



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 12, 2021 – 04:18 PM JST

PDB ID : 6LOH
Title : Crystal structure of the catalytic domain of human ubiquitin ligase AREL1
Authors : Chen, Z.Z.; Li, Z.H.; Shang, G.H.
Deposited on : 2020-01-05
Resolution : 3.21 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 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	:	2.16
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)	:	2.16

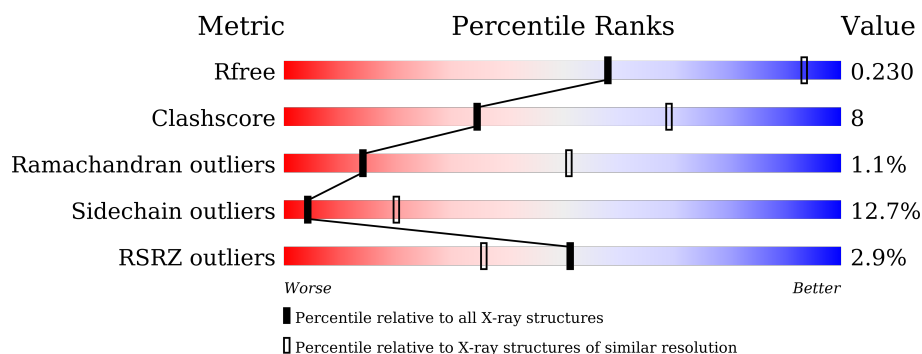
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.21 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	1133 (3.20-3.20)
Clashscore	141614	1253 (3.20-3.20)
Ramachandran outliers	138981	1234 (3.20-3.20)
Sidechain outliers	138945	1233 (3.20-3.20)
RSRZ outliers	127900	1095 (3.20-3.20)

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	389	<div> <div>2%</div> <div> <div></div> <div>76%</div> <div>18%</div> <div>• •</div> </div> </div>
1	B	389	<div> <div>2%</div> <div> <div></div> <div>70%</div> <div>25%</div> <div>• •</div> </div> </div>
1	C	389	<div> <div>4%</div> <div> <div></div> <div>66%</div> <div>25%</div> <div>• •</div> </div> </div>

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 8990 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 Apoptosis-resistant E3 ubiquitin protein ligase 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	373	Total	C	N	O	S	0	0	0
			3003	1921	517	551	14			
1	B	375	Total	C	N	O	S	0	0	0
			3002	1927	514	546	15			
1	C	372	Total	C	N	O	S	0	0	0
			2979	1907	509	548	15			

There are 36 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	423	HIS	-	expression tag	UNP O15033
A	424	HIS	-	expression tag	UNP O15033
A	425	HIS	-	expression tag	UNP O15033
A	426	HIS	-	expression tag	UNP O15033
A	427	HIS	-	expression tag	UNP O15033
A	428	HIS	-	expression tag	UNP O15033
A	429	GLU	-	expression tag	UNP O15033
A	430	ASN	-	expression tag	UNP O15033
A	431	LEU	-	expression tag	UNP O15033
A	432	TYR	-	expression tag	UNP O15033
A	433	PHE	-	expression tag	UNP O15033
A	434	GLN	-	expression tag	UNP O15033
B	423	HIS	-	expression tag	UNP O15033
B	424	HIS	-	expression tag	UNP O15033
B	425	HIS	-	expression tag	UNP O15033
B	426	HIS	-	expression tag	UNP O15033
B	427	HIS	-	expression tag	UNP O15033
B	428	HIS	-	expression tag	UNP O15033
B	429	GLU	-	expression tag	UNP O15033
B	430	ASN	-	expression tag	UNP O15033
B	431	LEU	-	expression tag	UNP O15033
B	432	TYR	-	expression tag	UNP O15033
B	433	PHE	-	expression tag	UNP O15033

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Chain	Residue	Modelled	Actual	Comment	Reference
B	434	GLN	-	expression tag	UNP O15033
C	423	HIS	-	expression tag	UNP O15033
C	424	HIS	-	expression tag	UNP O15033
C	425	HIS	-	expression tag	UNP O15033
C	426	HIS	-	expression tag	UNP O15033
C	427	HIS	-	expression tag	UNP O15033
C	428	HIS	-	expression tag	UNP O15033
C	429	GLU	-	expression tag	UNP O15033
C	430	ASN	-	expression tag	UNP O15033
C	431	LEU	-	expression tag	UNP O15033
C	432	TYR	-	expression tag	UNP O15033
C	433	PHE	-	expression tag	UNP O15033
C	434	GLN	-	expression tag	UNP O15033

- Molecule 2 is water.

