



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 1, 2026 – 03:18 AM UTC

PDB ID : 7YC2 / pdb_00007yc2
Title : Crystal structure of auxiliary protein in complex with human protein
Authors : Gao, X.; Cui, S.
Deposited on : 2022-06-30
Resolution : 2.90 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

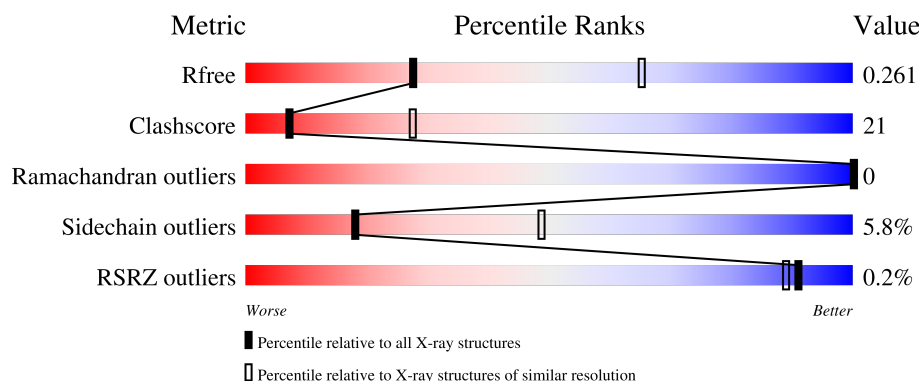
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.90 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	2481 (2.90-2.90)
Clashscore	190562	2690 (2.90-2.90)
Ramachandran outliers	187476	2623 (2.90-2.90)
Sidechain outliers	187428	2625 (2.90-2.90)
RSRZ outliers	180081	2481 (2.90-2.90)

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	240	 55% 42% .
1	B	240	 67% 30% ..
1	C	240	 55% 41% .
1	D	240	 55% 42% .
2	E	7	 14% 14% 43% 43%

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Mol	Chain	Length	Quality of chain
2	F	7	29% 43% 29%
2	G	7	14% 43% 43%
2	H	7	14% 71% 14%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 7885 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 Protein zyg-11 homolog B.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	239	Total	C	N	O	S	0	0	0
			1927	1233	327	357	10			
1	B	234	Total	C	N	O	S	0	0	0
			1884	1205	319	350	10			
1	C	239	Total	C	N	O	S	0	0	0
			1927	1233	327	357	10			
1	D	239	Total	C	N	O	S	0	0	0
			1927	1233	327	357	10			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	489	SER	-	expression tag	UNP Q9C0D3
B	489	SER	-	expression tag	UNP Q9C0D3
C	489	SER	-	expression tag	UNP Q9C0D3
D	489	SER	-	expression tag	UNP Q9C0D3

- Molecule 2 is a protein called ORF.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	E	7	Total	C	N	O	0	0	0
			55	38	8	9			
2	F	7	Total	C	N	O	0	0	0
			55	38	8	9			
2	G	7	Total	C	N	O	0	0	0
			55	38	8	9			
2	H	7	Total	C	N	O	0	0	0
			55	38	8	9			

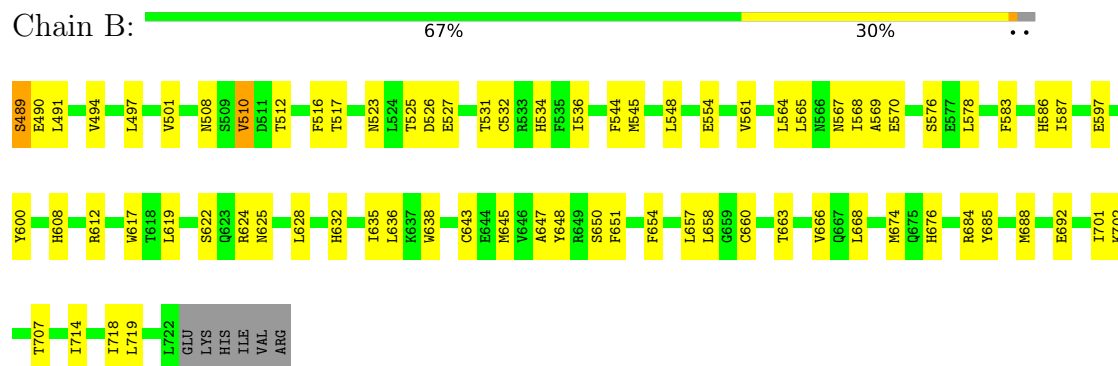
3 Residue-property plots [i](#)

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.

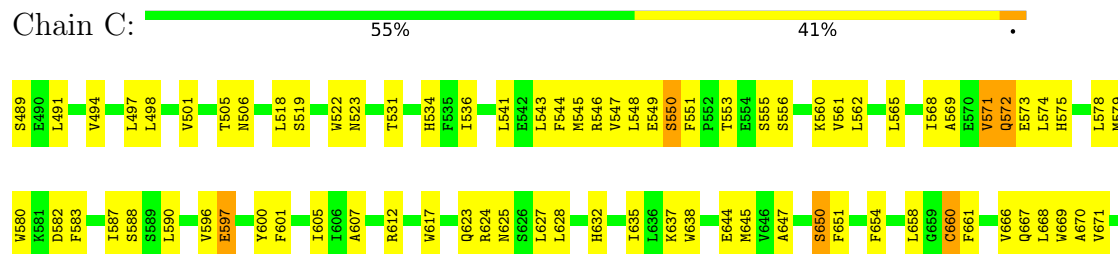
• Molecule 1: Protein zyg-11 homolog B



• Molecule 1: Protein zyg-11 homolog B



• Molecule 1: Protein zyg-11 homolog B





• Molecule 1: Protein zyg-11 homolog B

Chain D: 55% 42% .



