



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 03:22 PM UTC

PDB ID : 2XHS / pdb_00002xhs
Title : Crystal structure of the ligand binding domain of Fushi tarazu factor 1 of *Drosophila melanogaster*.
Authors : Yoo, J.H.; Cho, H.S.
Deposited on : 2010-06-21
Resolution : 2.80 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
Mogul	:	2022.3.0, CSD as543be (2022)
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

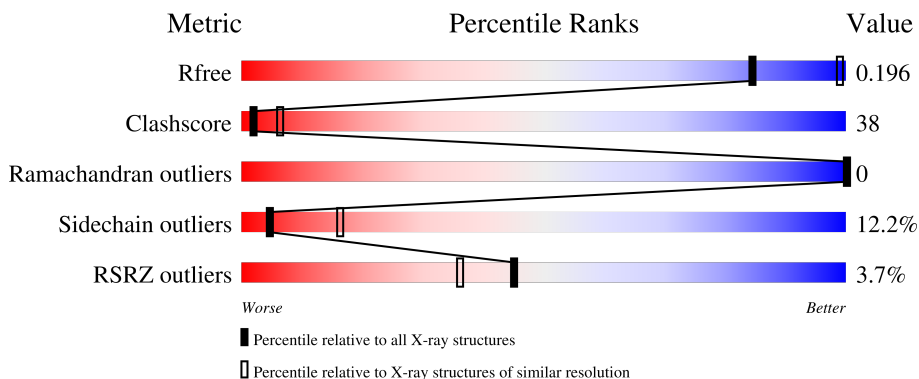
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 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	3866 (2.80-2.80)
Clashscore	190562	4276 (2.80-2.80)
Ramachandran outliers	187476	4196 (2.80-2.80)
Sidechain outliers	187428	4198 (2.80-2.80)
RSRZ outliers	180081	3869 (2.80-2.80)

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	245	2% 49% 40% 10%
2	B	9	33% 33% 33% 33%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 2042 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 NUCLEAR HORMONE RECEPTOR FTZ-F1.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	244	Total	C	N	O	S	Se	6	0	1
			1970	1250	342	363	4	11			

There are 7 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	783	GLY	-	expression tag	UNP P33244
A	784	SER	-	expression tag	UNP P33244
A	785	HIS	-	expression tag	UNP P33244
A	786	MSE	-	expression tag	UNP P33244
A	787	LEU	-	expression tag	UNP P33244
A	788	GLU	-	expression tag	UNP P33244
A	789	ASP	-	expression tag	UNP P33244

- Molecule 2 is a protein called SEGMENTATION PROTEIN FUSHI TARAZU.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	B	9	Total	C	N	O	0	0	0
			69	42	13	14			

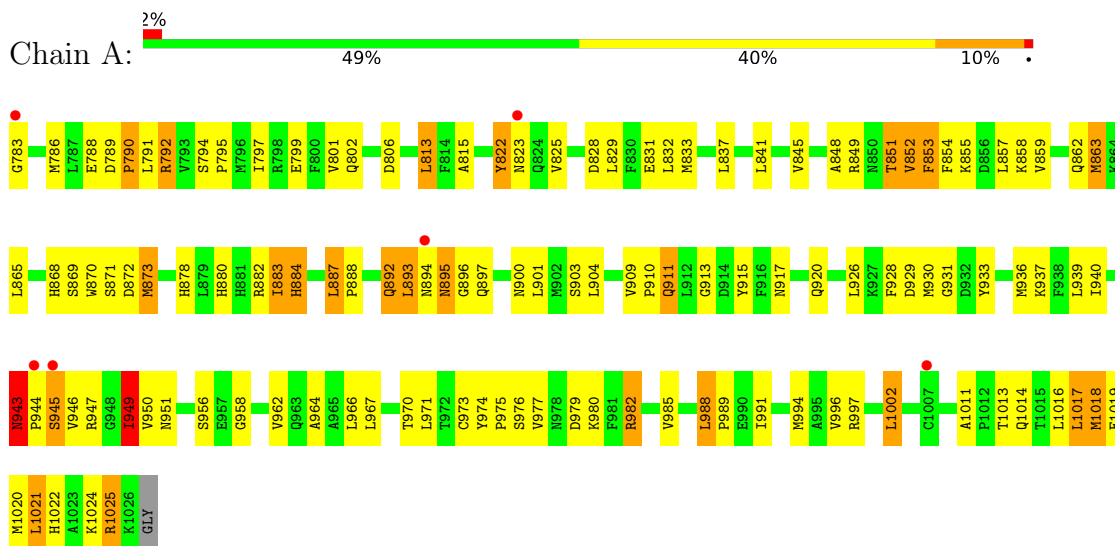
- Molecule 3 is water.