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

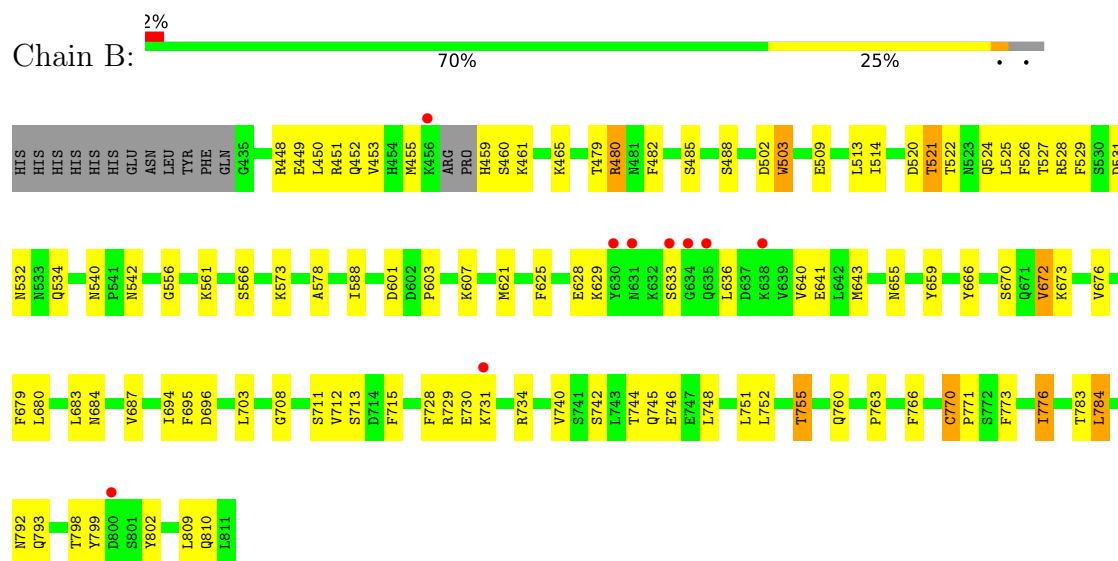
3 Residue-property plots [i](#)

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: Apoptosis-resistant E3 ubiquitin protein ligase 1