• Molecule 2: ORF

Chain E: 14% 14% 43% 43%



• Molecule 2: ORF

Chain F: 29% 43% 29%



• Molecule 2: ORF

Chain G: 14% 43% 43%



• Molecule 2: ORF

Chain H: 14% 71% 14%



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	106.30Å 72.80Å 127.50Å 90.00° 113.40° 90.00°	Depositor
Resolution (Å)	17.05 – 2.90 17.05 – 2.90	Depositor EDS
% Data completeness (in resolution range)	99.6 (17.05-2.90) 99.3 (17.05-2.90)	Depositor EDS
R_{merge}	0.27	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.28 (at 2.90Å)	Xtriage
Refinement program	PHENIX 1.19.2_4158	Depositor
R, R_{free}	0.250 , 0.262 0.244 , 0.261	Depositor DCC
R_{free} test set	985 reflections (4.91%)	wwPDB-VP
Wilson B-factor (Å ²)	72.3	Xtriage
Anisotropy	0.155	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 27.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.003 for h,-k,-h-l	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	7885	wwPDB-VP
Average B, all atoms (Å ²)	68.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 13.47% 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	0.35	0/1974	0.61	4/2684 (0.1%)
1	B	0.28	0/1930	0.47	0/2625
1	C	0.33	0/1974	0.67	3/2684 (0.1%)
1	D	0.33	0/1974	0.64	5/2684 (0.2%)
2	E	0.87	0/56	1.67	1/75 (1.3%)
2	F	0.90	0/56	1.67	2/75 (2.7%)
2	G	0.92	0/56	1.86	2/75 (2.7%)
2	H	0.72	0/56	1.94	2/75 (2.7%)
All	All	0.35	0/8076	0.67	19/10977 (0.2%)

There are no bond length outliers.

All (19) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	661	PHE	N-CA-C	-15.55	94.76	114.04
1	D	619	LEU	N-CA-C	-12.05	93.27	108.45
1	C	660	CYS	N-CA-C	-9.92	95.18	109.96
1	C	660	CYS	CB-CA-C	8.55	122.47	110.16
1	D	619	LEU	CB-CA-C	-6.80	97.45	113.02
1	A	595	GLU	N-CA-C	-6.44	100.78	110.24
1	A	596	VAL	N-CA-C	-6.38	103.35	112.35
1	A	614	GLU	CB-CG-CD	5.79	122.43	112.60
2	H	6	VAL	CA-C-N	-5.62	115.08	123.00
2	H	6	VAL	C-N-CA	-5.62	115.08	123.00
1	D	619	LEU	CA-C-N	5.42	129.05	121.02
1	D	619	LEU	C-N-CA	5.42	129.05	121.02
1	A	595	GLU	CB-CA-C	5.31	117.81	109.84
1	D	490	GLU	N-CA-C	-5.19	105.70	111.36
2	E	3	TYR	CB-CA-C	5.12	120.28	109.79
2	F	7	PHE	CA-C-N	5.08	130.85	121.70
2	F	7	PHE	C-N-CA	5.08	130.85	121.70
2	G	7	PHE	CA-C-N	5.00	130.71	121.70
2	G	7	PHE	C-N-CA	5.00	130.71	121.70

There are no chirality outliers.

There are no planarity outliers.

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	1927	0	1902	79	5
1	B	1884	0	1856	52	1
1	C	1927	0	1902	91	15
1	D	1927	0	1902	89	4
2	E	55	0	51	18	3
2	F	55	0	51	6	1
2	G	55	0	51	14	5
2	H	55	0	51	14	6
All	All	7885	0	7766	324	24

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 21.

All (324) 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:523:ASN:ND2	2:E:3:TYR:CE1	2.06	1.22
1:A:644:GLU:OE2	1:A:681:ASN:HB2	1.00	1.18
1:A:644:GLU:OE2	1:A:681:ASN:CB	1.91	1.17
1:D:644:GLU:HB3	2:H:5:ASN:HD22	0.99	1.13
1:C:644:GLU:O	2:E:5:ASN:ND2	1.82	1.12
1:C:523:ASN:ND2	2:E:3:TYR:CD1	2.19	1.10
1:D:523:ASN:OD1	2:H:4:ILE:HG22	1.55	1.06
1:C:660:CYS:O	1:C:667:GLN:OE1	1.73	1.05
2:E:4:ILE:HD12	2:E:4:ILE:H	1.18	1.04
1:A:523:ASN:OD1	2:F:4:ILE:HG22	1.62	0.99
2:G:4:ILE:HD12	2:G:4:ILE:H	1.27	0.99
1:D:644:GLU:HB3	2:H:5:ASN:ND2	1.79	0.98
1:C:523:ASN:ND2	2:E:3:TYR:HE1	1.58	0.94
1:D:644:GLU:CB	2:H:5:ASN:HD22	1.82	0.92
1:B:647:ALA:HB3	2:G:3:TYR:HB3	1.49	0.92