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

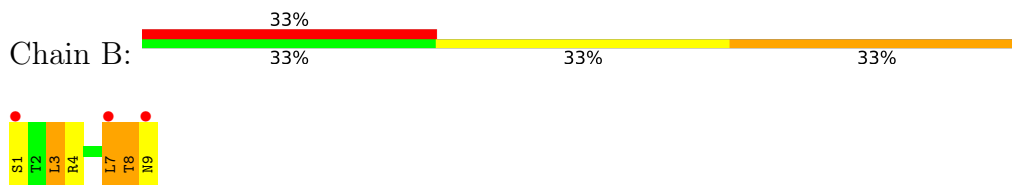
3 Residue-property plots

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: NUCLEAR HORMONE RECEPTOR FTZ-F1



• Molecule 2: SEGMENTATION PROTEIN FUSHI TARAZU



4 Data and refinement statistics

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, α , β , γ	97.93Å 97.93Å 123.47Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	18.60 – 2.80 18.60 – 2.80	Depositor EDS
% Data completeness (in resolution range)	100.0 (18.60-2.80) 99.5 (18.60-2.80)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.15 (at 2.79Å)	Xtriage
Refinement program	REFMAC 5.5.0109	Depositor
R, R_{free}	0.217 , 0.246 (Not available) , 0.196	Depositor DCC
R_{free} test set	875 reflections (5.06%)	wwPDB-VP
Wilson B-factor (Å ²)	49.7	Xtriage
Anisotropy	0.110	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 17.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	0.045 for -h,-k,l	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	2042	wwPDB-VP
Average B, all atoms (Å ²)	54.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.47% 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	2.25	15/2000 (0.8%)	1.22	15/2689 (0.6%)
2	B	1.75	0/68	1.18	0/90
All	All	2.24	15/2068 (0.7%)	1.22	15/2779 (0.5%)

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	2

All (15) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	1011	ALA	CA-CB	-9.63	1.42	1.53
1	A	797	ILE	CA-CB	-7.28	1.45	1.54
1	A	964	ALA	CA-CB	-6.49	1.43	1.53
1	A	991	ILE	CA-CB	-5.74	1.47	1.54
1	A	895	ASN	CA-C	-5.41	1.45	1.52
1	A	859	VAL	CA-CB	-5.40	1.47	1.54
1	A	996	VAL	CA-CB	-5.40	1.47	1.54
1	A	949	ILE	CA-CB	-5.38	1.47	1.54
1	A	988	LEU	C-O	-5.28	1.19	1.24
1	A	878	HIS	C-O	-5.14	1.18	1.24
1	A	871	SER	C-O	-5.12	1.18	1.24
1	A	945	SER	CA-C	-5.08	1.47	1.53
1	A	1018	MSE	SE-CE	-5.05	1.80	1.95
1	A	815	ALA	CA-CB	-5.02	1.45	1.53
1	A	863	MSE	C-O	-5.01	1.18	1.24

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	950	VAL	N-CA-C	9.01	119.00	110.53
1	A	896	GLY	N-CA-C	-7.56	104.79	115.30
1	A	883	ILE	CB-CA-C	-7.05	102.68	111.92
1	A	893	LEU	N-CA-C	6.49	121.04	113.18
1	A	943	ASN	CA-C-N	6.39	126.31	119.28
1	A	943	ASN	C-N-CA	6.39	126.31	119.28
1	A	991	ILE	CB-CA-C	-6.15	103.84	112.14
1	A	884	HIS	N-CA-C	6.13	118.04	111.36
1	A	788	GLU	N-CA-C	6.04	117.73	111.03
1	A	1025	ARG	O-C-N	-5.93	113.51	123.00
1	A	1022	HIS	N-CA-C	5.90	117.79	111.36
1	A	853	PHE	N-CA-C	5.55	117.41	111.36
1	A	997	ARG	N-CA-C	-5.37	105.43	111.28
1	A	790	PRO	N-CA-C	-5.31	106.12	113.81
1	A	868	HIS	N-CA-C	5.05	116.86	111.36

