



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 04:18 AM UTC

PDB ID : 9LMA / pdb_00009lma
Title : Crystal structure of Ornithine decarboxylase H346A mutant without PLP
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Deposited on : 2025-01-18
Resolution : 1.74 Å(reported)

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

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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

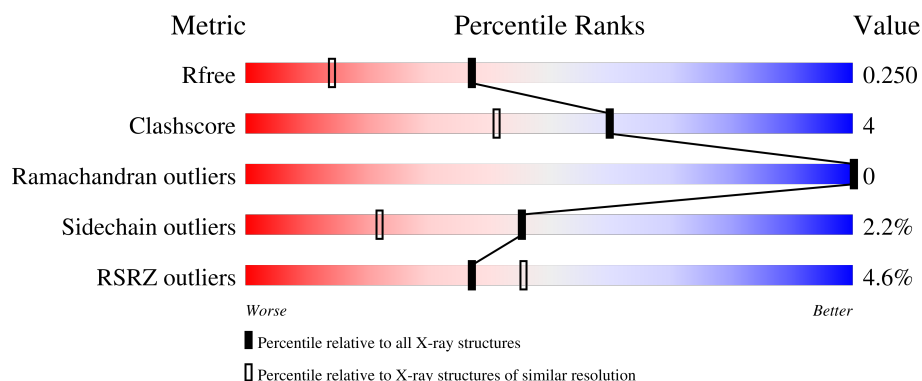
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.74 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	1187 (1.74-1.74)
Clashscore	190562	1207 (1.74-1.74)
Ramachandran outliers	187476	1200 (1.74-1.74)
Sidechain outliers	187428	1200 (1.74-1.74)
RSRZ outliers	180081	1188 (1.74-1.74)

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	693	4% 91% 9%
1	B	693	5% 86% 12% .

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 11177 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	693	Total	C	N	O	S	0	0	0
			5329	3413	913	995	8			
1	B	692	Total	C	N	O	S	0	0	0
			5322	3409	912	993	8			

There are 18 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	16	ALA	PRO	conflict	UNP A0AAN4BJQ6
A	17	ALA	ASN	conflict	UNP A0AAN4BJQ6
A	66	TYR	HIS	conflict	UNP A0AAN4BJQ6
A	102	GLU	GLY	conflict	UNP A0AAN4BJQ6
A	346	ALA	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	684	ALA	LYS	conflict	UNP A0AAN4BJQ6
B	16	ALA	PRO	conflict	UNP A0AAN4BJQ6
B	17	ALA	ASN	conflict	UNP A0AAN4BJQ6
B	66	TYR	HIS	conflict	UNP A0AAN4BJQ6
B	102	GLU	GLY	conflict	UNP A0AAN4BJQ6
B	346	ALA	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	684	ALA	LYS	conflict	UNP A0AAN4BJQ6

- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	258	Total	O	0	0
			258	258		

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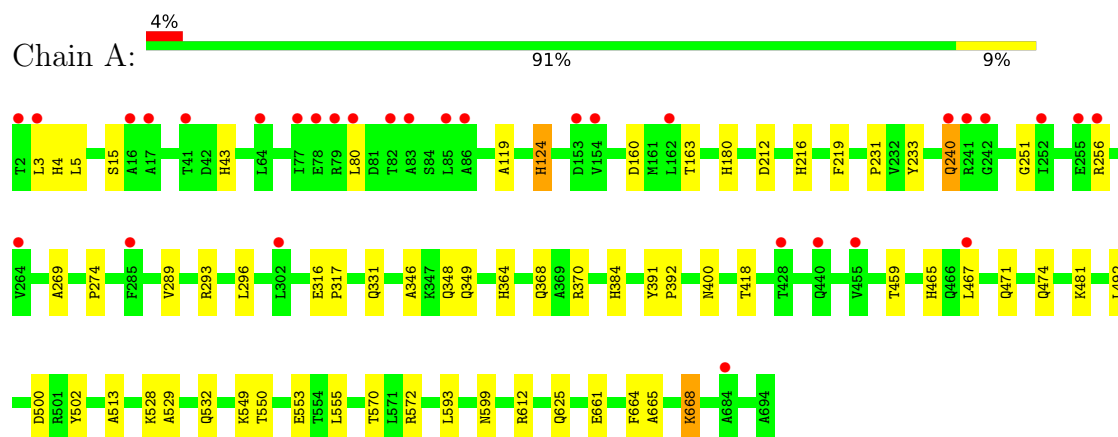
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	268	Total 268	O 268	0	0

