



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 20, 2026 – 01:59 PM UTC

PDB ID : 6P5L / pdb_00006p5l
Title : Crystal Structure of Ubl123 with an EZH2 peptide
Authors : Saridakis, V.
Deposited on : 2019-05-30
Resolution : 3.30 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

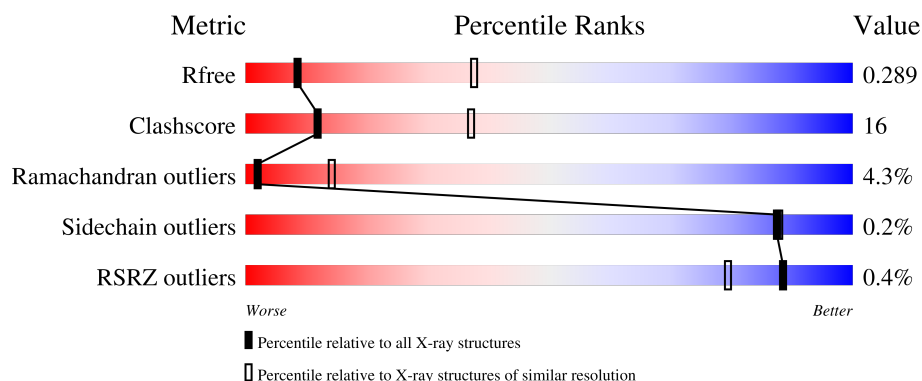
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.30 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	1169 (3.32-3.28)
Clashscore	190562	1209 (3.32-3.28)
Ramachandran outliers	187476	1188 (3.32-3.28)
Sidechain outliers	187428	1187 (3.32-3.28)
RSRZ outliers	180081	1169 (3.32-3.28)

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	356	
1	B	356	
2	D	8	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 5622 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 Ubiquitin carboxyl-terminal hydrolase 7.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	341	Total	C	N	O	S	0	0	0
			2784	1760	474	532	18			
1	B	340	Total	C	N	O	S	0	0	0
			2780	1757	477	530	16			

- Molecule 2 is a protein called PRO-ARG-LYS-LYS-LYS-ARG-LYS-HIS.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	D	6	Total	C	N	O	0	0	0
			58	36	16	6			

PRO	R490	R491	K492	K493	R494	K495	HIS
-----	------	------	------	------	------	------	-----

4 Data and refinement statistics

Property	Value	Source
Space group	P 42 21 2	Depositor
Cell constants a, b, c, α , β , γ	165.45Å 165.45Å 110.90Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.61 – 3.30 29.61 – 3.30	Depositor EDS
% Data completeness (in resolution range)	99.5 (29.61-3.30) 99.3 (29.61-3.30)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.13 (at 3.31Å)	Xtriage
Refinement program	PHENIX 1.13_2998	Depositor
R, R_{free}	0.231 , 0.289 0.240 , 0.289	Depositor DCC
R_{free} test set	2000 reflections (8.40%)	wwPDB-VP
Wilson B-factor (Å ²)	117.6	Xtriage
Anisotropy	0.133	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 69.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	5622	wwPDB-VP
Average B, all atoms (Å ²)	122.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.64% 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	0.41	0/2843	1.24	21/3840 (0.5%)
1	B	0.35	1/2838 (0.0%)	0.86	2/3832 (0.1%)
2	D	0.34	0/57	0.76	0/69
All	All	0.38	1/5738 (0.0%)	1.06	23/7741 (0.3%)

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 (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	597	LEU	CG-CD2	-5.64	1.33	1.52

All (23) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	645	ASN	CB-CG-OD1	28.98	178.76	120.80
1	A	645	ASN	CB-CG-ND2	-26.54	76.59	116.40
1	A	880	GLN	CG-CD-NE2	-22.18	83.14	116.40
1	A	645	ASN	OD1-CG-ND2	-19.98	102.62	122.60
1	A	880	GLN	CA-CB-CG	-10.43	93.23	114.10
1	A	880	GLN	OE1-CD-NE2	-9.90	112.69	122.60
1	B	597	LEU	CB-CG-CD1	7.56	133.38	110.70
1	A	880	GLN	CB-CG-CD	7.35	125.09	112.60
1	A	880	GLN	N-CA-CB	-6.62	99.21	109.48
1	B	544	LEU	CB-CG-CD2	-6.58	90.95	110.70

