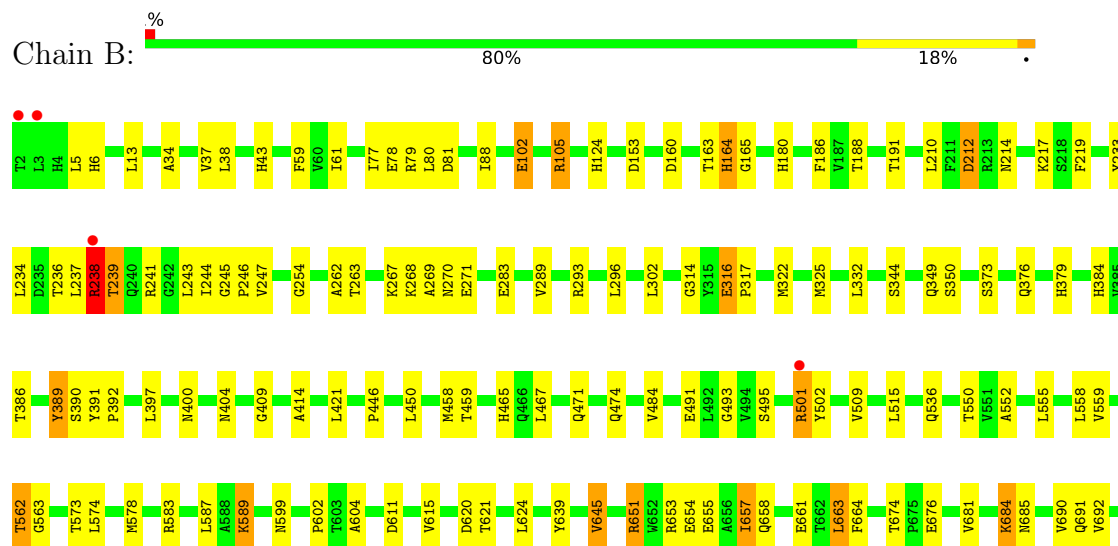
3 Residue-property plots

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: Ornithine decarboxylase



• Molecule 1: Ornithine decarboxylase



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4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	73.01Å 73.12Å 78.78Å 107.19° 114.42° 100.24°	Depositor
Resolution (Å)	44.54 – 2.64 44.54 – 2.64	Depositor EDS
% Data completeness (in resolution range)	97.7 (44.54-2.64) 97.7 (44.54-2.64)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.38 (at 2.65Å)	Xtriage
Refinement program	REFMAC 5.8.0352	Depositor
R, R_{free}	0.168 , 0.242 0.181 , 0.247	Depositor DCC
R_{free} test set	1917 reflections (4.92%)	wwPDB-VP
Wilson B-factor (Å ²)	42.2	Xtriage
Anisotropy	0.055	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 29.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.014 for k,h,-h-k-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	10713	wwPDB-VP
Average B, all atoms (Å ²)	44.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.92% 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

5.1 Standard geometry

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	0.69	4/5488 (0.1%)	1.07	7/7500 (0.1%)
1	B	0.69	5/5488 (0.1%)	1.07	5/7500 (0.1%)
All	All	0.69	9/10976 (0.1%)	1.07	12/15000 (0.1%)

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
1	B	0	1
All	All	0	2

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	379	HIS	CE1-NE2	6.78	1.39	1.32
1	A	346	HIS	CE1-NE2	6.62	1.39	1.32
1	A	180	HIS	CE1-NE2	6.04	1.38	1.32
1	B	379	HIS	CE1-NE2	5.98	1.38	1.32
1	B	180	HIS	CE1-NE2	5.70	1.38	1.32
1	B	124	HIS	CE1-NE2	5.56	1.38	1.32
1	A	63	HIS	CE1-NE2	5.51	1.38	1.32
1	B	384	HIS	CE1-NE2	5.49	1.38	1.32
1	B	6	HIS	CE1-NE2	5.28	1.37	1.32

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	674	THR	CB-CA-C	10.28	120.81	110.33
1	B	674	THR	CB-CA-C	10.01	120.54	110.33
1	B	645	VAL	N-CA-C	-6.61	106.35	112.96

