



Full wwPDB X-ray Structure Validation Report ⓘ

Oct 14, 2023 – 07:48 PM EDT

PDB ID : 7U1H
Title : Crystal structure of Lens culinaris vicilin
Authors : Robinson, K.A.; Bakestani, I.D.; Loewen, M.C.
Deposited on : 2022-02-21
Resolution : 2.50 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 : 4.02b-467
Xtriage (Phenix) : 1.13
EDS : 2.36
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.36

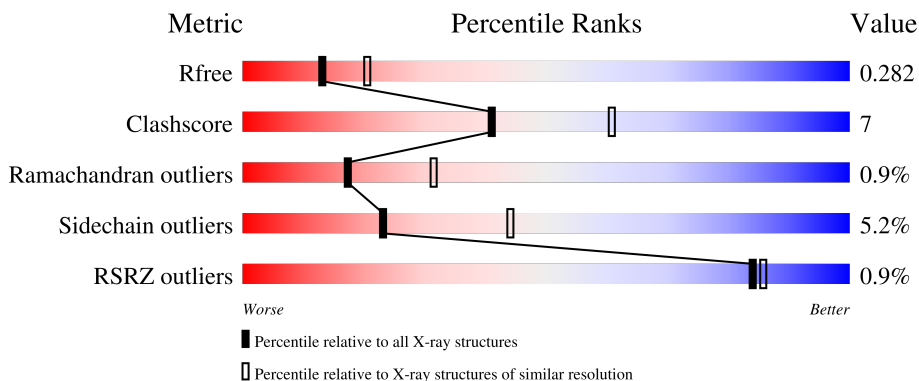
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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}	130704	4661 (2.50-2.50)
Clashscore	141614	5346 (2.50-2.50)
Ramachandran outliers	138981	5231 (2.50-2.50)
Sidechain outliers	138945	5233 (2.50-2.50)
RSRZ outliers	127900	4559 (2.50-2.50)

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	431	<div style="display: flex; align-items: center;"> <div style="width: 20px; height: 10px; background-color: red; margin-right: 5px;"></div> <div style="width: 100%; height: 20px; background: linear-gradient(to right, red 2%, orange 20%, yellow 36%, green 68%, grey 100%); position: relative;"> 68% 16% 15% </div> </div>
1	B	431	<div style="display: flex; align-items: center;"> <div style="width: 100%; height: 20px; background: linear-gradient(to right, green 65%, yellow 84%, orange 90%, grey 100%); position: relative;"> 65% 19% 15% </div> </div>
1	C	431	<div style="display: flex; align-items: center;"> <div style="width: 100%; height: 20px; background: linear-gradient(to right, green 68%, yellow 82%, orange 96%, grey 100%); position: relative;"> 68% 14% 16% </div> </div>

2 Entry composition i

There are 2 unique types of molecules in this entry. The entry contains 17579 atoms, of which 8775 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 Allergen Len c 1.0101.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	H	N	O			
1	A	365	5889	1858	2951	518	562	62	4	0
1	B	365	5857	1848	2930	516	563	61	4	0
1	C	363	5793	1832	2894	510	557	61	3	0

There are 57 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-17	MET	-	expression tag	UNP Q84UI1
A	-16	GLY	-	expression tag	UNP Q84UI1
A	-15	SER	-	expression tag	UNP Q84UI1
A	-14	SER	-	expression tag	UNP Q84UI1
A	-13	HIS	-	expression tag	UNP Q84UI1
A	-12	HIS	-	expression tag	UNP Q84UI1
A	-11	HIS	-	expression tag	UNP Q84UI1
A	-10	HIS	-	expression tag	UNP Q84UI1
A	-9	HIS	-	expression tag	UNP Q84UI1
A	-8	HIS	-	expression tag	UNP Q84UI1
A	-7	LEU	-	expression tag	UNP Q84UI1
A	-6	VAL	-	expression tag	UNP Q84UI1
A	-5	PRO	-	expression tag	UNP Q84UI1
A	-4	ARG	-	expression tag	UNP Q84UI1
A	-3	GLY	-	expression tag	UNP Q84UI1
A	-2	SER	-	expression tag	UNP Q84UI1
A	-1	HIS	-	expression tag	UNP Q84UI1
A	0	MET	-	expression tag	UNP Q84UI1
A	1	MET	-	expression tag	UNP Q84UI1
B	-17	MET	-	expression tag	UNP Q84UI1
B	-16	GLY	-	expression tag	UNP Q84UI1
B	-15	SER	-	expression tag	UNP Q84UI1
B	-14	SER	-	expression tag	UNP Q84UI1