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:523:ASN:HD21	2:G:4:ILE:H	0.99	0.91
1:A:510:VAL:HG21	1:A:554:GLU:HG3	1.50	0.91
1:C:523:ASN:CG	2:E:3:TYR:HD1	1.82	0.87
1:C:650:SER:HA	1:C:688:MET:HE2	1.55	0.85
1:D:490:GLU:OE1	1:D:490:GLU:N	2.10	0.84
1:B:491:LEU:HG	1:B:534:HIS:CE1	2.12	0.83
1:D:647:ALA:HB3	2:H:3:TYR:HB3	1.60	0.83
1:A:647:ALA:HB3	2:F:3:TYR:HB3	1.61	0.82
1:C:644:GLU:C	2:E:5:ASN:HD21	1.86	0.82
1:C:523:ASN:CG	2:E:3:TYR:CD1	2.58	0.80
1:D:676:HIS:CE1	1:D:680:LYS:HD2	2.19	0.78
1:C:565:LEU:HD13	1:C:583:PHE:HE1	1.49	0.77
1:D:641:PRO:O	1:D:680:LYS:HE3	1.85	0.76
1:D:523:ASN:CG	2:H:4:ILE:HG22	2.11	0.75
1:D:650:SER:HA	1:D:688:MET:SD	2.26	0.75
1:B:516:PHE:HE1	2:G:8:ALA:HB3	1.51	0.74
1:C:696:GLN:O	1:C:700:ASN:ND2	2.23	0.72
1:B:658:LEU:HD11	1:B:674:MET:HE1	1.72	0.72
1:D:628:LEU:HD21	1:D:664:PRO:HD2	1.72	0.71
1:C:660:CYS:O	1:C:667:GLN:CD	2.33	0.71
2:G:4:ILE:HD12	2:G:4:ILE:N	2.04	0.70
1:A:588:SER:HB2	1:A:627:LEU:HD21	1.73	0.70
1:A:712:GLN:O	1:A:716:VAL:HG23	1.93	0.68
1:A:644:GLU:HG2	1:A:680:LYS:HB3	1.74	0.68
1:D:702:LYS:HA	1:D:712:GLN:HE21	1.59	0.68
1:C:556:SER:O	1:C:560:LYS:HD3	1.94	0.67
1:B:523:ASN:ND2	2:G:4:ILE:H	1.84	0.67
1:A:503:GLN:O	1:A:507:GLN:HG3	1.96	0.66
1:A:695:LEU:HD11	1:A:726:ILE:HD11	1.78	0.64
1:A:658:LEU:O	1:A:667:GLN:NE2	2.25	0.64
1:B:569:ALA:O	1:B:612:ARG:NH2	2.30	0.64
1:C:701:ILE:HG23	1:C:707:THR:HG21	1.79	0.64
1:A:658:LEU:HD11	1:A:674:MET:HE1	1.79	0.64
1:B:625:ASN:HA	1:B:628:LEU:HD12	1.79	0.64
1:B:684:ARG:NE	2:F:7:PHE:CZ	2.66	0.63
1:C:654:PHE:HB3	1:C:674:MET:HE2	1.79	0.63
1:C:536:ILE:HD12	1:C:541:LEU:HD11	1.79	0.62
1:A:610:ILE:HG21	1:A:628:LEU:HD21	1.79	0.62
1:A:724:LYS:O	1:A:724:LYS:NZ	2.31	0.62
1:D:491:LEU:HD23	1:D:534:HIS:CG	2.34	0.62
1:C:518:LEU:HB3	1:C:560:LYS:HB3	1.82	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:588:SER:HB2	1:C:627:LEU:HD21	1.82	0.61
1:D:566:ASN:O	1:D:570:GLU:HG3	2.00	0.61
1:D:641:PRO:O	1:D:680:LYS:CE	2.48	0.61
1:D:701:ILE:HG23	1:D:707:THR:HG21	1.83	0.61
1:B:647:ALA:O	2:G:3:TYR:HB2	2.01	0.61
1:D:644:GLU:OE2	1:D:681:ASN:ND2	2.34	0.60
1:C:658:LEU:HD23	1:C:697:HIS:HB3	1.83	0.60
1:C:491:LEU:HD13	1:C:534:HIS:CG	2.37	0.60
1:D:536:ILE:HG13	1:D:574:LEU:HD22	1.83	0.59
1:D:544:PHE:CD2	1:D:561:VAL:HG13	2.38	0.59
1:D:658:LEU:HD12	1:D:697:HIS:HB3	1.85	0.59
1:D:490:GLU:H	1:D:490:GLU:CD	2.05	0.59
1:B:565:LEU:HD13	1:B:583:PHE:HE1	1.68	0.59
1:D:644:GLU:CB	2:H:5:ASN:ND2	2.51	0.59
1:D:644:GLU:OE1	2:H:5:ASN:ND2	2.33	0.59
1:A:600:TYR:OH	1:A:648:TYR:OH	2.10	0.58
1:B:647:ALA:O	2:G:3:TYR:N	2.37	0.58
1:C:579:MET:HE2	1:C:579:MET:HA	1.84	0.58
1:C:697:HIS:O	1:C:701:ILE:HD12	2.03	0.58
1:D:489:SER:HB2	1:D:492:PHE:HB3	1.85	0.58
1:C:650:SER:HA	1:C:688:MET:CE	2.30	0.58
1:A:506:ASN:OD1	1:A:546:ARG:NH2	2.30	0.58
1:C:494:VAL:HG21	1:C:531:THR:HG23	1.85	0.58
1:A:660:CYS:HB3	1:A:663:THR:OG1	2.03	0.58
1:C:562:LEU:HD11	1:C:590:LEU:HD13	1.84	0.58
1:D:601:PHE:HZ	1:D:646:VAL:HB	1.68	0.58
1:D:676:HIS:O	1:D:680:LYS:HG3	2.03	0.57
1:D:491:LEU:HD23	1:D:534:HIS:HB2	1.85	0.57
1:C:545:MET:HE2	1:C:583:PHE:HA	1.87	0.57
2:E:7:PHE:O	2:E:7:PHE:CD1	2.57	0.57
1:C:501:VAL:O	1:C:505:THR:OG1	2.18	0.57
1:D:489:SER:O	1:D:493:ILE:HG13	2.05	0.57
1:C:607:ALA:HA	1:C:666:VAL:HG22	1.87	0.56
1:A:497:LEU:HA	1:A:500:ILE:HD12	1.88	0.56
1:D:536:ILE:HD11	1:D:568:ILE:HG21	1.87	0.56
1:B:545:MET:HE1	1:B:586:HIS:HB2	1.87	0.56
1:A:600:TYR:HA	1:A:638:TRP:CH2	2.41	0.56
1:D:674:MET:HE2	1:D:689:LEU:HD21	1.87	0.56
1:C:617:TRP:CE2	1:C:624:ARG:HB2	2.41	0.55
1:A:687:SER:O	1:A:691:GLU:HG2	2.07	0.55
1:A:654:PHE:HB3	1:A:674:MET:HE2	1.89	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:698:LEU:HB2	1:C:719:LEU:HD21	1.89	0.54
1:D:569:ALA:HB2	1:D:605:ILE:HG23	1.89	0.54
1:B:490:GLU:O	1:B:494:VAL:HG23	2.07	0.54
1:A:494:VAL:HG21	1:A:531:THR:HG23	1.89	0.54
1:A:644:GLU:CG	1:A:680:LYS:HB3	2.38	0.54
1:C:506:ASN:OD1	1:C:546:ARG:NH2	2.38	0.54
1:D:491:LEU:HD23	1:D:534:HIS:CB	2.36	0.54
1:D:724:LYS:HA	1:D:727:VAL:HG22	1.90	0.54
1:C:522:TRP:CZ3	1:C:560:LYS:HA	2.43	0.54
1:C:569:ALA:O	1:C:612:ARG:NH2	2.20	0.54
1:C:596:VAL:HG22	1:C:645:MET:HE1	1.90	0.54
1:A:683:SER:O	2:G:6:VAL:HA	2.07	0.54
1:B:645:MET:HE2	1:B:676:HIS:CE1	2.43	0.54
1:B:660:CYS:HB3	1:B:663:THR:OG1	2.08	0.53
1:C:658:LEU:HG	1:C:701:ILE:HD11	1.90	0.53
1:D:667:GLN:O	1:D:671:VAL:HG22	2.08	0.53
2:G:4:ILE:H	2:G:4:ILE:CD1	1.99	0.53
1:A:565:LEU:HD13	1:A:583:PHE:CE1	2.44	0.53
1:B:565:LEU:HD13	1:B:583:PHE:CE1	2.44	0.53
1:C:523:ASN:HD22	2:E:3:TYR:HE1	1.47	0.53
1:B:648:TYR:CZ	1:B:654:PHE:HZ	2.27	0.53