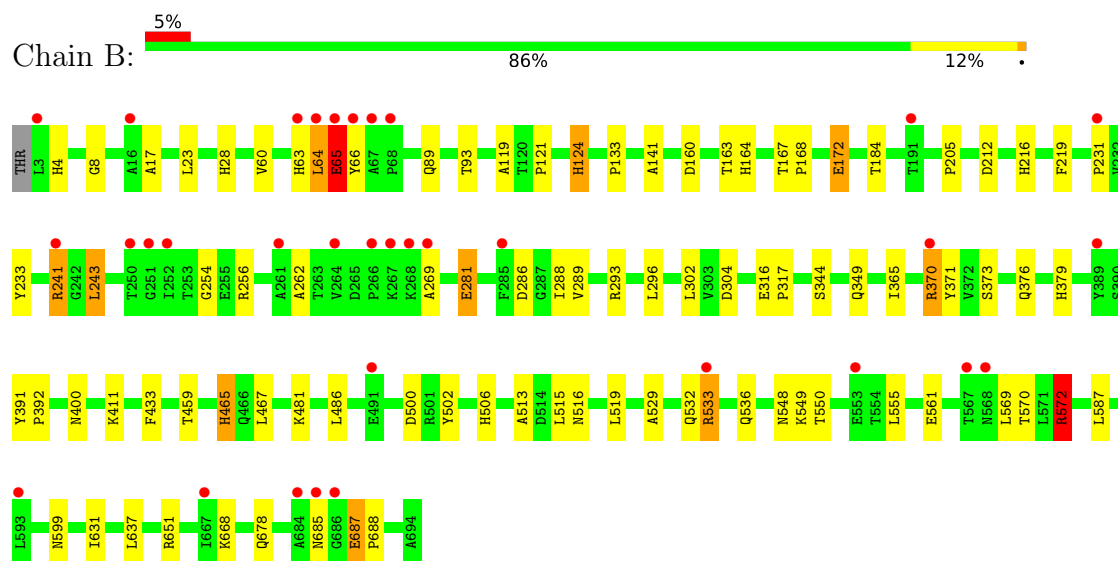
3 Residue-property plots [i](#)

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: Ornithine decarboxylase



• Molecule 1: Ornithine decarboxylase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	72.00Å 72.31Å 78.01Å 107.60° 114.56° 100.37°	Depositor
Resolution (Å)	29.29 – 1.74 29.29 – 1.74	Depositor EDS
% Data completeness (in resolution range)	96.3 (29.29-1.74) 96.3 (29.29-1.74)	Depositor EDS
R_{merge}	0.24	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.72 (at 1.75Å)	Xtriage
Refinement program	REFMAC 5.8.0352	Depositor
R, R_{free}	0.202 , 0.244 0.211 , 0.250	Depositor DCC
R_{free} test set	6230 reflections (4.78%)	wwPDB-VP
Wilson B-factor (Å ²)	22.4	Xtriage
Anisotropy	0.050	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 21.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.021 for k,h,-h-k-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	11177	wwPDB-VP
Average B, all atoms (Å ²)	26.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.65% 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	0.73	7/5466 (0.1%)	1.00	4/7473 (0.1%)
1	B	0.75	10/5459 (0.2%)	1.01	3/7463 (0.0%)
All	All	0.74	17/10925 (0.2%)	1.01	7/14936 (0.0%)

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	B	0	1

All (17) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	216	HIS	CE1-NE2	6.69	1.39	1.32
1	B	379	HIS	CE1-NE2	6.54	1.39	1.32
1	A	364	HIS	CE1-NE2	6.51	1.39	1.32
1	A	216	HIS	CE1-NE2	6.30	1.38	1.32
1	B	465	HIS	CE1-NE2	5.91	1.38	1.32
1	A	124	HIS	CE1-NE2	5.90	1.38	1.32
1	B	164	HIS	CE1-NE2	5.72	1.38	1.32
1	B	506	HIS	CE1-NE2	5.68	1.38	1.32
1	A	384	HIS	CE1-NE2	5.50	1.38	1.32
1	A	43	HIS	CE1-NE2	5.30	1.37	1.32
1	A	180	HIS	CE1-NE2	5.29	1.37	1.32
1	B	572	ARG	C-O	-5.24	1.18	1.24
1	B	124	HIS	CE1-NE2	5.23	1.37	1.32
1	A	4	HIS	CE1-NE2	5.19	1.37	1.32
1	B	4	HIS	CE1-NE2	5.19	1.37	1.32
1	B	28	HIS	CE1-NE2	5.17	1.37	1.32
1	B	344	SER	CA-CB	-5.10	1.46	1.53