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-13	HIS	-	expression tag	UNP Q84UI1
B	-12	HIS	-	expression tag	UNP Q84UI1
B	-11	HIS	-	expression tag	UNP Q84UI1
B	-10	HIS	-	expression tag	UNP Q84UI1
B	-9	HIS	-	expression tag	UNP Q84UI1
B	-8	HIS	-	expression tag	UNP Q84UI1
B	-7	LEU	-	expression tag	UNP Q84UI1
B	-6	VAL	-	expression tag	UNP Q84UI1
B	-5	PRO	-	expression tag	UNP Q84UI1
B	-4	ARG	-	expression tag	UNP Q84UI1
B	-3	GLY	-	expression tag	UNP Q84UI1
B	-2	SER	-	expression tag	UNP Q84UI1
B	-1	HIS	-	expression tag	UNP Q84UI1
B	0	MET	-	expression tag	UNP Q84UI1
B	1	MET	-	expression tag	UNP Q84UI1
C	-17	MET	-	expression tag	UNP Q84UI1
C	-16	GLY	-	expression tag	UNP Q84UI1
C	-15	SER	-	expression tag	UNP Q84UI1
C	-14	SER	-	expression tag	UNP Q84UI1
C	-13	HIS	-	expression tag	UNP Q84UI1
C	-12	HIS	-	expression tag	UNP Q84UI1
C	-11	HIS	-	expression tag	UNP Q84UI1
C	-10	HIS	-	expression tag	UNP Q84UI1
C	-9	HIS	-	expression tag	UNP Q84UI1
C	-8	HIS	-	expression tag	UNP Q84UI1
C	-7	LEU	-	expression tag	UNP Q84UI1
C	-6	VAL	-	expression tag	UNP Q84UI1
C	-5	PRO	-	expression tag	UNP Q84UI1
C	-4	ARG	-	expression tag	UNP Q84UI1
C	-3	GLY	-	expression tag	UNP Q84UI1
C	-2	SER	-	expression tag	UNP Q84UI1
C	-1	HIS	-	expression tag	UNP Q84UI1
C	0	MET	-	expression tag	UNP Q84UI1
C	1	MET	-	expression tag	UNP Q84UI1

- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	12	Total O 12 12	0	0
2	B	17	Total O 17 17	0	0

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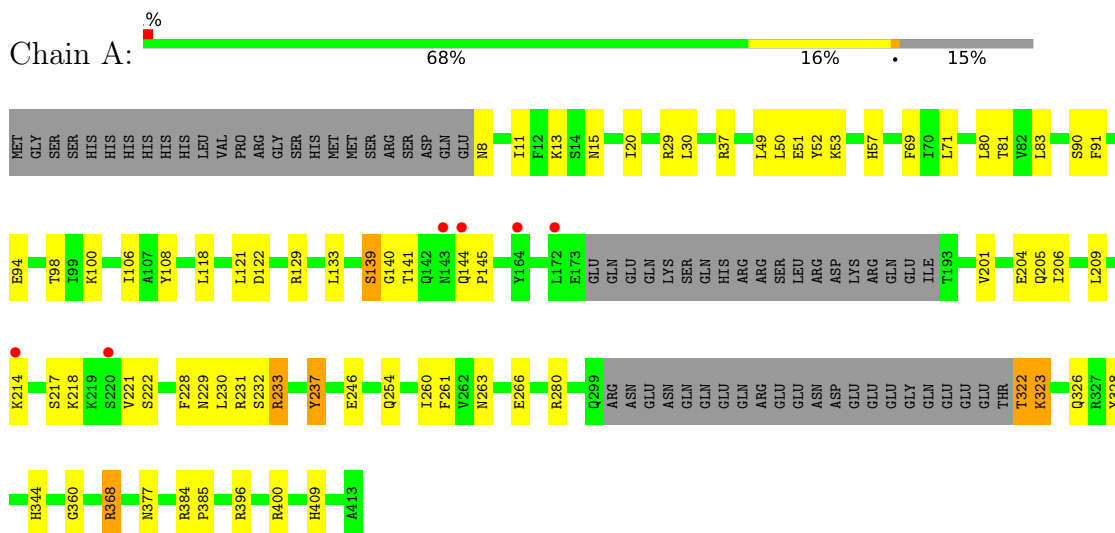
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	C	11	Total	O	0	0
			11	11		