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	792	ARG	Sidechain
1	A	982	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	1970	0	1947	144	0
2	B	69	0	78	20	0
3	A	3	0	0	0	0
All	All	2042	0	2025	156	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 38.

All (156) 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:880:HIS:CD2	1:A:884:HIS:HD2	1.41	1.38
2:B:3:LEU:O	2:B:7:LEU:CD2	1.73	1.34
1:A:1024:LYS:O	1:A:1025:ARG:HG2	1.32	1.28
1:A:791:LEU:CB	1:A:792:ARG:HH11	1.49	1.24
1:A:880:HIS:HD2	1:A:884:HIS:CD2	1.54	1.23
1:A:880:HIS:CD2	1:A:884:HIS:CD2	2.34	1.13
1:A:813:LEU:HD23	1:A:813:LEU:O	1.48	1.13
1:A:791:LEU:HB2	1:A:792:ARG:HH11	1.12	1.11
1:A:813:LEU:HD23	1:A:813:LEU:C	1.76	1.05
1:A:1024:LYS:C	1:A:1025:ARG:HG2	1.71	1.05
1:A:813:LEU:C	1:A:813:LEU:CD2	2.30	1.05
2:B:7:LEU:H	2:B:7:LEU:HD23	1.20	1.04
1:A:791:LEU:HB2	1:A:792:ARG:NH1	1.73	1.04
2:B:3:LEU:O	2:B:7:LEU:HD23	1.59	1.03
2:B:3:LEU:O	2:B:7:LEU:HD21	1.60	0.97
1:A:943:ASN:HD22	1:A:944:PRO:CD	1.77	0.96
1:A:943:ASN:O	1:A:946:VAL:HG23	1.66	0.96
1:A:791:LEU:CB	1:A:792:ARG:NH1	2.28	0.96
1:A:791:LEU:HB3	1:A:792:ARG:HH11	1.33	0.92
1:A:943:ASN:HD22	1:A:943:ASN:C	1.70	0.91
1:A:851:THR:HG22	1:A:854:PHE:H	1.39	0.84
1:A:926:LEU:O	1:A:980:LYS:HE3	1.76	0.84
1:A:943:ASN:C	1:A:943:ASN:ND2	2.30	0.83
1:A:893:LEU:O	1:A:894:ASN:HB2	1.79	0.82
1:A:822:TYR:CE2	1:A:823:ASN:ND2	2.48	0.82
1:A:943:ASN:HD22	1:A:944:PRO:HD2	1.43	0.81
1:A:951:ASN:C	1:A:951:ASN:HD22	1.85	0.81
1:A:943:ASN:ND2	1:A:944:PRO:HD2	1.98	0.78
2:B:8:THR:O	2:B:8:THR:CG2	2.30	0.78
1:A:1017:LEU:CD2	1:A:1017:LEU:C	2.58	0.77
2:B:7:LEU:CD2	2:B:7:LEU:H	1.98	0.77
1:A:943:ASN:HD22	1:A:944:PRO:N	1.82	0.76
1:A:943:ASN:O	1:A:946:VAL:CG2	2.34	0.76
1:A:851:THR:O	1:A:855:LYS:HG3	1.85	0.75
2:B:3:LEU:O	2:B:7:LEU:HD22	1.84	0.75
2:B:8:THR:O	2:B:8:THR:HG23	1.85	0.75
1:A:1016:LEU:HD11	1:A:1020:MSE:HE1	1.70	0.74
1:A:1024:LYS:O	1:A:1025:ARG:CG	2.25	0.74
1:A:1024:LYS:C	1:A:1025:ARG:CG	2.53	0.73
1:A:979:ASP:OD1	1:A:982:ARG:HD3	1.89	0.73
1:A:928:PHE:CE1	1:A:936:MSE:HE1	2.24	0.72
1:A:1017:LEU:CD2	1:A:1017:LEU:O	2.39	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:858:LYS:O	1:A:862:GLN:HG3	1.91	0.70
1:A:791:LEU:HD12	1:A:792:ARG:HH12	1.55	0.70
1:A:929:ASP:C	1:A:929:ASP:OD1	2.30	0.70
2:B:4:ARG:O	2:B:8:THR:HB	1.93	0.69
1:A:849:ARG:NH2	2:B:9:ASN:OXT	2.26	0.69
1:A:882:ARG:HD3	1:A:920:GLN:OE1	1.93	0.68
1:A:799:GLU:O	1:A:802:GLN:HB2	1.94	0.68
