



Full wwPDB X-ray Structure Validation Report ⓘ

May 16, 2020 – 02:49 am BST

PDB ID : 3OHN
Title : Crystal structure of the FimD translocation domain
Authors : Wang, T.; Li, H.
Deposited on : 2010-08-17
Resolution : 3.01 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	:	2.11
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)	:	2.11

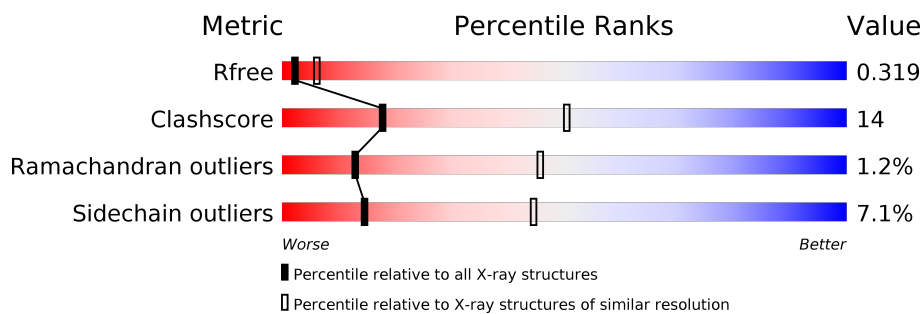
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.01 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	2399 (3.04-3.00)
Clashscore	141614	2734 (3.04-3.00)
Ramachandran outliers	138981	2640 (3.04-3.00)
Sidechain outliers	138945	2643 (3.04-3.00)

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 on 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 $\leq 5\%$.

Mol	Chain	Length	Quality of chain
1	A	558	
1	B	558	

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 7246 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 Outer membrane usher protein FimD.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	468	Total	C	N	O	S	0	0	0
			3636	2271	640	719	6			
1	B	463	Total	C	N	O	S	0	0	0
			3610	2259	634	711	6			

There are 36 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	664	GLY	-	EXPRESSION TAG	UNP C8U0R5
A	665	GLY	-	EXPRESSION TAG	UNP C8U0R5
A	666	PRO	-	EXPRESSION TAG	UNP C8U0R5
A	667	VAL	-	EXPRESSION TAG	UNP C8U0R5
A	668	ALA	-	EXPRESSION TAG	UNP C8U0R5
A	669	THR	-	EXPRESSION TAG	UNP C8U0R5
A	670	LEU	-	EXPRESSION TAG	UNP C8U0R5
A	671	VAL	-	EXPRESSION TAG	UNP C8U0R5
A	672	PRO	-	EXPRESSION TAG	UNP C8U0R5
A	673	ARG	-	EXPRESSION TAG	UNP C8U0R5
A	674	GLY	-	EXPRESSION TAG	UNP C8U0R5
A	675	SER	-	EXPRESSION TAG	UNP C8U0R5
A	676	HIS	-	EXPRESSION TAG	UNP C8U0R5
A	677	HIS	-	EXPRESSION TAG	UNP C8U0R5
A	678	HIS	-	EXPRESSION TAG	UNP C8U0R5
A	679	HIS	-	EXPRESSION TAG	UNP C8U0R5
A	680	HIS	-	EXPRESSION TAG	UNP C8U0R5
A	681	HIS	-	EXPRESSION TAG	UNP C8U0R5
B	664	GLY	-	EXPRESSION TAG	UNP C8U0R5
B	665	GLY	-	EXPRESSION TAG	UNP C8U0R5
B	666	PRO	-	EXPRESSION TAG	UNP C8U0R5
B	667	VAL	-	EXPRESSION TAG	UNP C8U0R5
B	668	ALA	-	EXPRESSION TAG	UNP C8U0R5
B	669	THR	-	EXPRESSION TAG	UNP C8U0R5
B	670	LEU	-	EXPRESSION TAG	UNP C8U0R5