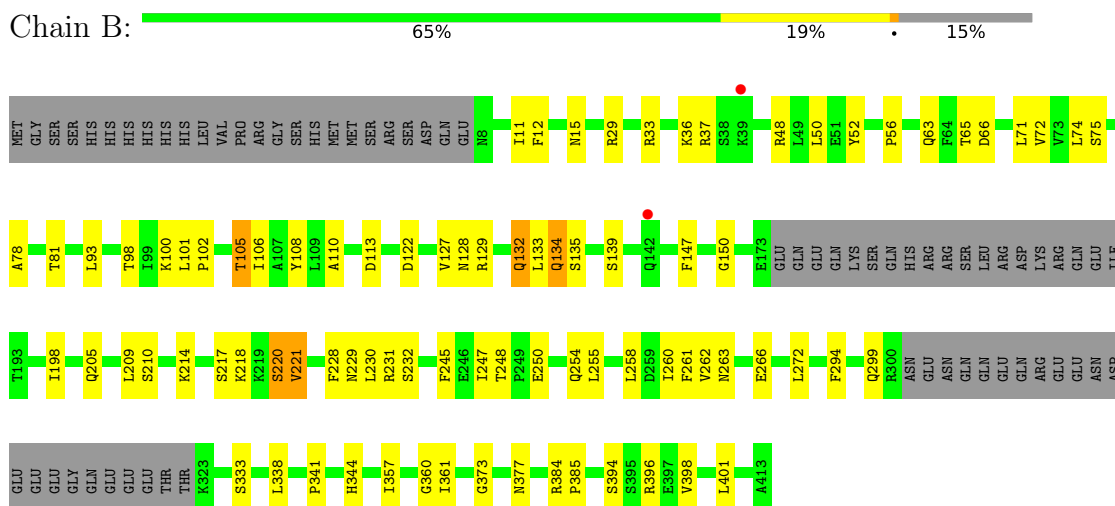
3 Residue-property plots [i](#)