1:A:928:PHE:HE1	1:A:936:MSE:CE	2.06	0.68
1:A:1017:LEU:C	1:A:1017:LEU:HD22	2.19	0.68
1:A:936:MSE:O	1:A:940:ILE:HG13	1.93	0.68
1:A:949:ILE:HG22	1:A:949:ILE:O	1.94	0.67
1:A:979:ASP:CG	1:A:982:ARG:HD3	2.20	0.67
1:A:813:LEU:O	1:A:813:LEU:CD2	2.30	0.67
1:A:928:PHE:HE1	1:A:936:MSE:HE1	1.58	0.67
1:A:813:LEU:C	1:A:813:LEU:HD22	2.18	0.67
1:A:1017:LEU:O	1:A:1017:LEU:HD23	1.95	0.66
2:B:3:LEU:C	2:B:7:LEU:HD21	2.20	0.66
1:A:828:ASP:OD2	1:A:831:GLU:HG3	1.96	0.66
1:A:801:VAL:HG12	1:A:801:VAL:O	1.96	0.65
1:A:851:THR:CG2	1:A:854:PHE:H	2.09	0.65
1:A:887:LEU:HD22	1:A:888:PRO:HD2	1.81	0.63
1:A:848:ALA:O	1:A:851:THR:HB	1.99	0.63
1:A:943:ASN:ND2	1:A:944:PRO:CD	2.54	0.62
1:A:795:PRO:O	1:A:799:GLU:HG3	2.01	0.61
1:A:974:TYR:N	1:A:975:PRO:HD3	2.15	0.61
2:B:7:LEU:HD23	2:B:7:LEU:N	2.02	0.61
1:A:988:LEU:N	1:A:989:PRO:HD2	2.15	0.61
1:A:789:ASP:O	1:A:792:ARG:N	2.30	0.60
1:A:863:MSE:HG2	2:B:3:LEU:CD1	2.32	0.60
1:A:863:MSE:HG2	2:B:3:LEU:HD13	1.85	0.59
1:A:783:GLY:HA2	1:A:786:MSE:HB2	1.83	0.59
1:A:791:LEU:CD1	1:A:792:ARG:NH1	2.66	0.59
1:A:895:ASN:HB3	1:A:897:GLN:HB2	1.83	0.59
1:A:951:ASN:C	1:A:951:ASN:ND2	2.53	0.59
1:A:852:VAL:HG13	1:A:853:PHE:CD2	2.38	0.58
1:A:873:MSE:HG2	1:A:937:LYS:HE2	1.86	0.58
1:A:851:THR:HG23	1:A:852:VAL:N	2.18	0.58
1:A:943:ASN:ND2	1:A:945:SER:H	2.02	0.56
1:A:1017:LEU:O	1:A:1017:LEU:HD22	2.03	0.56
1:A:1024:LYS:O	1:A:1024:LYS:HG3	2.05	0.56
1:A:929:ASP:OD1	1:A:931:GLY:N	2.39	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:873:MSE:HG2	1:A:937:LYS:CE	2.36	0.55
1:A:801:VAL:O	1:A:801:VAL:CG1	2.53	0.55
1:A:783:GLY:O	1:A:786:MSE:O	2.25	0.55
1:A:849:ARG:HH21	2:B:9:ASN:C	2.15	0.55
1:A:966:LEU:O	1:A:970:THR:HG23	2.06	0.55
1:A:1017:LEU:C	1:A:1017:LEU:HD23	2.30	0.55
1:A:1019:GLU:OE2	2:B:3:LEU:HB2	2.08	0.54
1:A:789:ASP:OD2	1:A:792:ARG:HD3	2.08	0.53
1:A:1016:LEU:HD11	1:A:1020:MSE:CE	2.38	0.53
1:A:979:ASP:OD1	1:A:982:ARG:CD	2.55	0.53
1:A:951:ASN:ND2	1:A:951:ASN:O	2.42	0.52
1:A:926:LEU:O	1:A:980:LYS:CE	2.54	0.52
2:B:7:LEU:CD2	2:B:7:LEU:N	2.65	0.52
1:A:900:ASN:ND2	1:A:903:SER:OG	2.39	0.52
1:A:913:GLY:O	1:A:917:ASN:ND2	2.43	0.52
1:A:865:LEU:HG	1:A:949:ILE:HD11	1.92	0.51
1:A:1016:LEU:CD1	1:A:1020:MSE:HE1	2.39	0.51
1:A:789:ASP:O	1:A:791:LEU:N	2.44	0.51
1:A:901:LEU:O	1:A:904:LEU:HB2	2.10	0.51
1:A:791:LEU:CD1	1:A:792:ARG:HH12	2.19	0.51
1:A:851:THR:HG23	1:A:853:PHE:H	1.76	0.50
1:A:1016:LEU:CD1	1:A:1020:MSE:CE	2.89	0.50
1:A:915:TYR:CD2	1:A:994:MSE:HG3	2.46	0.49