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	212	ASP	CA-CB-CG	7.06	119.66	112.60
1	B	212	ASP	CA-CB-CG	7.00	119.60	112.60
1	A	231	PRO	N-CA-C	6.39	124.27	111.69
1	B	231	PRO	N-CA-C	5.98	124.79	112.47
1	A	231	PRO	CB-CA-C	-5.58	101.38	110.25
1	A	418	THR	CA-CB-OG1	-5.49	101.36	109.60
1	B	65	GLU	N-CA-C	-5.19	102.17	109.96

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	241	ARG	Sidechain

5.2 Too-close contacts ⓘ

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	5329	0	5269	42	0
1	B	5322	0	5262	55	0
2	A	258	0	0	1	0
2	B	268	0	0	3	0
All	All	11177	0	10531	92	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 (92) 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:B:17:ALA:HB2	2:B:944:HOH:O	1.67	0.94
1:A:233:TYR:H	1:A:599:ASN:HD21	1.24	0.84
1:B:465:HIS:HD2	1:B:467:LEU:H	1.26	0.83
1:A:465:HIS:HD2	1:A:467:LEU:H	1.27	0.82
1:A:625:GLN:HG2	1:A:661:GLU:OE2	1.81	0.81

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:233:TYR:H	1:B:599:ASN:HD21	1.29	0.81
1:A:251:GLY:O	1:A:256:ARG:HD2	1.83	0.79
1:B:529:ALA:HA	1:B:532:GLN:HE21	1.55	0.69
1:B:241:ARG:NH1	1:B:516:ASN:HB2	2.06	0.69
1:A:349:GLN:NE2	1:A:400:ASN:HD21	1.91	0.68
1:B:349:GLN:NE2	1:B:400:ASN:HD21	1.93	0.67
1:B:631:ILE:CD1	1:B:651:ARG:HG2	2.25	0.66
1:B:631:ILE:HD11	1:B:651:ARG:NH1	2.09	0.66
1:A:368:GLN:NE2	1:A:370:ARG:HE	1.96	0.64
1:B:288:ILE:HD13	1:B:467:LEU:HD11	1.79	0.63
1:A:348:GLN:HE22	1:A:481:LYS:HZ2	1.45	0.62
1:B:262:ALA:HB2	1:B:269:ALA:CB	2.31	0.61
1:A:348:GLN:HE22	1:A:481:LYS:NZ	2.01	0.58
1:B:548:ASN:HD21	1:B:572:ARG:HD3	1.70	0.57
1:A:471:GLN:H	1:A:474:GLN:HE21	1.51	0.56
1:B:502:TYR:HA	1:B:555:LEU:HD21	1.88	0.56
1:B:65:GLU:O	1:B:66:TYR:HB2	2.06	0.56
1:B:433:PHE:CE1	1:B:486:LEU:HD21	2.42	0.55
1:B:631:ILE:HD11	1:B:651:ARG:CZ	2.38	0.54
1:B:515:LEU:HD11	1:B:587:LEU:HD23	1.89	0.54
1:B:631:ILE:HD13	1:B:651:ARG:HG2	1.89	0.54
1:A:664:PHE:O	1:A:668:LYS:HG2	2.08	0.53
1:B:241:ARG:NE	2:B:703:HOH:O	2.41	0.53
1:B:550:THR:HA	1:B:570:THR:HA	1.90	0.53
1:A:593:LEU:C	1:A:593:LEU:HD23	2.34	0.52
1:B:529:ALA:HA	1:B:532:GLN:NE2	2.25	0.52
1:A:550:THR:HA	1:A:570:THR:HA	1.91	0.52
1:B:316:GLU:N	1:B:317:PRO:CD	2.74	0.51
1:A:3:LEU:HD21	1:B:141:ALA:HA	1.91	0.51
1:B:365:ILE:HG13	1:B:371:TYR:HD1	1.76	0.51
1:B:119:ALA:O	1:B:124:HIS:HE1	1.94	0.51
1:A:502:TYR:HA	1:A:555:LEU:HD21	1.91	0.51
1:A:316:GLU:N	1:A:317:PRO:CD	2.73	0.51
1:B:373:SER:H	1:B:376:GLN:HE21	1.57	0.51
1:B:241:ARG:HH12	1:B:516:ASN:HB2	1.73	0.50
1:B:293:ARG:NH1	1:B:296:LEU:HD23	2.27	0.50
1:A:119:ALA:O	1:A:124:HIS:HE1	1.95	0.50