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: Allergen Len c 1.0101

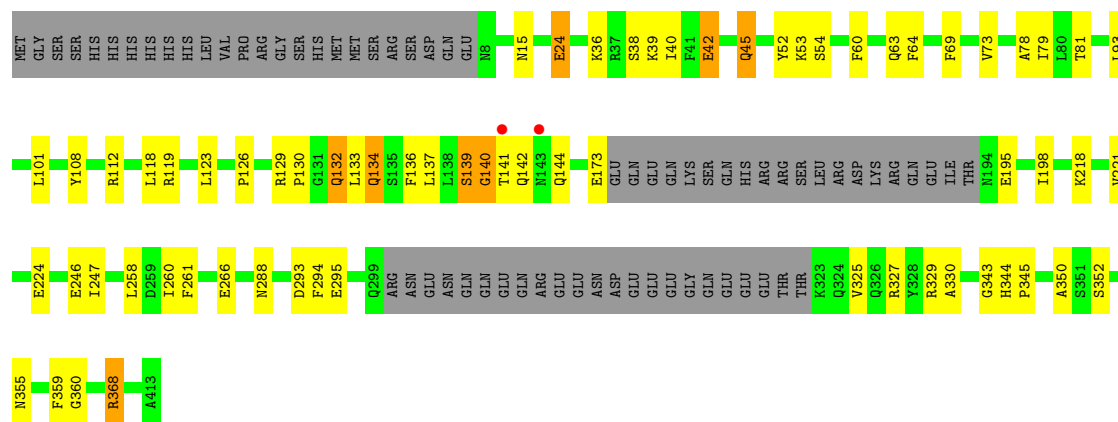


- Molecule 1: Allergen Len c 1.0101



- Molecule 1: Allergen Len c 1.0101





4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	86.44Å 92.50Å 143.74Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	77.91 – 2.50 77.79 – 2.50	Depositor EDS
% Data completeness (in resolution range)	98.2 (77.91-2.50) 97.9 (77.79-2.50)	Depositor EDS
R_{merge}	0.14	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.78 (at 2.51Å)	Xtrriage
Refinement program	REFMAC 5.8.0267	Depositor
R, R_{free}	0.199 , 0.284 0.202 , 0.282	Depositor DCC
R_{free} test set	2068 reflections (5.20%)	wwPDB-VP
Wilson B-factor (Å ²)	32.8	Xtrriage
Anisotropy	0.980	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.41 , 28.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	17579	wwPDB-VP
Average B, all atoms (Å ²)	38.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.61% 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	0.70	0/3000	0.93	3/4052 (0.1%)
1	B	0.71	0/2988	0.93	0/4037
1	C	0.70	0/2955	0.95	2/3994 (0.1%)
All	All	0.70	0/8943	0.94	5/12083 (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

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	368	ARG	NE-CZ-NH2	-7.20	116.70	120.30
1	A	368	ARG	NE-CZ-NH1	7.16	123.88	120.30
1	A	280	ARG	NE-CZ-NH1	6.31	123.45	120.30
1	C	368	ARG	NE-CZ-NH2	-6.29	117.15	120.30
1	C	368	ARG	NE-CZ-NH1	6.02	123.31	120.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	218	LYS	Peptide

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	2938	2951	2944	41	2
1	B	2927	2930	2919	50	0
1	C	2899	2894	2875	36	2
2	A	12	0	0	1	0
2	B	17	0	0	0	0
2	C	11	0	0	0	0
All	All	8804	8775	8738	124	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 7.