1:A:944:PRO:HD2	1:A:945:SER:H	1.77	0.49
1:A:849:ARG:NH2	2:B:9:ASN:C	2.69	0.48
1:A:863:MSE:HE3	2:B:3:LEU:HD12	1.96	0.47
1:A:900:ASN:O	1:A:903:SER:HB2	2.14	0.47
1:A:1017:LEU:HD22	1:A:1021:LEU:HD22	1.97	0.47
1:A:988:LEU:N	1:A:989:PRO:CD	2.78	0.47
1:A:893:LEU:O	1:A:894:ASN:CB	2.53	0.46
1:A:939:LEU:HA	1:A:939:LEU:HD23	1.54	0.46
1:A:854:PHE:CE2	1:A:862:GLN:HB3	2.51	0.46
1:A:870:TRP:CH2	1:A:1021:LEU:HD13	2.51	0.46
1:A:789:ASP:O	1:A:790:PRO:C	2.58	0.46
1:A:958:GLY:O	1:A:962:VAL:HG23	2.16	0.46
1:A:1002:LEU:HA	1:A:1002:LEU:HD12	1.70	0.45
1:A:904:LEU:HD23	1:A:904:LEU:HA	1.73	0.45
1:A:893:LEU:C	1:A:895:ASN:H	2.23	0.45
1:A:893:LEU:HB2	1:A:897:GLN:O	2.17	0.45
1:A:973:CYS:C	1:A:975:PRO:HD3	2.41	0.45
1:A:857:LEU:HA	1:A:857:LEU:HD23	1.74	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:806:ASP:OD1	1:A:880:HIS:HE1	1.99	0.45
1:A:979:ASP:OD2	1:A:982:ARG:NH1	2.50	0.45
1:A:789:ASP:C	1:A:791:LEU:N	2.75	0.44
1:A:892:GLN:HE21	1:A:892:GLN:HB3	1.54	0.44
1:A:822:TYR:CD2	1:A:823:ASN:ND2	2.85	0.44
1:A:841:LEU:O	1:A:845:VAL:HG23	2.17	0.44
1:A:985:VAL:O	1:A:985:VAL:HG12	2.16	0.44
1:A:1018:MSE:HE3	1:A:1021:LEU:HB3	2.00	0.44
1:A:1017:LEU:CD2	1:A:1021:LEU:HD22	2.48	0.43
1:A:883:ILE:HG22	1:A:884:HIS:N	2.28	0.43
1:A:971:LEU:HA	1:A:971:LEU:HD23	1.74	0.43
1:A:851:THR:HG22	1:A:851:THR:O	2.12	0.43
1:A:994:MSE:HB3	1:A:994:MSE:HE2	1.77	0.43
1:A:988:LEU:HB2	1:A:989:PRO:CD	2.49	0.43
1:A:863:MSE:HG2	2:B:7:LEU:HD11	2.00	0.42
1:A:870:TRP:CD1	1:A:1020:MSE:HG3	2.54	0.42
1:A:833:MSE:HE3	1:A:837:LEU:HD11	2.00	0.42
1:A:943:ASN:ND2	1:A:945:SER:N	2.67	0.42
1:A:928:PHE:HE1	1:A:936:MSE:HE3	1.83	0.42
1:A:911:GLN:H	1:A:911:GLN:HG2	1.06	0.41
1:A:970:THR:OG1	1:A:971:LEU:N	2.52	0.41
1:A:829:LEU:HD23	1:A:829:LEU:HA	1.88	0.41
1:A:967:LEU:O	1:A:970:THR:OG1	2.38	0.41
1:A:979:ASP:OD2	1:A:982:ARG:HD3	2.20	0.41
1:A:943:ASN:HA	1:A:944:PRO:HD3	1.62	0.41
1:A:928:PHE:CZ	1:A:933:TYR:HB2	2.56	0.41
1:A:801:VAL:HG13	1:A:930:MSE:HE3	2.03	0.41
1:A:792:ARG:N	1:A:792:ARG:HD2	2.36	0.41
1:A:909:VAL:HA	1:A:910:PRO:HD3	1.91	0.41
1:A:1021:LEU:HD12	1:A:1021:LEU:HA	1.82	0.41
1:A:870:TRP:CZ2	1:A:1021:LEU:HD13	2.56	0.40
1:A:854:PHE:O	1:A:857:LEU:HB2	2.20	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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	242/245 (99%)	240 (99%)	2 (1%)	0	100	100
2	B	7/9 (78%)	7 (100%)	0	0	100	100
All	All	249/254 (98%)	247 (99%)	2 (1%)	0	100	100