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	231	PRO	N-CA-C	6.41	124.31	111.69
1	A	212	ASP	CA-CB-CG	5.89	118.49	112.60
1	A	163	THR	N-CA-C	-5.86	100.95	110.20
1	A	580	ASP	CA-CB-CG	5.51	118.11	112.60
1	A	153	ASP	CA-CB-CG	5.51	118.11	112.60
1	A	255	GLU	CB-CG-CD	5.46	121.88	112.60
1	B	43	HIS	CA-CB-CG	5.42	119.22	113.80
1	B	153	ASP	CA-CB-CG	5.20	117.80	112.60
1	B	212	ASP	CA-CB-CG	5.08	117.68	112.60

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	230	ARG	Sidechain
1	B	238	ARG	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	5349	0	5285	75	0
1	B	5349	0	5285	77	0
2	A	8	0	0	0	0
2	B	7	0	0	0	0
All	All	10713	0	10570	148	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 7.

All (148) 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:239:THR:HG22	1:A:243:LEU:H	1.38	0.89
1:A:471:GLN:H	1:A:474:GLN:HE21	1.19	0.88

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:465:HIS:HD2	1:A:467:LEU:H	1.25	0.82
1:A:237:LEU:HD12	1:A:246:PRO:HG2	1.64	0.77
1:A:246:PRO:HG3	1:A:470:LEU:HD11	1.68	0.75
1:B:237:LEU:HD12	1:B:246:PRO:HG2	1.69	0.74
1:B:465:HIS:HD2	1:B:467:LEU:H	1.35	0.74
1:B:624:LEU:HD23	1:B:661:GLU:HG3	1.70	0.74
1:B:233:TYR:H	1:B:599:ASN:HD21	1.36	0.73
1:B:446:PRO:O	1:B:450:LEU:HD23	1.89	0.72
1:A:550:THR:HG22	1:A:552:ALA:H	1.55	0.71
1:B:550:THR:HG22	1:B:552:ALA:H	1.56	0.70
1:B:574:LEU:HG	1:B:578:MET:HE2	1.76	0.67
1:A:239:THR:HG23	1:A:241:ARG:H	1.59	0.67
1:B:160:ASP:OD2	1:B:163:THR:HB	1.95	0.67
1:A:558:LEU:O	1:A:562:THR:HB	1.95	0.66
1:A:598:THR:O	1:A:599:ASN:HB2	1.95	0.66
1:B:559:VAL:HA	1:B:562:THR:HG22	1.78	0.66
1:B:37:VAL:HG11	1:B:80:LEU:HD21	1.78	0.64
1:A:432:LEU:O	1:A:487:PRO:HD3	1.99	0.62
1:A:502:TYR:HA	1:A:555:LEU:HD21	1.82	0.61
1:A:160:ASP:OD2	1:A:163:THR:HB	2.00	0.60
1:B:163:THR:O	1:B:164:HIS:C	2.45	0.60
1:A:515:LEU:HD11	1:A:587:LEU:HD23	1.84	0.59
1:A:643:ILE:HD12	1:A:644:PHE:N	2.19	0.58
1:A:238:ARG:HE	1:A:591:GLN:NE2	2.02	0.58
1:B:676:GLU:OE2	1:B:676:GLU:HA	2.04	0.58
1:B:624:LEU:HG	1:B:657:ILE:HG23	1.86	0.57
1:B:404:ASN:O	1:B:409:GLY:HA3	2.05	0.57
1:B:373:SER:H	1:B:376:GLN:HE21	1.50	0.57
1:B:191:THR:HG23	1:B:344:SER:OG	2.04	0.57
1:A:459:THR:O	1:A:465:HIS:HE1	1.88	0.56
1:B:160:ASP:HB3	1:B:165:GLY:HA2	1.87	0.56
1:A:167:THR:OG1	1:A:168:PRO:HD3	2.06	0.56
1:B:373:SER:H	1:B:376:GLN:NE2	2.03	0.55
1:B:502:TYR:HA	1:B:555:LEU:HD21	1.89	0.55
1:A:246:PRO:CG	1:A:470:LEU:HD11	2.37	0.54
1:A:254:GLY:HA2	1:A:302:LEU:HD11	1.88	0.54
1:A:221:ASN:HA	1:A:225:VAL:HB	1.90	0.54