All (124) 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:C:288:ASN:HD22	1:C:355[B]:ASN:HD22	1.24	0.85
1:B:71:LEU:HD12	1:B:122:ASP:OD1	1.79	0.83
1:C:129:ARG:HB2	1:C:132:GLN:HG3	1.61	0.82
1:A:20:ILE:HD11	1:A:30:LEU:HB2	1.61	0.82
1:B:258:LEU:HB3	1:B:260:ILE:HD12	1.69	0.74
1:B:71:LEU:O	1:B:98:THR:HA	1.89	0.72
1:B:29:ARG:O	1:B:50:LEU:HA	1.93	0.68
1:B:71:LEU:CD1	1:B:122:ASP:OD1	2.43	0.66
1:C:343:GLY:O	1:C:345:PRO:HD3	1.97	0.65
1:B:48:ARG:HG3	1:B:133:LEU:CD2	2.28	0.64
1:B:394:SER:O	1:B:398:VAL:HG23	1.99	0.63
1:C:38:SER:OG	1:C:40:ILE:HB	1.99	0.62
1:A:246:GLU:CD	1:A:368:ARG:HH22	2.04	0.61
1:C:78:ALA:HB3	1:C:93:LEU:HB2	1.81	0.60
1:A:140:GLY:HA2	1:A:145:PRO:HA	1.84	0.60
1:C:288:ASN:HD22	1:C:355[B]:ASN:ND2	1.96	0.59
1:B:78:ALA:HA	1:B:110:ALA:O	2.02	0.59
1:B:11:ILE:O	1:B:37:ARG:NH2	2.36	0.58
1:A:322:THR:O	1:A:323:LYS:HB2	2.03	0.58
1:B:205:GLN:O	1:B:209:LEU:HG	2.04	0.58
1:B:81:THR:HB	1:B:108:TYR:CE2	2.39	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:29:ARG:O	1:A:50:LEU:HA	2.03	0.57
1:B:129[A]:ARG:NH1	1:B:132:GLN:HG2	2.19	0.57
1:C:39:LYS:O	1:C:42[B]:GLU:HG3	2.04	0.57
1:B:134:GLN:OE1	1:C:344:HIS:NE2	2.37	0.57
1:B:228:PHE:CE2	1:B:247:ILE:HG21	2.40	0.57
1:B:48:ARG:HG3	1:B:133:LEU:HD21	1.85	0.57
1:C:79:ILE:HD11	1:C:112:ARG:HD3	1.87	0.57
1:A:83:LEU:HD12	1:A:106:ILE:CG2	2.36	0.56
1:C:246:GLU:CD	1:C:368:ARG:HH22	2.08	0.56
1:B:260:ILE:HA	1:B:360:GLY:O	2.07	0.55
1:C:129:ARG:CB	1:C:132:GLN:HG3	2.36	0.53
1:A:384[B]:ARG:HB3	1:A:385:PRO:CD	2.38	0.53
1:A:326:GLN:HG2	1:A:328:TYR:CZ	2.44	0.52
1:C:73:VAL:HG21	1:C:93:LEU:HB3	1.92	0.51
1:A:57:HIS:HA	1:A:206:ILE:HD11	1.93	0.51
1:B:98:THR:HG21	1:B:230:LEU:HD21	1.93	0.51
1:B:220:SER:OG	1:B:221:VAL:N	2.44	0.51
1:A:13:LYS:HA	1:A:13:LYS:HE2	1.93	0.50
1:A:260:ILE:HA	1:A:360:GLY:O	2.12	0.50
1:C:295:GLU:HG2	1:C:329:ARG:HG2	1.92	0.50
1:A:221:VAL:HG11	1:A:233:ARG:HD2	1.93	0.49
1:A:229:ASN:HB3	1:A:232:SER:HB3	1.94	0.49
1:B:384:ARG:HB3	1:B:385:PRO:HD3	1.94	0.49
1:A:57:HIS:HA	1:A:206:ILE:CD1	2.42	0.49
1:C:81:THR:HB	1:C:108:TYR:CE2	2.46	0.49
1:C:294:PHE:CE2	1:C:330:ALA:HB3	2.47	0.49
1:A:69:PHE:CD1	1:A:122[B]:ASP:OD2	2.66	0.49
1:B:361:ILE:HD12	1:B:361:ILE:N	2.28	0.49
1:C:126:PRO:HB3	1:C:132:GLN:O	2.13	0.48
1:C:139:SER:O	1:C:140:GLY:O	2.30	0.48
1:B:220:SER:O	1:B:221:VAL:HB	2.13	0.48
1:B:74:LEU:CD2	1:B:357:ILE:HG13	2.43	0.47
1:B:71:LEU:HD13	1:B:101:LEU:HD13	1.97	0.47
1:B:100:LYS:HE2	1:B:254:GLN:O	2.15	0.47
1:A:8:ASN:N	2:A:502:HOH:O	2.47	0.47
1:B:294:PHE:CE2	1:B:338:LEU:HD13	2.50	0.47
1:B:229:ASN:HB3	1:B:232:SER:OG	2.14	0.46
1:B:12:PHE:CZ	1:B:37:ARG:HG2	2.50	0.46
1:A:71:LEU:HD22	1:A:122[A]:ASP:OD1	2.16	0.46
1:A:57:HIS:HB3	1:A:201:VAL:O	2.15	0.46