1:D:660:CYS:O	1:D:663:THR:O	2.26	0.53
1:C:632:HIS:HE1	1:C:710:HIS:HD2	1.56	0.53
1:C:522:TRP:CZ3	2:E:4:ILE:HG23	2.44	0.53
1:A:644:GLU:OE2	1:A:681:ASN:CG	2.52	0.52
1:B:523:ASN:HD21	2:G:4:ILE:N	1.84	0.52
1:A:574:LEU:HD12	1:A:577:GLU:HG3	1.91	0.52
1:C:580:TRP:HB3	1:C:583:PHE:HB3	1.90	0.52
1:D:601:PHE:CZ	1:D:646:VAL:HB	2.44	0.52
1:C:645:MET:HA	2:E:5:ASN:ND2	2.25	0.52
2:E:4:ILE:HD12	2:E:4:ILE:N	2.03	0.52
1:A:660:CYS:HB3	1:A:663:THR:HG1	1.75	0.52
1:B:651:PHE:HZ	1:B:685:TYR:HB3	1.75	0.52
1:D:648:TYR:HB2	1:D:685:TYR:HE1	1.72	0.52
1:C:544:PHE:CD2	1:C:561:VAL:HG13	2.45	0.52
1:D:700:ASN:O	1:D:704:HIS:HB2	2.10	0.52
1:C:651:PHE:CD2	1:C:688:MET:HG2	2.45	0.52
1:D:699:TYR:OH	1:D:723:GLU:OE2	2.22	0.52
1:B:532:CYS:HB3	1:B:568:ILE:HD13	1.91	0.51
1:D:556:SER:O	1:D:560:LYS:NZ	2.35	0.51
1:A:700:ASN:O	1:A:704:HIS:HB2	2.10	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:536:ILE:HG13	1:C:574:LEU:HD22	1.92	0.51
1:A:688:MET:O	1:A:692:GLU:HG3	2.10	0.51
1:D:579:MET:HA	1:D:579:MET:HE3	1.92	0.51
1:D:569:ALA:O	1:D:612:ARG:NH2	2.33	0.51
1:A:490:GLU:HA	1:A:493:ILE:HD12	1.92	0.51
1:A:614:GLU:H	1:A:614:GLU:CD	2.19	0.51
1:C:505:THR:HG23	1:C:551:PHE:CE2	2.46	0.51
1:C:518:LEU:HD13	1:C:561:VAL:HG23	1.93	0.51
1:A:699:TYR:CE1	1:A:719:LEU:HD13	2.46	0.51
1:C:543:LEU:O	1:C:547:VAL:HG23	2.11	0.51
1:D:597:GLU:HA	1:D:645:MET:SD	2.51	0.51
1:B:544:PHE:CD2	1:B:561:VAL:HG13	2.46	0.50
1:C:580:TRP:CE2	1:C:582:ASP:HB2	2.47	0.50
1:C:716:VAL:HA	1:C:719:LEU:HD12	1.94	0.50
1:D:643:CYS:O	1:D:680:LYS:CD	2.59	0.50
1:D:645:MET:HE2	1:D:676:HIS:NE2	2.27	0.50
2:E:3:TYR:HE1	2:E:8:ALA:HA	1.75	0.50
1:D:494:VAL:HG21	1:D:531:THR:HG23	1.93	0.50
1:D:523:ASN:ND2	2:H:4:ILE:CG2	2.75	0.49
1:B:583:PHE:O	1:B:587:ILE:HG12	2.13	0.49
1:D:606:ILE:O	1:D:610:ILE:HG13	2.12	0.49
1:A:614:GLU:N	1:A:614:GLU:OE1	2.45	0.49
1:C:667:GLN:HE21	1:C:711:VAL:HG11	1.77	0.49
1:C:688:MET:HE3	1:C:692:GLU:OE2	2.12	0.49
1:D:676:HIS:HE1	1:D:680:LYS:HD2	1.69	0.49
1:B:635:ILE:HG21	1:B:668:LEU:HG	1.93	0.49
1:A:641:PRO:HG2	1:A:676:HIS:CE1	2.48	0.49
1:D:565:LEU:HA	1:D:568:ILE:HG12	1.94	0.49
1:C:522:TRP:CE3	1:C:560:LYS:HA	2.48	0.48
1:A:504:LYS:HB2	1:A:514:LEU:HD13	1.95	0.48
1:A:652:ASN:HA	1:A:655:PHE:CD2	2.48	0.48
1:B:490:GLU:OE1	1:B:490:GLU:N	2.37	0.48
1:D:704:HIS:ND1	1:D:705:GLU:O	2.47	0.48
1:B:497:LEU:HD22	1:B:517:THR:HG23	1.95	0.48
1:D:617:TRP:CD1	1:D:619:LEU:O	2.66	0.48
2:E:4:ILE:H	2:E:4:ILE:CD1	1.96	0.48
1:A:520:ALA:O	1:A:524:LEU:HG	2.14	0.48
1:C:667:GLN:O	1:C:671:VAL:HG22	2.14	0.48
1:D:600:TYR:HA	1:D:638:TRP:CH2	2.48	0.48
1:D:542:GLU:OE1	1:D:542:GLU:N	2.41	0.47
1:A:628:LEU:HD23	1:A:628:LEU:HA	1.47	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:617:TRP:CG	1:B:624:ARG:HD2	2.49	0.47
1:C:670:ALA:O	1:C:674:MET:HG3	2.13	0.47
1:D:545:MET:HA	1:D:548:LEU:HD12	1.96	0.47
1:A:671:VAL:O	1:A:718:ILE:HD11	2.14	0.47
1:A:683:SER:HB2	2:G:5:ASN:ND2	2.29	0.47
1:A:617:TRP:CG	1:A:624:ARG:HD2	2.50	0.47
1:A:580:TRP:CE3	1:A:583:PHE:HB2	2.50	0.47
1:A:703:ASP:OD1	1:A:703:ASP:N	2.47	0.47
1:C:568:ILE:O	1:C:571:VAL:HG13	2.15	0.47
1:D:583:PHE:O	1:D:587:ILE:HG12	2.15	0.47
1:D:644:GLU:CD	1:D:681:ASN:ND2	2.73	0.47
1:D:660:CYS:HB3	1:D:663:THR:OG1	2.14	0.47
1:B:526:ASP:OD1	1:B:527:GLU:HG2	2.15	0.47
1:B:490:GLU:N	1:B:490:GLU:CD	2.73	0.47
1:A:565:LEU:HD13	1:A:583:PHE:HE1	1.79	0.47
1:D:658:LEU:HD11	1:D:674:MET:HE1	1.97	0.47
1:C:575:HIS:ND1	1:C:612:ARG:HG3	2.30	0.46
1:D:523:ASN:CG	2:H:4:ILE:CG2	2.83	0.46
1:A:617:TRP:CB	1:A:624:ARG:HD2	2.46	0.46
1:B:578:LEU:HG	1:B:583:PHE:CE2	2.51	0.46
1:C:597:GLU:H	1:C:597:GLU:HG2	1.46	0.46
1:D:654:PHE:HB3	1:D:674:MET:HE3	1.97	0.46
1:A:515:LYS:HG2	1:A:557:ILE:HD11	1.98	0.46
1:B:657:LEU:HB3	1:B:666:VAL:HG12	1.96	0.46
1:D:523:ASN:ND2	2:H:4:ILE:HG22	2.30	0.46
1:A:606:ILE:HG22	1:A:631:LEU:HD22	1.97	0.46
1:C:708:ASP:HB3	1:C:711:VAL:HB	1.96	0.46
1:D:497:LEU:HD22	1:D:517:THR:HG23	1.98	0.46
1:A:510:VAL:HG23	1:A:551:PHE:CD1	2.51	0.46
1:B:688:MET:O	1:B:692:GLU:HG3	2.16	0.46
1:B:688:MET:SD	2:F:7:PHE:HB3	2.56	0.46
1:C:719:LEU:O	1:C:723:GLU:HG2	2.15	0.46
1:D:581:LYS:NZ	1:D:585:ASP:OD2	2.31	0.45
1:B:516:PHE:CE1	2:G:8:ALA:HB3	2.41	0.45
1:A:551:PHE:HB3	1:A:554:GLU:HB2	1.98	0.45
1:C:660:CYS:O	1:C:667:GLN:HG3	2.17	0.45
1:B:490:GLU:CD	1:B:490:GLU:H	2.21	0.45
1:B:532:CYS:O	1:B:536:ILE:HG12	2.16	0.45
1:B:684:ARG:HG2	1:B:685:TYR:CE2	2.52	0.45
1:D:596:VAL:HB	1:D:639:PRO:HD2	1.98	0.45
1:D:673:ALA:O	1:D:677:VAL:HG22	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:F:4:ILE:HG13	2:F:6:VAL:H	1.81	0.45