1:B:515:LEU:HD11	1:B:587:LEU:HD23	1.88	0.53
1:A:253:THR:HG22	1:A:255:GLU:H	1.72	0.53
1:A:236:THR:HG23	1:A:595:PHE:HB3	1.90	0.53
1:B:238:ARG:HH21	1:B:245:GLY:H	1.57	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:484:VAL:HG22	1:A:518:LEU:HB2	1.91	0.52
1:A:598:THR:O	1:A:599:ASN:CB	2.57	0.52
1:B:655:GLU:H	1:B:655:GLU:CD	2.17	0.52
1:B:471:GLN:H	1:B:474:GLN:HE21	1.58	0.51
1:B:493:GLY:HA3	1:B:583:ARG:HH11	1.76	0.51
1:B:639:TYR:CD1	1:B:676:GLU:HB3	2.45	0.51
1:A:212:ASP:OD1	1:A:214:ASN:HB2	2.11	0.51
1:B:212:ASP:OD1	1:B:214:ASN:HB2	2.11	0.51
1:A:322:MET:HE3	1:A:325:MET:HE2	1.94	0.50
1:A:354:GLN:HG2	1:B:386:THR:HG21	1.93	0.50
1:A:465:HIS:CD2	1:A:467:LEU:H	2.17	0.50
1:A:54:LEU:CD2	1:A:408:ARG:HD2	2.42	0.50
1:A:471:GLN:H	1:A:474:GLN:NE2	2.00	0.50
1:B:684:LYS:HE3	1:B:684:LYS:HA	1.93	0.49
1:B:391:TYR:N	1:B:392:PRO:CD	2.75	0.49
1:A:484:VAL:CG2	1:A:518:LEU:HB2	2.43	0.49
1:A:215:ASN:O	1:A:643:ILE:HD13	2.12	0.49
1:B:79:ARG:HG2	1:B:81:ASP:OD1	2.13	0.49
1:A:162:LEU:HD11	1:A:387:THR:HA	1.95	0.49
1:B:59:PHE:CD2	1:B:88:ILE:HG12	2.47	0.49
1:B:263:THR:HG22	1:B:263:THR:O	2.13	0.49
1:A:354:GLN:HE22	1:B:188:THR:CB	2.26	0.48
1:B:38:LEU:C	1:B:38:LEU:HD12	2.38	0.48
1:A:373:SER:H	1:A:376:GLN:HE21	1.61	0.48
1:B:13:LEU:HD23	1:B:80:LEU:HD11	1.94	0.48
1:B:322:MET:HE3	1:B:325:MET:HE2	1.96	0.48
1:B:552:ALA:HA	1:B:559:VAL:HG21	1.94	0.48
1:B:459:THR:O	1:B:465:HIS:HE1	1.97	0.48
1:B:254:GLY:HA2	1:B:302:LEU:HD11	1.94	0.47
1:A:186:PHE:HB3	1:A:397:LEU:HD13	1.96	0.47
1:A:294:GLN:O	1:A:298:LEU:HG	2.14	0.47
1:A:373:SER:H	1:A:376:GLN:NE2	2.12	0.47
1:A:354:GLN:HE21	1:B:390:SER:H	1.62	0.47
1:B:191:THR:CG2	1:B:344:SER:OG	2.63	0.47
1:B:501:ARG:HD2	1:B:558:LEU:HD13	1.97	0.47
1:B:414:ALA:HA	1:B:450:LEU:HD11	1.97	0.46
1:B:186:PHE:HB3	1:B:397:LEU:HD13	1.96	0.46
1:B:654:GLU:O	1:B:658:GLN:HG3	2.14	0.46
1:A:5:LEU:HB2	1:A:34:ALA:HB2	1.97	0.46
1:B:316:GLU:N	1:B:317:PRO:CD	2.79	0.46
1:A:202:LEU:HD11	1:A:340:ILE:HD12	1.97	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:64:LEU:HD22	1:A:74:ILE:HG21	1.98	0.45
1:A:391:TYR:N	1:A:392:PRO:CD	2.79	0.45
1:B:268:LYS:O	1:B:271:GLU:HG2	2.17	0.45
1:B:314:GLY:HA3	1:B:349:GLN:NE2	2.31	0.45
1:B:501:ARG:HD2	1:B:558:LEU:HB2	1.98	0.45
1:A:292:VAL:HG23	1:A:308:PHE:CZ	2.52	0.45
1:A:514:ASP:HB2	1:A:517:SER:OG	2.17	0.45
1:A:550:THR:HG22	1:A:552:ALA:N	2.28	0.44
1:B:102:GLU:HB3	1:B:105:ARG:CZ	2.46	0.44
1:A:316:GLU:N	1:A:317:PRO:CD	2.80	0.44
1:A:562:THR:HG21	1:A:566:TYR:CD2	2.52	0.44