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	879	GLN	CA-C-N	6.57	130.45	121.05
1	A	879	GLN	C-N-CA	6.57	130.45	121.05
1	A	830	LEU	CA-CB-CG	6.29	138.32	116.30
1	A	880	GLN	CB-CA-C	6.19	119.13	109.84
1	A	774	GLU	CA-CB-CG	-6.03	102.03	114.10
1	A	880	GLN	CG-CD-OE1	5.99	132.79	120.80
1	A	547	GLU	CA-CB-CG	-5.97	102.16	114.10
1	A	871	ARG	N-CA-CB	5.78	120.26	110.49
1	A	871	ARG	CB-CA-C	-5.71	99.06	110.42
1	A	784	LYS	CD-CE-NZ	5.68	130.07	111.90
1	A	847	ASP	N-CA-C	-5.62	106.08	113.17
1	A	880	GLN	N-CA-C	5.31	118.04	110.24
1	A	875	LYS	CD-CE-NZ	-5.22	95.19	111.90

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	880	GLN	Sidechain
1	B	607	GLN	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	2784	0	2727	94	0
1	B	2780	0	2725	91	0
2	D	58	0	77	4	0
All	All	5622	0	5529	176	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 16.

All (176) 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:719:VAL:HG12	1:B:723:ARG:HD2	1.50	0.91
1:B:569:ILE:HG21	1:B:613:MET:HE1	1.56	0.85
1:B:578:HIS:NE2	1:B:580:GLY:O	2.10	0.85
1:A:869:LYS:HB3	1:A:870:PRO:HD2	1.63	0.80
1:A:704:HIS:O	1:A:723:ARG:NH2	2.16	0.79
1:B:647:THR:N	1:B:650:GLU:OE2	2.22	0.72
1:B:541:VAL:HG13	1:B:544:LEU:HD12	1.72	0.71
1:B:672:SER:O	1:B:674:ALA:N	2.24	0.70
1:A:779:GLU:N	1:A:779:GLU:OE1	2.25	0.69
1:B:621:ARG:HD2	1:B:676:LEU:HD21	1.74	0.68
1:B:815:ASN:HB3	1:B:862:ARG:HB2	1.76	0.68
1:A:671:ALA:O	1:A:673:GLY:N	2.25	0.67
1:B:616:PRO:O	1:B:618:ASP:N	2.28	0.67
1:B:542:GLU:O	1:B:545:GLN:N	2.28	0.66
1:B:548:LYS:O	1:B:552:ALA:N	2.29	0.65
1:A:616:PRO:O	1:A:618:ASP:N	2.30	0.65
1:A:642:ALA:C	1:A:644:GLY:H	2.04	0.64
1:A:671:ALA:C	1:A:673:GLY:H	2.05	0.64
1:A:691:MET:HE2	1:A:700:ASN:HB2	1.79	0.63
1:A:761:MET:SD	2:D:495:LYS:HE2	2.38	0.63
1:A:652:SER:OG	1:A:655:GLU:O	2.16	0.63
1:A:646:LYS:HB2	1:A:651:LEU:HD21	1.81	0.62
1:A:691:MET:HG2	1:A:692:TYR:N	2.14	0.62
1:B:713:ILE:HG12	1:B:753:LEU:HD23	1.81	0.61
1:B:818:ASN:ND2	1:B:821:GLN:HG3	2.14	0.61
1:A:835:MET:HE2	1:B:834:PRO:O	2.01	0.61
1:B:555:ARG:HA	1:B:558:ARG:NE	2.16	0.60
1:B:760:LEU:HD12	1:B:761:MET:H	1.66	0.60
1:B:758:ASP:OD1	1:B:758:ASP:N	2.33	0.60
1:A:843:GLN:O	1:A:845:TYR:N	2.34	0.59
1:B:609:LEU:HD11	1:B:613:MET:HE3	1.84	0.59
1:A:679:PHE:HD1	1:A:706:TYR:HD2	1.50	0.59
1:A:602:LEU:HA	1:A:648:MET:HE1	1.84	0.59
1:A:819:TYR:HE2	1:B:835:MET:HE1	1.67	0.58
1:A:759:GLU:HG2	2:D:492:LYS:HD2	1.85	0.58
1:A:852:PRO:HB2	1:B:854:ARG:HH22	1.67	0.58
1:A:639:ASP:OD2	1:A:642:ALA:HB3	2.04	0.58
1:B:652:SER:OG	1:B:655:GLU:N	2.37	0.58
1:A:566:GLN:N	1:A:656:ASN:OD1	2.31	0.57
1:A:639:ASP:HB3	1:A:643:ASP:H	1.70	0.57
1:A:620:ILE:HA	1:A:676:LEU:HD13	1.86	0.57
1:A:797:ILE:CD1	1:A:873:PRO:HB2	2.35	0.56