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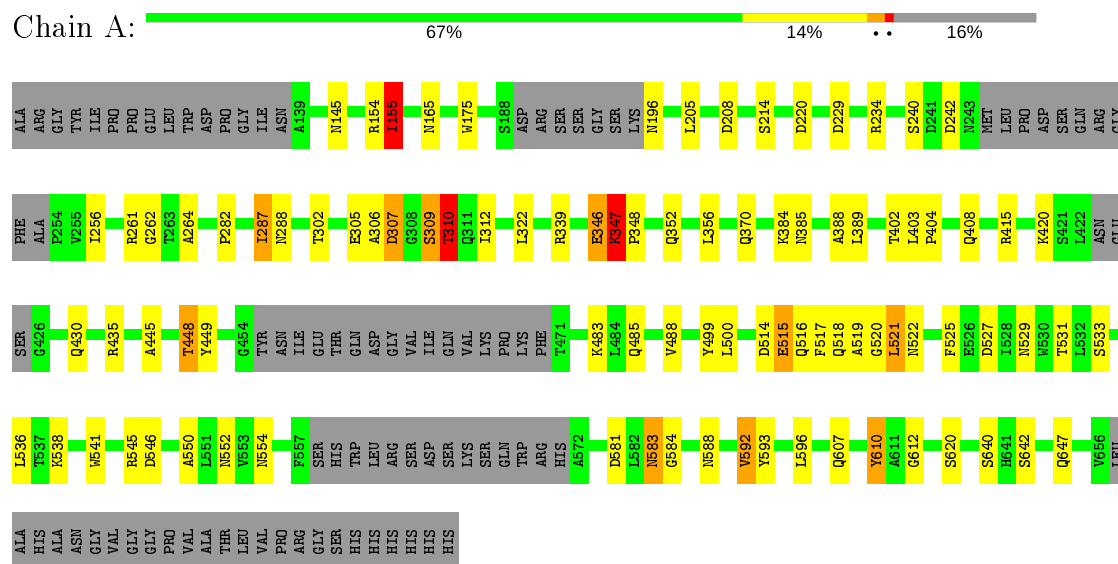
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Chain	Residue	Modelled	Actual	Comment	Reference
B	671	VAL	-	EXPRESSION TAG	UNP C8U0R5
B	672	PRO	-	EXPRESSION TAG	UNP C8U0R5
B	673	ARG	-	EXPRESSION TAG	UNP C8U0R5
B	674	GLY	-	EXPRESSION TAG	UNP C8U0R5
B	675	SER	-	EXPRESSION TAG	UNP C8U0R5
B	676	HIS	-	EXPRESSION TAG	UNP C8U0R5
B	677	HIS	-	EXPRESSION TAG	UNP C8U0R5
B	678	HIS	-	EXPRESSION TAG	UNP C8U0R5
B	679	HIS	-	EXPRESSION TAG	UNP C8U0R5
B	680	HIS	-	EXPRESSION TAG	UNP C8U0R5
B	681	HIS	-	EXPRESSION TAG	UNP C8U0R5