1:B:533:ARG:NH1	1:B:536:GLN:HB2	2.27	0.49
1:A:348:GLN:NE2	1:A:481:LYS:NZ	2.61	0.49
1:B:373:SER:H	1:B:376:GLN:NE2	2.10	0.49
1:A:331:GLN:NE2	2:A:703:HOH:O	2.29	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:368:GLN:HE22	1:A:370:ARG:HH21	1.62	0.48
1:A:269:ALA:HA	1:A:274:PRO:HB3	1.96	0.47
1:B:160:ASP:OD2	1:B:163:THR:OG1	2.24	0.47
1:A:465:HIS:CD2	1:A:467:LEU:H	2.18	0.47
1:B:65:GLU:HG2	1:B:66:TYR:CD1	2.49	0.47
1:A:5:LEU:HD11	1:B:133:PRO:HB3	1.96	0.47
1:A:593:LEU:HD23	1:A:593:LEU:O	2.13	0.47
1:A:346:ALA:O	1:B:121:PRO:HA	2.15	0.47
1:B:281:GLU:O	1:B:281:GLU:HG2	2.15	0.47
1:B:481:LYS:HB3	1:B:519:LEU:HD11	1.97	0.46
1:B:500:ASP:HB2	1:B:513:ALA:HB2	1.98	0.46
1:A:465:HIS:CD2	1:A:467:LEU:HB2	2.51	0.45
1:B:687:GLU:HG2	1:B:688:PRO:HD2	1.98	0.45
1:A:80:LEU:N	1:A:80:LEU:HD12	2.31	0.45
1:A:549:LYS:HD2	1:A:553:GLU:OE1	2.16	0.45
1:B:548:ASN:ND2	1:B:572:ARG:HD3	2.32	0.45
1:A:528:LYS:O	1:A:532:GLN:HG3	2.17	0.45
1:A:529:ALA:HA	1:A:532:GLN:HE21	1.81	0.45
1:B:172:GLU:HG2	1:B:184:THR:O	2.17	0.45
1:A:549:LYS:HD2	1:A:553:GLU:CD	2.41	0.45
1:B:60:VAL:HG11	1:B:64:LEU:HD11	1.99	0.44
1:A:471:GLN:H	1:A:474:GLN:NE2	2.13	0.44
1:B:459:THR:O	1:B:465:HIS:HE1	2.01	0.44
1:A:665:ALA:HA	1:A:668:LYS:HE2	2.00	0.44
1:B:533:ARG:HH12	1:B:536:GLN:HB2	1.83	0.44
1:A:3:LEU:CD2	1:B:141:ALA:HA	2.47	0.43
1:A:459:THR:O	1:A:465:HIS:HE1	2.01	0.43
1:B:262:ALA:CA	1:B:269:ALA:HB2	2.49	0.43
1:A:293:ARG:NH1	1:A:296:LEU:HD23	2.33	0.43
1:A:492:LEU:HD13	1:A:572:ARG:NH2	2.34	0.43
1:B:241:ARG:CZ	2:B:703:HOH:O	2.65	0.43
1:B:167:THR:N	1:B:168:PRO:CD	2.82	0.43
1:B:243:LEU:HD13	1:B:286:ASP:HB3	1.99	0.43
1:B:304:ASP:HA	1:B:370:ARG:HH11	1.84	0.43
1:B:262:ALA:HB2	1:B:269:ALA:HB3	2.00	0.42
1:B:637:LEU:HD23	1:B:678:GLN:OE1	2.20	0.42
1:A:160:ASP:OD2	1:A:163:THR:OG1	2.27	0.42
1:A:612:ARG:NH2	1:B:205:PRO:HG2	2.35	0.42
1:B:89:GLN:O	1:B:93:THR:HG23	2.20	0.41
1:A:240:GLN:CD	1:A:240:GLN:H	2.27	0.41
1:B:8:GLY:HA2	1:B:23:LEU:O	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:500:ASP:HB2	1:A:513:ALA:HB2	2.02	0.41
1:B:254:GLY:HA2	1:B:302:LEU:HD11	2.02	0.41
1:A:233:TYR:N	1:A:599:ASN:HD21	2.05	0.40
1:A:391:TYR:N	1:A:392:PRO:CD	2.84	0.40
1:B:391:TYR:N	1:B:392:PRO:CD	2.84	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	691/693 (100%)	680 (98%)	11 (2%)	0	100	100
1	B	690/693 (100%)	674 (98%)	16 (2%)	0	100	100
All	All	1381/1386 (100%)	1354 (98%)	27 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	552/552 (100%)	547 (99%)	5 (1%)	70	59
1	B	551/552 (100%)	532 (97%)	19 (3%)	32	10
All	All	1103/1104 (100%)	1079 (98%)	24 (2%)	45	22