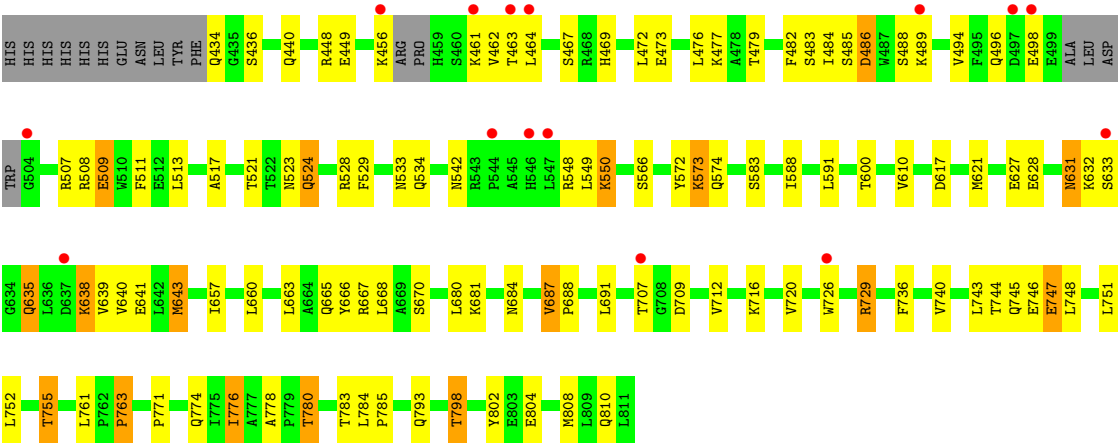


- Molecule 1: Apoptosis-resistant E3 ubiquitin protein ligase 1



- Molecule 1: Apoptosis-resistant E3 ubiquitin protein ligase 1





4 Data and refinement statistics

Property	Value	Source
Space group	P 65 2 2	Depositor
Cell constants a, b, c, α , β , γ	150.23Å 150.23Å 342.64Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	49.22 – 3.21 49.17 – 3.21	Depositor EDS
% Data completeness (in resolution range)	99.8 (49.22-3.21) 99.9 (49.17-3.21)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.81 (at 3.19Å)	Xtriage
Refinement program	REFMAC 5.8.0267	Depositor
R, R_{free}	0.211 , 0.225 0.213 , 0.230	Depositor DCC
R_{free} test set	1878 reflections (4.91%)	wwPDB-VP
Wilson B-factor (Å ²)	71.9	Xtriage
Anisotropy	0.030	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 71.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	8990	wwPDB-VP
Average B, all atoms (Å ²)	74.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.56% 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	0.67	0/3075	0.87	0/4158
1	B	0.68	0/3075	0.85	0/4160
1	C	0.71	0/3048	0.88	0/4119
All	All	0.69	0/9198	0.86	0/12437

There are no bond length outliers.

There are no bond angle outliers.

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	3003	0	2935	36	0
1	B	3002	0	2924	57	0
1	C	2979	0	2900	54	0
2	A	3	0	0	0	0
2	B	3	0	0	0	0
All	All	8990	0	8759	147	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 8.