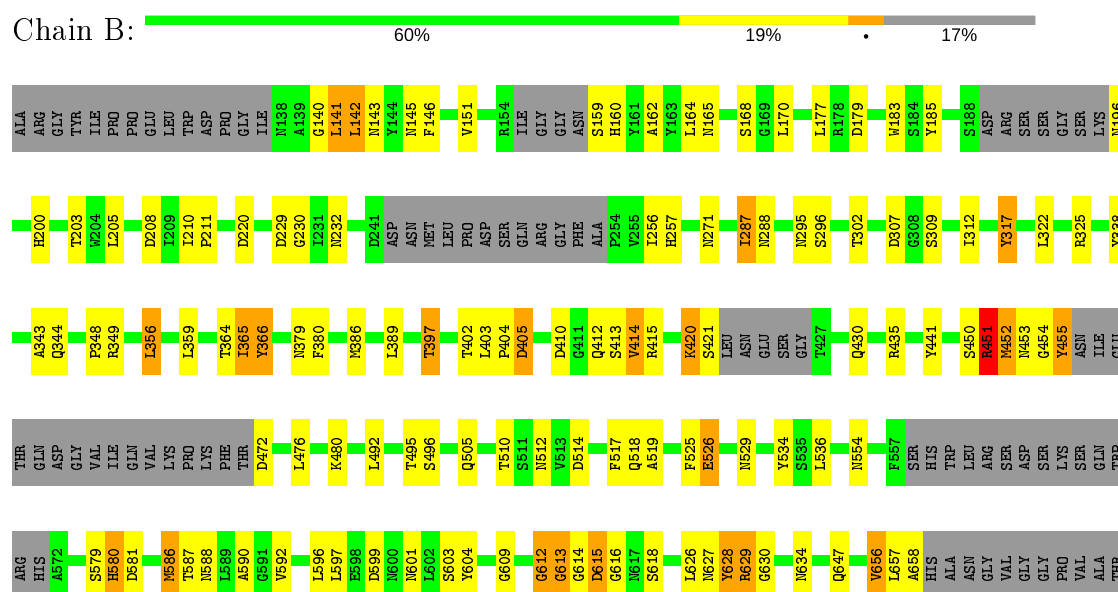
3 Residue-property plots

These plots are drawn for all protein, RNA and DNA 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. 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. 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: Outer membrane usher protein FimD



• Molecule 1: Outer membrane usher protein FimD



LEU
VAL
PRO
ARG
GLY
SER
HIS
HIS
HIS
HIS
HIS

4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	50.02Å 87.13Å 95.88Å 63.92° 88.43° 76.94°	Depositor
Resolution (Å)	28.89 – 3.01 28.89 – 3.01	Depositor EDS
% Data completeness (in resolution range)	98.0 (28.89-3.01) 98.2 (28.89-3.01)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.95 (at 3.00Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.6.1_357)	Depositor
R, R_{free}	0.230 , 0.305 0.309 , 0.319	Depositor DCC
R_{free} test set	1371 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	57.4	Xtriage
Anisotropy	0.629	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 36.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.85	EDS
Total number of atoms	7246	wwPDB-VP
Average B, all atoms (Å ²)	61.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.62% 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	0.52	0/3715	0.57	3/5038 (0.1%)
1	B	0.80	5/3689 (0.1%)	0.64	4/5002 (0.1%)
All	All	0.67	5/7404 (0.1%)	0.60	7/10040 (0.1%)

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	414	VAL	CB-CG1	-7.25	1.37	1.52
1	B	366	TYR	CD2-CE2	-5.90	1.30	1.39
1	B	366	TYR	CE1-CZ	-5.50	1.31	1.38
1	B	366	TYR	CE2-CZ	-5.42	1.31	1.38
1	B	366	TYR	CD1-CE1	-5.32	1.31	1.39

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	612	GLY	N-CA-C	-8.56	91.71	113.10
1	A	310	THR	N-CA-C	6.87	129.55	111.00
1	B	140	GLY	N-CA-C	-6.76	96.19	113.10
1	A	347	LYS	C-N-CD	-6.38	106.56	120.60
1	B	613	GLY	N-CA-C	-6.24	97.51	113.10
1	A	309	SER	C-N-CA	-5.74	107.36	121.70
1	B	307	ASP	N-CA-C	-5.62	95.81	111.00

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	3636	0	3426	63	0
1	B	3610	0	3402	130	0
All	All	7246	0	6828	192	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 14.