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:567:VAL:HG11	1:B:605:PHE:CE2	2.40	0.56
1:A:835:MET:HG3	1:B:834:PRO:HB2	1.87	0.55
1:A:852:PRO:HB2	1:B:854:ARG:NH2	2.21	0.55
1:B:727:ILE:O	1:B:730:THR:OG1	2.21	0.55
1:A:779:GLU:H	1:A:779:GLU:CD	2.14	0.55
1:A:800:ASP:N	1:A:806:ASP:OD2	2.32	0.55
1:A:714:ARG:HD2	1:A:750:ASP:OD1	2.07	0.55
1:B:543:ARG:O	1:B:547:GLU:HG3	2.07	0.55
1:A:616:PRO:C	1:A:618:ASP:H	2.15	0.54
1:A:767:VAL:HG21	1:A:787:PHE:HE2	1.72	0.54
1:A:577:GLY:O	1:A:579:GLN:NE2	2.41	0.53
1:B:597:LEU:O	1:B:600:SER:OG	2.26	0.53
1:B:811:VAL:HG22	1:B:829:ARG:NH2	2.25	0.52
1:A:647:THR:O	1:A:650:GLU:HG2	2.09	0.52
1:B:554:LYS:O	1:B:558:ARG:NE	2.41	0.52
1:A:679:PHE:HD1	1:A:706:TYR:CD2	2.28	0.52
1:A:800:ASP:HB2	1:A:878:TYR:CZ	2.44	0.52
1:B:621:ARG:HD2	1:B:676:LEU:CD2	2.38	0.52
1:B:817:MET:HA	1:B:821:GLN:OE1	2.09	0.52
1:B:572:GLU:HB2	1:B:661:PHE:HE1	1.75	0.52
1:A:819:TYR:CE2	1:B:835:MET:HE1	2.44	0.51
1:A:758:ASP:N	1:A:758:ASP:OD1	2.43	0.51
1:B:809:PHE:CZ	1:B:830:LEU:HD11	2.45	0.51
1:A:847:ASP:O	1:A:877:TYR:OH	2.23	0.51
1:B:680:ASP:OD1	1:B:683:HIS:N	2.43	0.51
2:D:494:ARG:HG3	2:D:495:LYS:H	1.76	0.51
1:B:634:ARG:NH2	1:B:738:VAL:O	2.43	0.51
1:B:651:LEU:C	1:B:653:ASP:H	2.19	0.51
1:A:572:GLU:OE2	1:A:704:HIS:ND1	2.44	0.51
1:A:602:LEU:HA	1:A:648:MET:CE	2.41	0.51
1:A:797:ILE:HD12	1:A:873:PRO:HB2	1.93	0.51
1:B:741:ASN:HB2	1:B:791:TYR:CD2	2.46	0.50
1:A:665:VAL:HG23	1:A:676:LEU:HD12	1.93	0.50
1:A:838:GLN:OE1	1:B:856:ASN:HB3	2.11	0.50
1:A:560:GLU:O	1:A:563:LEU:HB2	2.12	0.49
1:B:734:LEU:HD13	1:B:749:TYR:CZ	2.47	0.49
1:A:843:GLN:C	1:A:845:TYR:H	2.19	0.49
1:A:642:ALA:C	1:A:644:GLY:N	2.71	0.49
1:B:564:TYR:CD2	1:B:595:LYS:HE2	2.47	0.49
1:B:578:HIS:CD2	1:B:580:GLY:H	2.30	0.49
1:B:645:ASN:O	1:B:646:LYS:HG2	2.12	0.49