All (147) 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:770:CYS:CB	1:B:771:PRO:HD3	1.71	1.16
1:B:776:ILE:HD11	1:B:793:GLN:OE1	1.56	1.04
1:B:776:ILE:CD1	1:B:793:GLN:OE1	2.13	0.97
1:C:804:GLU:O	1:C:808:MET:HG2	1.72	0.89
1:B:770:CYS:CB	1:B:771:PRO:CD	2.52	0.84
1:A:676:VAL:O	1:A:680:LEU:CD2	2.26	0.83
1:A:676:VAL:O	1:A:680:LEU:HD23	1.88	0.74
1:A:533:ASN:ND2	1:A:533:ASN:O	2.21	0.74
1:B:520:ASP:O	1:B:522:THR:N	2.26	0.68
1:A:508:ARG:HD3	1:A:574:GLN:HE22	1.63	0.63
1:C:524:GLN:HE21	1:C:524:GLN:HA	1.64	0.62
1:C:784:LEU:HA	1:C:808:MET:CE	2.29	0.62
1:B:751:LEU:O	1:B:755:THR:HG23	2.00	0.61
1:A:684:ASN:HA	1:A:687:VAL:O	2.01	0.61
1:A:534:GLN:HG2	1:A:534:GLN:O	1.99	0.61
1:B:684:ASN:HA	1:B:687:VAL:O	2.01	0.60
1:A:804:GLU:O	1:A:808:MET:HG2	2.02	0.60
1:A:512:GLU:O	1:A:516:LYS:HG3	2.01	0.59
1:C:550:LYS:HE3	1:C:550:LYS:H	1.67	0.59
1:C:534:GLN:HG2	1:C:534:GLN:O	2.02	0.59
1:C:810:GLN:HE21	1:C:810:GLN:HA	1.66	0.58
1:B:776:ILE:HD11	1:B:793:GLN:CD	2.23	0.58
1:C:684:ASN:HA	1:C:687:VAL:O	2.04	0.58
1:B:728:PHE:CD1	1:B:798:THR:HG22	2.39	0.58
1:B:603:PRO:O	1:B:607:LYS:HG3	2.04	0.58
1:C:785:PRO:HD3	1:C:808:MET:HE2	1.86	0.58
1:C:810:GLN:HA	1:C:810:GLN:NE2	2.19	0.58
1:C:533:ASN:OD1	1:C:763:PRO:HG2	2.04	0.58
1:B:695:PHE:CZ	1:B:703:LEU:HD22	2.39	0.57
1:A:521:THR:OG1	1:A:526:PHE:O	2.20	0.56
1:B:459:HIS:O	1:B:461:LYS:NZ	2.38	0.56
1:C:778:ALA:O	1:C:798:THR:HG22	2.05	0.56
1:A:534:GLN:HG3	1:A:762:PRO:CG	2.36	0.56
1:A:676:VAL:O	1:A:680:LEU:HD22	2.06	0.55
1:B:460:SER:C	1:B:461:LYS:HD3	2.26	0.55
1:C:588:ILE:HA	1:C:680:LEU:HD11	1.88	0.55
1:B:672:VAL:HG13	1:B:676:VAL:HG23	1.87	0.55
1:B:532:ASN:C	1:B:534:GLN:H	2.11	0.54
1:C:483:SER:OG	1:C:484:ILE:N	2.40	0.54
1:C:550:LYS:CE	1:C:550:LYS:H	2.20	0.54
1:A:613:ILE:O	1:A:656:LYS:HD2	2.07	0.54
1:C:785:PRO:HD3	1:C:808:MET:CE	2.37	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:631:ASN:C	1:C:633:SER:H	2.11	0.54
1:C:583:SER:OG	1:C:668:LEU:HB3	2.08	0.54
1:C:712:VAL:HG13	1:C:740:VAL:HG12	1.89	0.54
1:B:449:GLU:O	1:B:453:VAL:HG23	2.08	0.53
1:C:631:ASN:HD21	1:C:635:GLN:HG2	1.72	0.53
1:C:524:GLN:HA	1:C:524:GLN:NE2	2.23	0.53
1:B:802:TYR:C	1:B:802:TYR:CD2	2.81	0.53
1:C:524:GLN:HE21	1:C:524:GLN:CA	2.21	0.53
1:A:680:LEU:HD23	1:A:680:LEU:H	1.72	0.53
1:B:485:SER:O	1:B:488:SER:OG	2.25	0.53
1:C:627:GLU:OE2	1:C:667:ARG:NH1	2.42	0.52
1:C:521:THR:HG22	1:C:521:THR:O	2.10	0.52
1:C:529:PHE:N	1:C:529:PHE:CD2	2.77	0.52
1:C:707:THR:HB	1:C:746:GLU:HA	1.91	0.51
1:B:728:PHE:HD1	1:B:798:THR:HG22	1.75	0.51
1:A:618:MET:HA	1:A:621:MET:HG3	1.92	0.51
1:A:549:LEU:HB3	1:A:678:HIS:CD2	2.46	0.51
1:C:566:SER:HA	1:C:573:LYS:HA	1.93	0.51
1:B:521:THR:HG22	1:B:524:GLN:NE2	2.25	0.50
1:B:527:THR:HG22	1:B:540:ASN:HA	1.94	0.50
1:B:566:SER:HA	1:B:573:LYS:HA	1.94	0.50
1:B:776:ILE:CG1	1:B:793:GLN:OE1	2.60	0.50
1:C:726:TRP:HA	1:C:729:ARG:HG3	1.94	0.50
1:A:761:LEU:HD21	1:A:766:PHE:CD2	2.46	0.49
1:C:747:GLU:OE2	1:C:810:GLN:NE2	2.45	0.49
1:C:610:VAL:HG13	1:C:660:LEU:HD22	1.95	0.49
1:B:521:THR:HG22	1:B:524:GLN:HE22	1.78	0.49
1:C:802:TYR:C	1:C:802:TYR:CD2	2.87	0.49
1:B:521:THR:CG2	1:B:524:GLN:HE22	2.26	0.48
1:A:566:SER:HA	1:A:573:LYS:HA	1.95	0.48
1:C:740:VAL:HA	1:C:743:LEU:HD12	1.95	0.48
1:C:688:PRO:HD2	1:C:691:LEU:HD12	1.95	0.48
1:C:709:ASP:N	1:C:709:ASP:OD2	2.47	0.48