1:B:48:ARG:CG	1:B:133:LEU:HD21	2.46	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:218:LYS:O	1:A:221:VAL:HG12	2.15	0.46
1:A:53:LYS:HA	1:A:118:LEU:O	2.16	0.45
1:B:66:ASP:HB2	1:B:128:ASN:ND2	2.32	0.45
1:A:377:ASN:OD1	1:A:377:ASN:C	2.54	0.45
1:B:98:THR:OG1	1:B:228:PHE:CE2	2.70	0.45
1:C:60:PHE:CD1	1:C:198:ILE:HG12	2.52	0.44
1:C:266:GLU:HG2	1:C:355[A]:ASN:ND2	2.32	0.44
1:B:102:PRO:O	1:B:105:THR:OG1	2.25	0.44
1:A:29:ARG:NH1	1:A:51:GLU:OE1	2.49	0.44
1:A:217:SER:OG	1:A:218:LYS:N	2.49	0.44
1:C:350:ALA:HB1	1:C:352:SER:O	2.18	0.44
1:C:260:ILE:HA	1:C:360:GLY:O	2.18	0.44
1:B:255:LEU:HB3	1:B:260:ILE:O	2.18	0.43
1:B:401:LEU:C	1:B:401:LEU:HD23	2.39	0.43
1:A:52:TYR:CE2	1:A:122[A]:ASP:OD2	2.72	0.43
1:B:230:LEU:HD21	1:B:262:VAL:HG12	2.00	0.43
1:A:81:THR:HB	1:A:108:TYR:CE2	2.53	0.43
1:A:91:PHE:HA	1:A:214:LYS:O	2.19	0.43
1:A:221:VAL:HG11	1:A:233:ARG:CD	2.48	0.43
1:B:245:PHE:O	1:B:263:ASN:HB2	2.18	0.43
1:C:54:SER:HB3	1:C:60:PHE:CZ	2.53	0.43
1:A:11:ILE:O	1:A:37:ARG:NH2	2.36	0.43
1:B:341:PRO:O	1:B:344:HIS:HB2	2.19	0.43
1:B:78:ALA:HB3	1:B:93:LEU:HB2	2.00	0.43
1:A:71:LEU:HD22	1:A:122[A]:ASP:CG	2.39	0.43
1:A:83:LEU:HD12	1:A:106:ILE:HG21	2.00	0.42
1:B:150:GLY:O	1:C:327:ARG:NE	2.50	0.42
1:B:373:GLY:O	1:B:377:ASN:HB2	2.19	0.42
1:C:129:ARG:HB2	1:C:132:GLN:CG	2.42	0.42
1:B:56:PRO:HB3	1:B:113:ASP:O	2.19	0.42
1:C:258:LEU:HB3	1:C:260:ILE:HD12	2.01	0.42
1:C:141:THR:N	1:C:144:GLN:O	2.53	0.42
1:C:24:GLU:OE2	1:C:195:GLU:HG3	2.20	0.42
1:A:100:LYS:HE2	1:A:254:GLN:O	2.19	0.42
1:B:248:THR:OG1	1:B:250[B]:GLU:HG2	2.19	0.42
1:A:71:LEU:HD23	1:A:71:LEU:N	2.34	0.42
1:A:71:LEU:O	1:A:98:THR:HA	2.20	0.42
1:C:221:VAL:HG21	1:C:247:ILE:HD12	2.02	0.42
1:C:53:LYS:HA	1:C:118:LEU:O	2.20	0.41
1:B:48:ARG:HG3	1:B:133:LEU:HD23	2.03	0.41
1:B:384:ARG:N	1:B:385:PRO:HD2	2.35	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:52:TYR:O	1:C:119:ARG:HA	2.21	0.41
1:B:33:ARG:HB2	1:B:36:LYS:HG2	2.03	0.41
1:B:384:ARG:HB3	1:B:385:PRO:CD	2.51	0.41
1:C:123:LEU:HD22	1:C:359:PHE:HB3	2.02	0.41
1:B:66:ASP:CB	1:B:128:ASN:HD22	2.34	0.41
1:B:72:VAL:HG22	1:B:98:THR:HG22	2.03	0.41
1:A:246:GLU:HG3	1:A:263:ASN:HB3	2.02	0.41
1:B:231:ARG:NH2	1:B:266:GLU:OE2	2.54	0.41
1:A:80:LEU:O	1:A:90:SER:HA	2.21	0.41
1:A:205:GLN:O	1:A:209:LEU:HG	2.20	0.41
1:B:65:THR:HG21	1:B:133:LEU:HD12	2.03	0.41
1:C:64:PHE:CE1	1:C:136:PHE:HB2	2.56	0.41
1:C:69:PHE:HB2	1:C:101:LEU:HB3	2.03	0.41
1:B:147:PHE:HA	1:C:325:VAL:HG11	2.03	0.41
1:A:98:THR:OG1	1:A:230:LEU:HD21	2.22	0.40
1:A:15:ASN:OD1	1:A:15:ASN:C	2.60	0.40
1:A:221:VAL:HA	1:A:228:PHE:HB2	2.03	0.40
1:A:237:TYR:HB3	1:A:409:HIS:CG	2.57	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:344:HIS:HE2	1:C:134:GLN:OE1[2_555]	1.58	0.02
1:A:344:HIS:NE2	1:C:134:GLN:OE1[2_555]	2.19	0.01