1:A:544:PHE:CD2	1:A:561:VAL:HG13	2.51	0.45
1:C:545:MET:O	1:C:549:GLU:N	2.43	0.45
1:C:565:LEU:HD13	1:C:583:PHE:CE1	2.40	0.45
1:C:667:GLN:NE2	1:C:711:VAL:HG11	2.32	0.45
1:C:704:HIS:HB3	1:C:707:THR:OG1	2.16	0.45
1:A:664:PRO:HB3	1:A:708:ASP:HB2	1.99	0.44
1:B:701:ILE:HG23	1:B:707:THR:HG21	1.99	0.44
1:C:587:ILE:H	1:C:587:ILE:HG12	1.68	0.44
1:A:601:PHE:HZ	1:A:646:VAL:HB	1.83	0.44
1:C:546:ARG:O	1:C:550:SER:OG	2.28	0.44
1:C:635:ILE:HB	1:C:668:LEU:HD23	2.00	0.44
1:A:714:ILE:O	1:A:718:ILE:HG13	2.18	0.44
1:D:714:ILE:O	1:D:718:ILE:HG13	2.18	0.44
1:B:654:PHE:CB	1:B:674:MET:HE3	2.48	0.44
1:B:536:ILE:HD11	1:B:568:ILE:HG21	1.98	0.43
1:C:601:PHE:O	1:C:605:ILE:HG13	2.18	0.43
1:D:654:PHE:CB	1:D:674:MET:HE3	2.48	0.43
2:E:4:ILE:N	2:E:4:ILE:CD1	2.73	0.43
1:A:606:ILE:HD13	1:A:627:LEU:HD22	2.00	0.43
1:D:640:THR:HG22	1:D:680:LYS:HE2	1.99	0.43
1:D:498:LEU:HD22	1:D:543:LEU:HD23	1.99	0.43
1:D:619:LEU:HB3	1:D:620:SER:H	1.62	0.43
1:A:518:LEU:HD13	1:A:561:VAL:HG23	2.00	0.43
1:B:714:ILE:O	1:B:718:ILE:HG13	2.18	0.43
1:D:718:ILE:HG22	1:D:722:LEU:HG	2.00	0.43
1:A:498:LEU:HD22	1:A:543:LEU:HD23	2.00	0.43
1:C:497:LEU:O	1:C:501:VAL:HG23	2.18	0.43
1:C:637:LYS:HE2	1:C:637:LYS:HB2	1.72	0.43
1:C:687:SER:HA	1:C:690:ILE:HD12	2.00	0.43
1:A:621:ARG:H	1:A:621:ARG:HD3	1.84	0.43
1:C:632:HIS:CE1	1:C:710:HIS:HD2	2.35	0.43
1:C:660:CYS:O	1:C:667:GLN:CG	2.66	0.43
1:C:695:LEU:HD11	1:C:726:ILE:HD11	2.01	0.43
1:A:545:MET:HE3	1:A:549:GLU:HG3	2.01	0.43
1:D:491:LEU:HB3	1:D:534:HIS:CD2	2.54	0.43
1:D:523:ASN:HD21	2:H:4:ILE:CG2	2.32	0.43
1:D:632:HIS:CE1	1:D:636:LEU:HD11	2.54	0.43
1:B:600:TYR:HA	1:B:638:TRP:CH2	2.54	0.43
1:A:571:VAL:HG11	1:A:574:LEU:HD23	2.01	0.42
1:D:527:GLU:HB2	1:D:649:ARG:HH22	1.84	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:498:LEU:HD23	1:C:498:LEU:HA	1.82	0.42
1:D:522:TRP:CE3	1:D:560:LYS:HA	2.54	0.42
1:A:554:GLU:O	1:A:558:GLN:HG3	2.19	0.42
1:C:519:SER:O	1:C:522:TRP:HB3	2.19	0.42
1:D:675:GLN:O	1:D:679:SER:HB3	2.18	0.42
1:C:600:TYR:HA	1:C:638:TRP:CH2	2.55	0.42
1:C:625:ASN:HA	1:C:628:LEU:HD12	2.01	0.42
1:D:515:LYS:HB3	1:D:560:LYS:HE2	2.01	0.42
1:A:497:LEU:O	1:A:501:VAL:HG23	2.19	0.42
1:B:632:HIS:CE1	1:B:636:LEU:HD11	2.55	0.42
1:B:508:ASN:CG	1:B:508:ASN:O	2.61	0.42
1:A:588:SER:O	1:A:591:LEU:HB3	2.20	0.42
1:C:541:LEU:HD21	1:C:578:LEU:HD23	2.01	0.42
1:C:583:PHE:O	1:C:587:ILE:HG12	2.20	0.42
1:D:644:GLU:OE1	1:D:681:ASN:ND2	2.53	0.42
1:D:647:ALA:CB	2:H:3:TYR:HB3	2.42	0.42
1:C:654:PHE:HE1	1:C:669:TRP:CH2	2.38	0.42
1:C:689:LEU:HD23	1:C:722:LEU:HD11	2.02	0.42
1:A:544:PHE:HD2	1:A:561:VAL:HG13	1.85	0.41
1:A:629:ASP:OD1	1:A:629:ASP:N	2.53	0.41
1:D:605:ILE:O	1:D:609:LEU:HG	2.19	0.41
1:C:578:LEU:HD22	1:C:583:PHE:CZ	2.55	0.41
1:C:689:LEU:HD12	1:C:689:LEU:HA	1.92	0.41
1:B:548:LEU:HG	1:B:561:VAL:HG11	2.02	0.41
1:B:564:LEU:O	1:B:567:ASN:HB2	2.20	0.41
1:C:647:ALA:HA	1:C:685:TYR:OH	2.20	0.41
1:A:584:ILE:O	1:A:588:SER:HB3	2.20	0.41
1:C:541:LEU:O	1:C:545:MET:HG2	2.21	0.41
1:A:680:LYS:HD3	1:A:680:LYS:HA	1.84	0.41
1:B:510:VAL:HG11	1:B:554:GLU:OE2	2.20	0.41
1:B:702:LYS:HD3	1:B:719:LEU:CD1	2.51	0.41
2:G:6:VAL:O	2:G:6:VAL:HG12	2.20	0.41
1:C:587:ILE:HD12	1:C:605:ILE:HD12	2.02	0.41
1:D:524:LEU:HB3	1:D:531:THR:HG21	2.03	0.41
1:D:647:ALA:HA	1:D:685:TYR:OH	2.20	0.41
2:E:4:ILE:HG22	2:E:6:VAL:HG23	2.03	0.41
1:B:497:LEU:O	1:B:501:VAL:HG23	2.20	0.41
1:A:526:ASP:OD2	1:A:649:ARG:HD3	2.20	0.41
1:A:548:LEU:HG	1:A:561:VAL:HG11	2.03	0.41
1:A:723:GLU:OE1	1:A:723:GLU:HA	2.21	0.41
1:B:570:GLU:HG2	1:B:608:HIS:CE1	2.56	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:617:TRP:NE1	1:B:619:LEU:HB2	2.36	0.41
1:C:697:HIS:O	1:C:700:ASN:HB2	2.21	0.41
2:E:4:ILE:CG2	2:E:6:VAL:HG23	2.51	0.41
1:A:523:ASN:CG	2:F:4:ILE:HG22	2.39	0.41
1:A:544:PHE:CE2	1:A:564:LEU:HD13	2.56	0.41
1:A:627:LEU:HA	1:A:627:LEU:HD23	1.79	0.41
1:A:667:GLN:O	1:A:671:VAL:HG22	2.21	0.41
1:D:502:LYS:HE2	1:D:503:GLN:N	2.36	0.41
1:A:601:PHE:O	1:A:605:ILE:HD12	2.21	0.40
1:C:562:LEU:HB3	1:C:601:PHE:HB2	2.03	0.40
1:D:702:LYS:HA	1:D:712:GLN:NE2	2.32	0.40
1:C:498:LEU:HD22	1:C:543:LEU:HD23	2.03	0.40
1:D:648:TYR:HB2	1:D:685:TYR:CE1	2.55	0.40
1:A:536:ILE:HD13	1:A:536:ILE:HA	1.93	0.40
1:A:549:GLU:O	1:A:552:PRO:HD3	2.21	0.40
1:A:667:GLN:HB3	1:A:711:VAL:HG11	2.03	0.40
1:B:525:THR:HG22	1:B:531:THR:HG22	2.02	0.40
1:C:548:LEU:HG	1:C:561:VAL:HG11	2.02	0.40
1:D:523:ASN:OD1	2:H:4:ILE:CG2	2.46	0.40
1:C:632:HIS:HE1	1:C:710:HIS:CD2	2.38	0.40
1:D:651:PHE:O	1:D:654:PHE:N	2.44	0.40
1:A:645:MET:HE3	1:A:645:MET:HB2	1.78	0.40