1:B:712:VAL:HG13	1:B:740:VAL:HG12	1.96	0.47
1:A:534:GLN:HG3	1:A:762:PRO:HG3	1.96	0.47
1:A:521:THR:HG23	1:A:521:THR:O	2.14	0.47
1:C:736:PHE:O	1:C:740:VAL:HG23	2.14	0.47
1:B:771:PRO:HD2	1:B:792:ASN:ND2	2.29	0.47
1:C:631:ASN:ND2	1:C:635:GLN:HG2	2.30	0.47
1:A:710:ILE:H	1:A:745:GLN:HE22	1.62	0.46
1:C:479:THR:HA	1:C:482:PHE:CD1	2.50	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:479:THR:HA	1:B:482:PHE:CD1	2.50	0.46
1:B:588:ILE:HA	1:B:680:LEU:HD21	1.97	0.46
1:A:556:GLY:HA3	1:A:679:PHE:O	2.16	0.45
1:A:698:ASN:HB3	1:A:753:GLN:CD	2.36	0.45
1:B:556:GLY:HA3	1:B:679:PHE:O	2.16	0.45
1:A:605:PHE:HZ	1:A:663:LEU:HD21	1.81	0.45
1:B:509:GLU:O	1:B:513:LEU:HG	2.15	0.45
1:C:641:GLU:HB3	1:C:643:MET:O	2.17	0.45
1:B:521:THR:HA	1:B:526:PHE:O	2.16	0.45
1:C:709:ASP:HA	1:C:745:GLN:NE2	2.31	0.45
1:A:784:LEU:HA	1:A:784:LEU:HD12	1.88	0.45
1:B:715:PHE:CD1	1:B:766:PHE:CE1	3.05	0.45
1:A:736:PHE:CD1	1:A:809:LEU:HD21	2.52	0.44
1:B:527:THR:CG2	1:B:540:ASN:HA	2.48	0.44
1:B:520:ASP:C	1:B:522:THR:H	2.18	0.44
1:C:666:TYR:HA	1:C:670:SER:HB2	1.98	0.44
1:C:486:ASP:OD1	1:C:486:ASP:N	2.51	0.44
1:B:625:PHE:CZ	1:B:659:TYR:HB2	2.53	0.44
1:C:751:LEU:O	1:C:755:THR:HG23	2.17	0.44
1:B:655:ASN:C	1:B:655:ASN:OD1	2.56	0.44
1:A:762:PRO:HD3	1:A:769:LEU:HD21	2.00	0.44
1:C:780:THR:HG21	1:C:783:THR:CG2	2.48	0.44
1:C:449:GLU:HA	1:C:449:GLU:OE1	2.17	0.44
1:B:641:GLU:HB3	1:B:643:MET:O	2.18	0.43
1:C:660:LEU:O	1:C:663:LEU:HB3	2.18	0.43
1:A:774:GLN:CG	1:A:793:GLN:HG3	2.48	0.43
1:C:638:LYS:HD2	1:C:640:VAL:HG22	1.99	0.43
1:A:534:GLN:O	1:A:579:ARG:NH1	2.50	0.43
1:C:509:GLU:O	1:C:513:LEU:HG	2.18	0.43
1:C:517:ALA:O	1:C:523:ASN:ND2	2.52	0.43
1:B:532:ASN:C	1:B:534:GLN:N	2.71	0.43
1:B:695:PHE:CE2	1:B:703:LEU:HD22	2.53	0.43
1:C:784:LEU:HA	1:C:808:MET:HE3	2.00	0.43
1:B:666:TYR:HA	1:B:670:SER:HB2	2.01	0.42
1:C:440:GLN:OE1	1:C:440:GLN:HA	2.18	0.42
1:A:521:THR:OG1	1:A:527:THR:HA	2.20	0.42
1:A:778:ALA:HB3	1:A:797:PRO:HA	2.01	0.42
1:B:514:ILE:HD12	1:B:514:ILE:HA	1.91	0.42
1:C:751:LEU:O	1:C:755:THR:CG2	2.66	0.42
1:B:521:THR:O	1:B:522:THR:C	2.57	0.42
1:B:784:LEU:HD12	1:B:784:LEU:HA	1.83	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:776:ILE:HD11	1:C:793:GLN:CG	2.50	0.42
1:A:774:GLN:HG2	1:A:793:GLN:HG3	2.02	0.42
1:B:450:LEU:HA	1:B:450:LEU:HD12	1.92	0.42
1:A:688:PRO:HD2	1:A:691:LEU:HD12	2.01	0.42
1:B:448:ARG:HG2	1:B:451:ARG:HH21	1.85	0.42
1:B:708:GLY:O	1:B:745:GLN:NE2	2.54	0.41
1:B:773:PHE:C	1:B:773:PHE:CD1	2.94	0.41
1:B:625:PHE:CE2	1:B:659:TYR:HA	2.55	0.41
1:B:728:PHE:HB2	1:B:798:THR:HG22	2.03	0.41
1:C:720:VAL:O	1:C:774:GLN:HA	2.21	0.41
1:A:617:ASP:OD1	1:A:619:SER:OG	2.33	0.41
1:A:762:PRO:CD	1:A:769:LEU:CD2	2.98	0.41
1:B:480:ARG:HH11	1:B:480:ARG:CB	2.34	0.41
1:B:503:TRP:CE3	1:B:503:TRP:HA	2.55	0.41
1:B:525:LEU:HD12	1:B:525:LEU:HA	1.81	0.41
1:B:730:GLU:OE1	1:B:734:ARG:NH2	2.54	0.41
1:C:496:GLN:OE1	1:C:496:GLN:N	2.53	0.41
1:B:783:THR:O	1:B:799:TYR:OH	2.27	0.40
1:C:511:PHE:CD2	1:C:574:GLN:HG3	2.56	0.40
1:B:578:ALA:O	1:B:760:GLN:NE2	2.52	0.40
1:A:530:SER:C	1:A:532:ASN:H	2.24	0.40
1:A:710:ILE:HB	1:A:748:LEU:HD12	2.03	0.40
1:B:529:PHE:HD2	1:B:601:ASP:O	2.04	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	369/389 (95%)	336 (91%)	29 (8%)	4 (1%)	14 51
1	B	371/389 (95%)	338 (91%)	28 (8%)	5 (1%)	12 47