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	363/431 (84%)	324 (89%)	36 (10%)	3 (1%)	19 35

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	363/431 (84%)	333 (92%)	27 (7%)	3 (1%)	19	35
1	C	360/431 (84%)	332 (92%)	24 (7%)	4 (1%)	14	26
All	All	1086/1293 (84%)	989 (91%)	87 (8%)	10 (1%)	17	31

All (10) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	323	LYS
1	B	15	ASN
1	B	220	SER
1	B	221	VAL
1	C	15	ASN
1	C	140	GLY
1	A	139	SER
1	C	142	GLN
1	A	144	GLN
1	C	130	PRO

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	330/391 (84%)	313 (95%)	17 (5%)	23	44
1	B	328/391 (84%)	308 (94%)	20 (6%)	18	36
1	C	323/391 (83%)	307 (95%)	16 (5%)	24	46
All	All	981/1173 (84%)	928 (95%)	53 (5%)	23	42

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

Mol	Chain	Res	Type
1	A	94	GLU
1	A	121	LEU
1	A	129	ARG
1	A	133	LEU

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Mol	Chain	Res	Type
1	A	139	SER
1	A	141	THR
1	A	204	GLU
1	A	222	SER
1	A	231	ARG
1	A	233	ARG
1	A	237	TYR
1	A	261	PHE
1	A	266	GLU
1	A	322	THR
1	A	396[A]	ARG
1	A	396[B]	ARG
1	A	400	ARG
1	B	52	TYR
1	B	63	GLN
1	B	75	SER
1	B	105	THR
1	B	106	ILE
1	B	127	VAL
1	B	132	GLN
1	B	134	GLN
1	B	135	SER
1	B	139[A]	SER
1	B	139[B]	SER
1	B	198	ILE
1	B	210	SER
1	B	214	LYS
1	B	217	SER
1	B	261	PHE
1	B	272	LEU
1	B	299	GLN
1	B	333	SER
1	B	396	ARG
1	C	24	GLU
1	C	36	LYS
1	C	42[A]	GLU
1	C	42[B]	GLU
1	C	45	GLN
1	C	63	GLN
1	C	132	GLN
1	C	133	LEU
1	C	134	GLN

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Mol	Chain	Res	Type
1	C	137	LEU
1	C	139	SER
1	C	173	GLU
1	C	218	LYS
1	C	224	GLU
1	C	261	PHE
1	C	293	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	412	ASN
1	C	132	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 monosaccharides 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 [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	365/431 (84%)	0.07	6 (1%) 72 74	21, 35, 69, 116	0
1	B	365/431 (84%)	0.04	2 (0%) 91 91	20, 34, 58, 93	0
1	C	363/431 (84%)	0.03	2 (0%) 89 90	23, 35, 60, 99	0
All	All	1093/1293 (84%)	0.05	10 (0%) 84 86	20, 35, 62, 116	0

All (10) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	220	SER	7.5
1	A	144	GLN	2.7
1	A	172	LEU	2.6
1	C	143	ASN	2.3
1	A	214	LYS	2.2
1	A	164	TYR	2.1
1	C	141	THR	2.1
1	A	143	ASN	2.1
1	B	142	GLN	2.0
1	B	39	LYS	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 monosaccharides in this entry.

6.4 Ligands

There are no ligands in this entry.

6.5 Other polymers

There are no such residues in this entry.