All (24) 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:489:SER:N	2:G:8:ALA:C[2_555]	0.20	2.00
1:C:572:GLN:CD	1:C:572:GLN:OE1[2_556]	0.56	1.64
1:C:572:GLN:CD	1:C:572:GLN:CD[2_556]	0.68	1.52
1:C:572:GLN:OE1	1:C:572:GLN:NE2[2_556]	1.02	1.18
1:C:489:SER:N	2:H:8:ALA:O[2_546]	1.10	1.10
1:C:489:SER:CA	2:H:8:ALA:O[2_546]	1.18	1.02
1:C:572:GLN:CG	1:C:572:GLN:OE1[2_556]	1.42	0.78
1:A:489:SER:N	2:G:8:ALA:O[2_555]	1.43	0.77
1:A:489:SER:N	2:G:8:ALA:CA[2_555]	1.47	0.73
1:A:489:SER:CA	2:G:8:ALA:C[2_555]	1.56	0.64
1:C:489:SER:CB	2:H:8:ALA:O[2_546]	1.67	0.53
1:C:572:GLN:OE1	1:C:572:GLN:OE1[2_556]	1.78	0.42
1:C:572:GLN:CD	1:C:572:GLN:NE2[2_556]	1.81	0.39

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:490:GLU:N	2:G:8:ALA:O[2_555]	1.84	0.36
1:C:489:SER:N	2:H:8:ALA:CA[2_546]	1.89	0.31
1:C:572:GLN:CG	1:C:572:GLN:CD[2_556]	1.92	0.28
1:C:489:SER:N	2:H:8:ALA:CB[2_546]	1.97	0.23
1:C:489:SER:CA	2:H:8:ALA:C[2_546]	2.06	0.14
1:D:687:SER:C	2:E:7:PHE:CE2[1_565]	2.07	0.13
1:D:688:MET:N	2:E:7:PHE:CE2[1_565]	2.08	0.12
1:B:489:SER:N	2:F:8:ALA:O[2_555]	2.11	0.09
1:C:572:GLN:CB	1:C:572:GLN:OE1[2_556]	2.11	0.09
1:C:684:ARG:NH2	1:D:684:ARG:NH2[1_545]	2.15	0.05
1:D:688:MET:CA	2:E:7:PHE:CZ[1_565]	2.18	0.02