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	C	366/389 (94%)	337 (92%)	26 (7%)	3 (1%)	19	58
All	All	1106/1167 (95%)	1011 (91%)	83 (8%)	12 (1%)	14	51

All (12) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	498	GLU
1	B	521	THR
1	B	633	SER
1	B	636	LEU
1	A	636	LEU
1	C	632	LYS
1	A	763	PRO
1	C	436	SER
1	C	771	PRO
1	B	770	CYS
1	B	763	PRO
1	A	770	CYS

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	326/346 (94%)	292 (90%)	34 (10%)	7	28
1	B	321/346 (93%)	288 (90%)	33 (10%)	7	29
1	C	321/346 (93%)	265 (83%)	56 (17%)	2	10
All	All	968/1038 (93%)	845 (87%)	123 (13%)	4	20

All (123) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	448	ARG
1	A	469	HIS
1	A	480	ARG

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Mol	Chain	Res	Type
1	A	508	ARG
1	A	509	GLU
1	A	521	THR
1	A	524	GLN
1	A	533	ASN
1	A	534	GLN
1	A	543	ARG
1	A	548	ARG
1	A	600	THR
1	A	604	GLU
1	A	627	GLU
1	A	628	GLU
1	A	635	GLN
1	A	637	ASP
1	A	667	ARG
1	A	674	GLU
1	A	680	LEU
1	A	702	LEU
1	A	709	ASP
1	A	711	SER
1	A	729	ARG
1	A	746	GLU
1	A	748	LEU
1	A	750	ARG
1	A	755	THR
1	A	761	LEU
1	A	763	PRO
1	A	772	SER
1	A	784	LEU
1	A	800	ASP
1	A	811	LEU
1	B	452	GLN
1	B	455	MET
1	B	465	LYS
1	B	480	ARG
1	B	502	ASP
1	B	503	TRP
1	B	528	ARG
1	B	531	ASP
1	B	542	ASN
1	B	561	LYS
1	B	621	MET