There are no Ramachandran outliers to report.

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	222/212 (105%)	198 (89%)	24 (11%)	6	21
2	B	8/8 (100%)	4 (50%)	4 (50%)	0	0
All	All	230/220 (104%)	202 (88%)	28 (12%)	5	16

All (28) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	794	SER
1	A	813	LEU
1	A	822	TYR
1	A	825	VAL
1	A	832	LEU
1	A	851	THR
1	A	852	VAL
1	A	869	SER
1	A	872	ASP
1	A	873	MSE
1	A	887	LEU
1	A	892	GLN
1	A	911	GLN
1	A	943	ASN

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Mol	Chain	Res	Type
1	A	947	ARG
1	A	949	ILE
1	A	956	SER
1	A	976	SER
1	A	977	VAL
1	A	1002	LEU
1	A	1013	THR
1	A	1014	GLN
1	A	1017	LEU
1	A	1021	LEU
2	B	1	SER
2	B	3	LEU
2	B	7	LEU
2	B	8	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	810	GLN
1	A	850	ASN
1	A	867	GLN
1	A	878	HIS
1	A	880	HIS
1	A	884	HIS
1	A	892	GLN
1	A	900	ASN
1	A	924	GLN
1	A	943	ASN
1	A	951	ASN

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 [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	233/245 (95%)	-0.17	6 (2%) 57 47	35, 50, 76, 92	5 (2%)
2	B	9/9 (100%)	1.09	3 (33%) 1 1	73, 74, 89, 95	0
All	All	242/254 (95%)	-0.12	9 (3%) 45 36	35, 51, 77, 95	5 (2%)

All (9) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	783	GLY	3.0
1	A	823	ASN	2.8
1	A	945	SER	2.5
2	B	1	SER	2.4
1	A	894	ASN	2.3
1	A	944	PRO	2.3
2	B	7	LEU	2.2
2	B	9	ASN	2.2
1	A	1007	CYS	2.1

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.