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	237/240 (99%)	232 (98%)	5 (2%)	0	100	100
1	B	232/240 (97%)	226 (97%)	6 (3%)	0	100	100
1	C	237/240 (99%)	233 (98%)	4 (2%)	0	100	100
1	D	237/240 (99%)	229 (97%)	8 (3%)	0	100	100
2	E	5/7 (71%)	5 (100%)	0	0	100	100
2	F	5/7 (71%)	5 (100%)	0	0	100	100
2	G	5/7 (71%)	5 (100%)	0	0	100	100
2	H	5/7 (71%)	5 (100%)	0	0	100	100
All	All	963/988 (98%)	940 (98%)	23 (2%)	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	219/220 (100%)	205 (94%)	14 (6%)	16	44
1	B	214/220 (97%)	206 (96%)	8 (4%)	30	64
1	C	219/220 (100%)	206 (94%)	13 (6%)	18	48
1	D	219/220 (100%)	210 (96%)	9 (4%)	27	61
2	E	5/5 (100%)	3 (60%)	2 (40%)	0	0
2	F	5/5 (100%)	4 (80%)	1 (20%)	1	4
2	G	5/5 (100%)	2 (40%)	3 (60%)	0	0
2	H	5/5 (100%)	3 (60%)	2 (40%)	0	0
All	All	891/900 (99%)	839 (94%)	52 (6%)	18	49