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Mol	Chain	Res	Type
1	B	628	GLU
1	B	629	LYS
1	B	640	VAL
1	B	672	VAL
1	B	673	LYS
1	B	683	LEU
1	B	694	ILE
1	B	696	ASP
1	B	711	SER
1	B	713	SER
1	B	729	ARG
1	B	731	LYS
1	B	742	SER
1	B	744	THR
1	B	746	GLU
1	B	748	LEU
1	B	752	LEU
1	B	755	THR
1	B	776	ILE
1	B	784	LEU
1	B	809	LEU
1	B	810	GLN
1	C	434	GLN
1	C	448	ARG
1	C	456	LYS
1	C	461	LYS
1	C	462	VAL
1	C	463	THR
1	C	464	LEU
1	C	467	SER
1	C	469	HIS
1	C	472	LEU
1	C	473	GLU
1	C	476	LEU
1	C	477	LYS
1	C	485	SER
1	C	486	ASP
1	C	488	SER
1	C	489	LYS
1	C	494	VAL
1	C	498	GLU
1	C	507	ARG

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Mol	Chain	Res	Type
1	C	508	ARG
1	C	509	GLU
1	C	524	GLN
1	C	528	ARG
1	C	542	ASN
1	C	548	ARG
1	C	549	LEU
1	C	550	LYS
1	C	572	TYR
1	C	573	LYS
1	C	591	LEU
1	C	600	THR
1	C	617	ASP
1	C	621	MET
1	C	628	GLU
1	C	631	ASN
1	C	635	GLN
1	C	638	LYS
1	C	639	VAL
1	C	643	MET
1	C	657	ILE
1	C	665	GLN
1	C	681	LYS
1	C	687	VAL
1	C	716	LYS
1	C	729	ARG
1	C	744	THR
1	C	747	GLU
1	C	748	LEU
1	C	752	LEU
1	C	755	THR
1	C	761	LEU
1	C	763	PRO
1	C	776	ILE
1	C	780	THR
1	C	798	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	481	ASN
1	A	490	ASN

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Mol	Chain	Res	Type
1	A	574	GLN
1	A	678	HIS
1	A	745	GLN
1	A	793	GLN
1	B	524	GLN
1	B	806	HIS
1	B	810	GLN
1	C	452	GLN
1	C	469	HIS
1	C	524	GLN
1	C	635	GLN
1	C	792	ASN
1	C	806	HIS
1	C	810	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

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	373/389 (95%)	0.04	8 (2%) 63 49	41, 64, 112, 141	0
1	B	375/389 (96%)	0.10	9 (2%) 59 44	44, 69, 113, 146	0
1	C	372/389 (95%)	0.21	15 (4%) 38 25	40, 74, 128, 174	0
All	All	1120/1167 (95%)	0.11	32 (2%) 51 36	40, 69, 121, 174	0

All (32) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	504	GLY	4.9
1	C	497	ASP	4.4
1	B	731	LYS	3.9
1	C	633	SER	3.4
1	B	456	LYS	3.3
1	A	457	ARG	3.2
1	A	458	PRO	3.2
1	B	635	GLN	3.2
1	B	634	GLY	3.1
1	A	435	GLY	3.1
1	C	463	THR	3.0
1	A	709	ASP	2.9
1	C	489	LYS	2.9
1	B	630	TYR	2.8
1	C	547	LEU	2.7
1	A	456	LYS	2.7
1	C	637	ASP	2.7
1	C	707	THR	2.6
1	B	638	LYS	2.6
1	A	534	GLN	2.5
1	C	461	LYS	2.5
1	C	456	LYS	2.4
1	C	464	LEU	2.4

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Mol	Chain	Res	Type	RSRZ
1	C	546	HIS	2.4
1	B	633	SER	2.4
1	C	726	TRP	2.3
1	B	800	ASP	2.3
1	C	544	PRO	2.3
1	A	765	GLY	2.1
1	C	498	GLU	2.1
1	A	731	LYS	2.1
1	B	631	ASN	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 [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.