All (24) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	15	SER
1	A	219	PHE
1	A	240	GLN
1	A	289	VAL
1	A	668	LYS
1	B	63	HIS
1	B	64	LEU
1	B	65	GLU
1	B	172	GLU
1	B	219	PHE
1	B	243	LEU
1	B	256	ARG
1	B	281	GLU
1	B	289	VAL
1	B	370	ARG
1	B	411	LYS
1	B	533	ARG
1	B	549	LYS
1	B	561	GLU
1	B	569	LEU
1	B	572	ARG
1	B	668	LYS
1	B	685	ASN
1	B	687	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (38) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	4	HIS
1	A	6	HIS
1	A	28	HIS
1	A	124	HIS
1	A	139	HIS
1	A	195	ASN
1	A	215	ASN
1	A	226	GLN
1	A	259	GLN
1	A	348	GLN
1	A	349	GLN
1	A	357	GLN
1	A	368	GLN

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Mol	Chain	Res	Type
1	A	405	GLN
1	A	465	HIS
1	A	471	GLN
1	A	474	GLN
1	A	532	GLN
1	A	599	ASN
1	A	600	ASN
1	A	608	GLN
1	A	658	GLN
1	B	195	ASN
1	B	215	ASN
1	B	226	GLN
1	B	349	GLN
1	B	357	GLN
1	B	368	GLN
1	B	376	GLN
1	B	440	GLN
1	B	465	HIS
1	B	532	GLN
1	B	548	ASN
1	B	584	GLN
1	B	599	ASN
1	B	600	ASN
1	B	608	GLN
1	B	658	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

There are no ligands in this entry.

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	693/693 (100%)	0.39	31 (4%) 38 47	12, 25, 44, 76	0
1	B	692/693 (99%)	0.39	33 (4%) 35 43	11, 23, 48, 111	0
All	All	1385/1386 (99%)	0.39	64 (4%) 37 46	11, 24, 46, 111	0

All (64) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	66	TYR	5.0
1	B	67	ALA	4.9
1	B	684	ALA	4.7
1	A	302	LEU	4.2
1	B	16	ALA	4.1
1	B	64	LEU	4.0
1	B	231	PRO	3.8
1	B	269	ALA	3.8
1	B	68	PRO	3.7
1	A	80	LEU	3.7
1	A	684	ALA	3.6
1	A	17	ALA	3.5
1	B	593	LEU	3.5
1	A	82	THR	3.5
1	B	63	HIS	3.4
1	B	568	ASN	3.4
1	A	241	ARG	3.4
1	A	41	THR	3.3
1	A	64	LEU	3.3
1	A	162	LEU	3.2
1	B	3	LEU	3.1
1	B	241	ARG	3.0
1	B	370	ARG	2.9
1	A	2	THR	2.8

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Mol	Chain	Res	Type	RSRZ
1	B	533	ARG	2.8
1	B	285	PHE	2.8
1	A	285	PHE	2.8
1	B	266	PRO	2.8
1	B	267	LYS	2.7
1	B	250	THR	2.7
1	B	252	ILE	2.7
1	A	16	ALA	2.7
1	B	389	TYR	2.6
1	A	83	ALA	2.6
1	A	79	ARG	2.6
1	B	686	GLY	2.5
1	A	428	THR	2.5
1	A	240	GLN	2.5
1	A	242	GLY	2.5
1	B	264	VAL	2.5
1	A	78	GLU	2.5
1	A	85	LEU	2.4
1	A	255	GLU	2.4
1	A	467	LEU	2.4
1	A	77	ILE	2.3
1	B	251	GLY	2.3
1	A	440	GLN	2.3
1	A	252	ILE	2.2
1	A	153	ASP	2.2
1	B	685	ASN	2.2
1	B	65	GLU	2.2
1	A	3	LEU	2.2
1	B	667	ILE	2.2
1	B	567	THR	2.1
1	B	491	GLU	2.1
1	A	256	ARG	2.1
1	B	191	THR	2.1
1	B	268	LYS	2.1
1	A	86	ALA	2.0
1	B	553	GLU	2.0
1	B	261	ALA	2.0
1	A	154	VAL	2.0
1	A	264	VAL	2.0
1	A	455	VAL	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 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.