1:B:458:MET:HE2	1:B:465:HIS:CB	2.46	0.44
1:A:314:GLY:HA3	1:A:349:GLN:NE2	2.33	0.44
1:B:611:ASP:O	1:B:615:VAL:HG23	2.17	0.44
1:A:18:ILE:HG22	1:A:80:LEU:HD13	2.00	0.44
1:B:238:ARG:HH21	1:B:245:GLY:N	2.16	0.44
1:B:458:MET:HE2	1:B:465:HIS:HB2	2.00	0.44
1:A:54:LEU:HD21	1:A:408:ARG:HD2	1.99	0.44
1:A:643:ILE:HD12	1:A:644:PHE:H	1.83	0.44
1:B:624:LEU:CD2	1:B:661:GLU:HG3	2.45	0.44
1:A:500:ASP:HB2	1:A:513:ALA:HB2	2.01	0.43
1:B:550:THR:HG22	1:B:552:ALA:N	2.30	0.43
1:A:508:ILE:HG21	1:A:534:LEU:HD13	1.99	0.43
1:A:655:GLU:OE2	1:A:655:GLU:N	2.45	0.43
1:A:559:VAL:HA	1:A:562:THR:HG22	2.01	0.43
1:B:210:LEU:HG	1:B:234:LEU:HD11	2.00	0.43
1:A:167:THR:N	1:A:168:PRO:CD	2.82	0.43
1:A:239:THR:HG23	1:A:241:ARG:N	2.31	0.43
1:B:471:GLN:H	1:B:474:GLN:NE2	2.17	0.43
1:B:589:LYS:HE3	1:B:589:LYS:HB3	1.37	0.43
1:A:552:ALA:HA	1:A:559:VAL:HG21	2.01	0.43
1:B:262:ALA:HA	1:B:269:ALA:HB2	2.00	0.43
1:B:350:SER:HB2	1:B:400:ASN:OD1	2.18	0.43
1:B:604:ALA:HB2	1:B:651:ARG:CZ	2.49	0.42
1:B:316:GLU:H	1:B:316:GLU:CD	2.27	0.42
1:B:61:ILE:HG12	1:B:77:ILE:HG13	2.01	0.42
1:B:293:ARG:NH2	1:B:296:LEU:HD23	2.35	0.42
1:B:389:TYR:CD1	1:B:389:TYR:N	2.88	0.42
1:A:241:ARG:HH22	1:A:485:LEU:HD22	1.84	0.42
1:A:627:ALA:O	1:A:628:ALA:C	2.63	0.42
1:B:493:GLY:HA3	1:B:583:ARG:NH1	2.34	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:654:GLU:HA	1:B:657:ILE:HG13	2.02	0.42
1:A:652:TRP:CZ2	1:A:690:VAL:HG21	2.54	0.42
1:B:602:PRO:O	1:B:651:ARG:HG3	2.19	0.42
1:A:38:LEU:C	1:A:38:LEU:HD12	2.44	0.42
1:A:611:ASP:O	1:A:615:VAL:HG23	2.19	0.41
1:B:236:THR:HB	1:B:238:ARG:HH22	1.86	0.41
1:B:239:THR:HG22	1:B:241:ARG:H	1.84	0.41
1:B:501:ARG:CD	1:B:558:LEU:HB2	2.50	0.41
1:B:238:ARG:NE	1:B:244:ILE:HA	2.36	0.41
1:A:210:LEU:HG	1:A:234:LEU:HD11	2.03	0.41
1:A:316:GLU:H	1:A:316:GLU:CD	2.28	0.41
1:A:639:TYR:HB2	1:A:676:GLU:HG2	2.03	0.41
1:B:270:ASN:HD22	1:B:270:ASN:HA	1.61	0.41
1:A:276:ARG:O	1:A:304:ASP:HB2	2.21	0.41
1:A:389:TYR:CD1	1:A:389:TYR:N	2.89	0.41
1:A:486:LEU:HD21	1:A:499:VAL:HG21	2.03	0.41
1:B:663:LEU:HA	1:B:663:LEU:HD12	1.76	0.41
1:A:78:GLU:HG2	1:A:79:ARG:HG3	2.02	0.41
1:A:209:VAL:HG12	1:A:277:LEU:HB3	2.03	0.41
1:A:614:PHE:HB2	1:A:633:VAL:HB	2.03	0.40
1:B:5:LEU:HB2	1:B:34:ALA:HB2	2.02	0.40
1:A:354:GLN:NE2	1:B:390:SER:H	2.18	0.40
1:A:30:LEU:HD13	1:A:52:SER:HB3	2.03	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	692/694 (100%)	665 (96%)	26 (4%)	1 (0%)	48 64
1	B	692/694 (100%)	671 (97%)	19 (3%)	2 (0%)	36 50