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:547:GLU:HG2	1:A:550:ILE:HD11	1.94	0.49
1:A:759:GLU:HA	2:D:492:LYS:HE3	1.94	0.49
1:B:704:HIS:O	1:B:723:ARG:NH2	2.46	0.49
1:B:735:TYR:HD2	1:B:745:ARG:HA	1.78	0.48
1:B:837:LEU:HD13	1:B:839:PHE:CZ	2.48	0.48
1:A:602:LEU:O	1:A:606:VAL:HG23	2.14	0.48
1:B:671:ALA:C	1:B:673:GLY:H	2.21	0.48
1:B:797:ILE:N	1:B:874:LYS:O	2.34	0.48
1:B:600:SER:HB2	1:B:604:GLU:OE1	2.14	0.47
1:B:824:LYS:CD	1:B:834:PRO:HG3	2.44	0.47
1:A:597:LEU:O	1:A:600:SER:HB3	2.15	0.47
1:A:854:ARG:NH1	1:B:854:ARG:HD2	2.29	0.47
1:A:616:PRO:C	1:A:618:ASP:N	2.72	0.47
1:A:709:ILE:HG22	1:A:754:ASP:HB2	1.97	0.47
1:A:830:LEU:O	1:A:831:ASN:C	2.56	0.47
1:B:843:GLN:HG3	1:B:848:GLY:O	2.14	0.47
1:A:639:ASP:CB	1:A:643:ASP:H	2.27	0.46
1:B:688:PHE:CE2	1:B:704:HIS:HB3	2.50	0.46
1:A:824:LYS:O	1:A:828:GLN:HG3	2.16	0.46
1:A:800:ASP:OD1	1:A:802:THR:OG1	2.22	0.46
1:B:811:VAL:HG22	1:B:829:ARG:CZ	2.45	0.46
1:A:709:ILE:CG2	1:A:754:ASP:HB2	2.45	0.46
1:B:814:SER:C	1:B:816:ARG:H	2.23	0.46
1:A:568:GLN:O	1:A:659:THR:HA	2.15	0.46
1:A:868:PHE:O	1:A:869:LYS:HD3	2.16	0.45
1:A:568:GLN:HB3	1:A:591:TYR:CD2	2.51	0.45
1:B:554:LYS:O	1:B:558:ARG:NH2	2.50	0.45
1:B:645:ASN:C	1:B:646:LYS:HG2	2.42	0.45
1:A:592:THR:HG21	1:A:613:MET:HE3	1.99	0.45
1:B:568:GLN:HB3	1:B:591:TYR:CD2	2.52	0.45
1:B:739:LYS:HG2	1:B:742:LEU:HB3	1.98	0.45
1:A:620:ILE:CA	1:A:676:LEU:HD13	2.46	0.45
1:B:634:ARG:CZ	1:B:765:ILE:HD12	2.47	0.45
1:A:833:ASP:OD2	1:B:824:LYS:NZ	2.37	0.45
1:B:651:LEU:O	1:B:653:ASP:N	2.50	0.45
1:B:733:ILE:HA	1:B:749:TYR:HH	1.82	0.45
1:A:621:ARG:HD2	1:A:706:TYR:OH	2.16	0.45
1:B:727:ILE:HG23	1:B:728:GLN:O	2.17	0.45
1:B:794:VAL:N	1:B:813:LEU:O	2.47	0.44
1:B:820:PHE:O	1:B:824:LYS:HG2	2.17	0.44
1:A:780:LEU:HD13	1:A:786:TYR:HA	2.00	0.44

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:572:GLU:HB2	1:B:661:PHE:CE1	2.53	0.44
1:B:581:ASN:O	1:B:582:ASP:HB2	2.17	0.44
1:B:751:VAL:CG2	1:B:756:ALA:HB2	2.48	0.44
1:A:870:PRO:O	1:A:871:ARG:C	2.60	0.44
1:A:605:PHE:O	1:A:609:LEU:N	2.49	0.44
1:B:671:ALA:O	1:B:673:GLY:N	2.49	0.44
1:B:837:LEU:HD23	1:B:837:LEU:HA	1.73	0.43
1:A:549:ARG:HA	1:A:552:ALA:HB3	2.00	0.43
1:A:597:LEU:O	1:A:598:LYS:C	2.61	0.43
1:A:749:TYR:O	1:A:751:VAL:N	2.49	0.43
1:A:669:LEU:HB3	1:A:674:ALA:CB	2.49	0.43
1:A:683:HIS:O	1:A:709:ILE:HG13	2.19	0.43
1:B:609:LEU:HD11	1:B:613:MET:CE	2.47	0.43
1:B:616:PRO:HB2	1:B:619:GLN:HG2	2.01	0.43
1:B:824:LYS:HD2	1:B:834:PRO:HG3	2.01	0.43
1:B:581:ASN:HA	1:B:737:GLU:OE2	2.19	0.42
1:A:550:ILE:O	1:A:554:LYS:HB2	2.19	0.42
1:B:556:LYS:C	1:B:558:ARG:H	2.28	0.42
1:A:656:ASN:HA	1:A:657:PRO:C	2.45	0.42
1:B:600:SER:O	1:B:648:MET:N	2.49	0.42
1:A:854:ARG:NH1	1:B:852:PRO:HB2	2.34	0.42
1:B:685:VAL:HA	1:B:762:ASP:OD1	2.19	0.42
1:A:668:GLU:C	1:A:670:ALA:N	2.78	0.42
1:B:575:PHE:N	1:B:575:PHE:CD1	2.87	0.42
1:B:687:LEU:HD23	1:B:687:LEU:HA	1.85	0.42
1:A:663:GLU:OE1	1:A:706:TYR:OH	2.34	0.42
1:B:554:LYS:HA	1:B:554:LYS:HD3	1.78	0.42
1:B:759:GLU:O	1:B:761:MET:HG3	2.20	0.42
1:A:613:MET:HB3	1:A:613:MET:HE2	1.84	0.42
1:B:542:GLU:O	1:B:544:LEU:N	2.52	0.42
1:B:609:LEU:HD11	1:B:613:MET:SD	2.59	0.42
1:B:735:TYR:CD2	1:B:745:ARG:HA	2.55	0.42
1:B:739:LYS:HG2	1:B:742:LEU:CB	2.50	0.42
1:A:751:VAL:HG23	1:A:752:SER:O	2.20	0.42
1:A:774:GLU:O	1:A:774:GLU:HG2	2.19	0.41
1:A:803:ILE:HB	1:A:806:ASP:HB2	2.00	0.41
1:A:600:SER:OG	1:A:601:SER:N	2.52	0.41
1:A:814:SER:HB3	1:A:817:MET:HG3	2.02	0.41
1:A:679:PHE:CD1	1:A:706:TYR:CD2	3.07	0.41
1:B:864:LEU:HA	1:B:864:LEU:HD23	1.84	0.41
1:B:580:GLY:O	1:B:581:ASN:O	2.38	0.41