All (52) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	491	LEU
1	A	530	THR
1	A	553	THR
1	A	574	LEU
1	A	588	SER
1	A	596	VAL
1	A	614	GLU
1	A	620	SER
1	A	621	ARG
1	A	623	GLN
1	A	675	GLN
1	A	703	ASP
1	A	712	GLN
1	A	713	GLN
1	B	489	SER
1	B	510	VAL
1	B	512	THR
1	B	576	SER
1	B	597	GLU

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Mol	Chain	Res	Type
1	B	622	SER
1	B	643	CYS
1	B	650	SER
1	C	550	SER
1	C	553	THR
1	C	555	SER
1	C	571	VAL
1	C	572	GLN
1	C	573	GLU
1	C	597	GLU
1	C	623	GLN
1	C	650	SER
1	C	675	GLN
1	C	683	SER
1	C	716	VAL
1	C	725	HIS
1	D	490	GLU
1	D	594	VAL
1	D	615	GLN
1	D	643	CYS
1	D	662	THR
1	D	684	ARG
1	D	706	HIS
1	D	707	THR
1	D	726	ILE
2	E	4	ILE
2	E	6	VAL
2	F	4	ILE
2	G	3	TYR
2	G	4	ILE
2	G	6	VAL
2	H	4	ILE
2	H	7	PHE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (28) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	496	GLN
1	A	652	ASN
1	A	681	ASN
1	A	712	GLN
1	B	499	GLN

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Mol	Chain	Res	Type
1	B	503	GLN
1	B	508	ASN
1	B	523	ASN
1	B	534	HIS
1	B	700	ASN
1	B	706	HIS
1	C	499	GLN
1	C	539	GLN
1	C	572	GLN
1	C	592	HIS
1	C	625	ASN
1	C	667	GLN
1	C	675	GLN
1	C	696	GLN
1	C	700	ASN
1	C	710	HIS
1	D	496	GLN
1	D	559	GLN
1	D	681	ASN
1	D	712	GLN
1	D	713	GLN
2	G	5	ASN
2	H	5	ASN

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	239/240 (99%)	-0.09	0 100 100	43, 56, 76, 89	0
1	B	234/240 (97%)	-0.07	0 100 100	46, 60, 75, 95	0
1	C	239/240 (99%)	-0.08	0 100 100	61, 74, 92, 108	0
1	D	239/240 (99%)	0.00	1 (0%) 88 85	60, 76, 99, 109	0
2	E	7/7 (100%)	1.41	1 (14%) 6 5	67, 81, 94, 98	0
2	F	7/7 (100%)	0.70	0 100 100	51, 62, 67, 72	0
2	G	7/7 (100%)	0.53	0 100 100	53, 64, 70, 74	0
2	H	7/7 (100%)	0.39	0 100 100	71, 82, 89, 90	0
All	All	979/988 (99%)	-0.04	2 (0%) 91 89	43, 67, 91, 109	0

All (2) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	527	GLU	2.1
2	E	7	PHE	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 ⓘ

There are no such residues in this entry.