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
All	All	1384/1388 (100%)	1336 (96%)	45 (3%)	3 (0%)	43	58

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	164	HIS
1	A	599	ASN
1	B	563	GLY

5.3.2 Protein sidechains [i](#)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	556/556 (100%)	515 (93%)	41 (7%)	13	20
1	B	556/556 (100%)	517 (93%)	39 (7%)	14	23
All	All	1112/1112 (100%)	1032 (93%)	80 (7%)	13	21

All (80) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	3	LEU
1	A	12	SER
1	A	13	LEU
1	A	82	THR
1	A	191	THR
1	A	217	LYS
1	A	219	PHE
1	A	224	LEU
1	A	226	GLN
1	A	236	THR
1	A	240	GLN
1	A	241	ARG
1	A	247	VAL
1	A	283	GLU
1	A	289	VAL

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Mol	Chain	Res	Type
1	A	293	ARG
1	A	316	GLU
1	A	332	LEU
1	A	339	ILE
1	A	368	GLN
1	A	389	TYR
1	A	408	ARG
1	A	455	VAL
1	A	459	THR
1	A	463	SER
1	A	470	LEU
1	A	474	GLN
1	A	558	LEU
1	A	562	THR
1	A	573	THR
1	A	621	THR
1	A	622	ILE
1	A	624	LEU
1	A	643	ILE
1	A	657	ILE
1	A	676	GLU
1	A	681	VAL
1	A	684	LYS
1	A	685	ASN
1	A	690	VAL
1	A	692	VAL
1	B	78	GLU
1	B	102	GLU
1	B	105	ARG
1	B	217	LYS
1	B	219	PHE
1	B	238	ARG
1	B	239	THR
1	B	243	LEU
1	B	247	VAL
1	B	267	LYS
1	B	283	GLU
1	B	289	VAL
1	B	316	GLU
1	B	332	LEU
1	B	389	TYR
1	B	421	LEU

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Mol	Chain	Res	Type
1	B	484	VAL
1	B	491	GLU
1	B	495	SER
1	B	501	ARG
1	B	509	VAL
1	B	536	GLN
1	B	562	THR
1	B	573	THR
1	B	589	LYS
1	B	620	ASP
1	B	621	THR
1	B	645	VAL
1	B	651	ARG
1	B	653	ARG
1	B	657	ILE
1	B	663	LEU
1	B	664	PHE
1	B	681	VAL
1	B	684	LYS
1	B	685	ASN
1	B	690	VAL
1	B	691	GLN
1	B	692	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (34) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	4	HIS
1	A	50	ASN
1	A	189	ASN
1	A	215	ASN
1	A	226	GLN
1	A	270	ASN
1	A	354	GLN
1	A	368	GLN
1	A	376	GLN
1	A	465	HIS
1	A	471	GLN
1	A	474	GLN
1	A	536	GLN
1	A	591	GLN
1	A	600	ASN

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Mol	Chain	Res	Type
1	A	608	GLN
1	B	28	HIS
1	B	44	GLN
1	B	87	GLN
1	B	189	ASN
1	B	215	ASN
1	B	226	GLN
1	B	270	ASN
1	B	343	GLN
1	B	368	GLN
1	B	376	GLN
1	B	405	GLN
1	B	465	HIS
1	B	471	GLN
1	B	474	GLN
1	B	540	GLN
1	B	591	GLN
1	B	599	ASN
1	B	691	GLN

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 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 ⓘ

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	694/694 (100%)	-0.22	3 (0%) 88 87	22, 43, 68, 105	0
1	B	694/694 (100%)	-0.22	4 (0%) 85 83	21, 40, 73, 107	0
All	All	1388/1388 (100%)	-0.22	7 (0%) 87 85	21, 42, 70, 107	0

All (7) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	225	VAL	4.9
1	A	241	ARG	3.7
1	B	2	THR	2.5
1	B	501	ARG	2.5
1	B	3	LEU	2.4
1	B	238	ARG	2.2
1	A	285	PHE	2.2

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](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.