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:550:ILE:HD13	1:A:550:ILE:HG21	1.81	0.41
1:A:562:HIS:CE1	1:A:563:LEU:HG	2.55	0.41
1:A:606:VAL:HG12	1:A:617:GLN:HG3	2.03	0.41
1:A:795:ASP:OD1	1:A:812:THR:OG1	2.29	0.41
1:A:838:GLN:HB2	1:A:881:LEU:HD11	2.03	0.41
1:A:679:PHE:CD1	1:A:706:TYR:HD2	2.35	0.40
1:A:578:HIS:CE1	1:A:580:GLY:H	2.40	0.40
1:A:638:LEU:HD12	1:A:638:LEU:HA	1.86	0.40
1:A:739:LYS:HG2	1:A:742:LEU:HB3	2.03	0.40
1:B:606:VAL:HG12	1:B:617:GLN:HG3	2.01	0.40
1:B:564:TYR:HD2	1:B:595:LYS:HG2	1.85	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	339/356 (95%)	293 (86%)	35 (10%)	11 (3%)	3	19
1	B	336/356 (94%)	272 (81%)	46 (14%)	18 (5%)	1	10
2	D	4/8 (50%)	2 (50%)	2 (50%)	0	100	100
All	All	679/720 (94%)	567 (84%)	83 (12%)	29 (4%)	2	14

All (29) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	617	GLN
1	A	645	ASN
1	A	672	SER
1	A	750	ASP
1	A	870	PRO
1	B	541	VAL

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	B	543	ARG
1	B	557	GLU
1	B	581	ASN
1	B	617	GLN
1	B	652	SER
1	B	673	GLY
1	B	674	ALA
1	B	599	ASN
1	B	672	SER
1	B	815	ASN
1	A	643	ASP
1	A	815	ASN
1	A	835	MET
1	A	869	LYS
1	B	542	GLU
1	B	654	ASN
1	B	578	HIS
1	A	829	ARG
1	B	668	GLU
1	A	844	GLY
1	B	647	THR
1	B	738	VAL
1	B	649	ILE

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	310/327 (95%)	310 (100%)	0	100	100
1	B	309/327 (94%)	308 (100%)	1 (0%)	86	86
2	D	6/8 (75%)	6 (100%)	0	100	100
All	All	625/662 (94%)	624 (100%)	1 (0%)	87	88

All (1) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	597	LEU

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	562	HIS
1	A	581	ASN
1	A	645	ASN
1	A	843	GLN
1	B	566	GLN
1	B	617	GLN
1	B	728	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 [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	341/356 (95%)	-0.29	2 (0%) 85 73	70, 106, 166, 199	0
1	B	340/356 (95%)	-0.21	1 (0%) 90 82	83, 129, 180, 211	0
2	D	6/8 (75%)	0.28	0 100 100	124, 127, 147, 156	0
All	All	687/720 (95%)	-0.24	3 (0%) 88 79	70, 116, 176, 211	0

All (3) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	847	ASP	3.9
1	B	540	LEU	3.3
1	A	816	ARG	2.8

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.