All (192) 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:628:TYR:CE2	1:B:630:GLY:N	1.87	1.38
1:B:142:LEU:HD12	1:B:143:ASN:N	1.39	1.36
1:B:141:LEU:HD12	1:B:142:LEU:N	1.43	1.30
1:B:365:ILE:O	1:B:365:ILE:HD12	1.35	1.24
1:B:141:LEU:C	1:B:141:LEU:HD12	1.53	1.21
1:B:142:LEU:HD12	1:B:142:LEU:C	1.53	1.17
1:B:365:ILE:HD12	1:B:365:ILE:C	1.68	1.13
1:B:451:ARG:HA	1:B:451:ARG:HE	1.02	1.10
1:B:403:LEU:HB3	1:B:404:PRO:CD	1.83	1.06
1:B:389:LEU:O	1:B:420:LYS:HD2	1.56	1.05
1:B:403:LEU:HB3	1:B:404:PRO:HD2	1.38	1.03
1:B:455:TYR:HE2	1:B:472:ASP:HB2	1.20	1.02
1:A:583:ASN:HD22	1:A:583:ASN:C	1.60	1.01
1:B:141:LEU:HB2	1:B:657:LEU:CD2	1.90	1.01
1:B:295:ASN:HB2	1:B:317:TYR:OH	1.60	1.00
1:A:521:LEU:C	1:A:521:LEU:CD2	2.29	0.99
1:B:628:TYR:CD2	1:B:629:ARG:N	2.30	0.99
1:B:587:THR:H	1:B:613:GLY:HA2	1.25	0.98
1:B:365:ILE:CD1	1:B:365:ILE:C	2.30	0.97
1:B:141:LEU:CD1	1:B:141:LEU:C	2.30	0.96
1:B:454:GLY:O	1:B:455:TYR:HB3	1.59	0.96
1:B:628:TYR:C	1:B:628:TYR:HD2	1.70	0.96
1:B:141:LEU:CD1	1:B:142:LEU:N	2.30	0.95
1:B:628:TYR:HE2	1:B:630:GLY:N	1.52	0.95
1:B:141:LEU:HB2	1:B:657:LEU:HD23	1.46	0.95
1:B:455:TYR:CE2	1:B:472:ASP:HB2	2.01	0.94
1:B:142:LEU:C	1:B:142:LEU:CD1	2.29	0.94
1:A:521:LEU:C	1:A:521:LEU:HD22	1.88	0.94
1:A:448:THR:O	1:A:448:THR:HG23	1.68	0.94

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:451:ARG:HA	1:B:451:ARG:NE	1.84	0.93
1:B:628:TYR:CD2	1:B:628:TYR:C	2.36	0.92
1:B:450:SER:O	1:B:451:ARG:HB2	1.70	0.92
1:B:141:LEU:HD12	1:B:142:LEU:CA	2.01	0.90
1:A:521:LEU:CD2	1:A:522:ASN:N	2.34	0.90
1:B:455:TYR:HD2	1:B:455:TYR:C	1.74	0.90
1:A:309:SER:O	1:A:310:THR:HG23	1.73	0.88
1:A:521:LEU:HD23	1:A:522:ASN:N	1.88	0.88
1:A:592:VAL:HG11	1:A:607:GLN:HA	1.55	0.86
1:B:159:SER:C	1:B:160:HIS:HD2	1.80	0.83
1:B:455:TYR:CD2	1:B:455:TYR:C	2.47	0.83
1:B:628:TYR:O	1:B:634:ASN:HB2	1.79	0.82
1:B:142:LEU:CD1	1:B:143:ASN:N	2.34	0.82
1:B:451:ARG:CA	1:B:451:ARG:HE	1.85	0.82
1:B:601:ASN:O	1:B:629:ARG:NE	2.14	0.80
1:B:452:MET:HG2	1:B:452:MET:O	1.81	0.79
1:A:499:TYR:CZ	1:A:520:GLY:HA3	2.18	0.78
1:B:159:SER:C	1:B:160:HIS:CD2	2.57	0.78
1:B:403:LEU:HD22	1:B:476:LEU:HD23	1.64	0.78
1:B:405:ASP:OD1	1:B:451:ARG:NH2	2.16	0.78
1:B:601:ASN:O	1:B:629:ARG:CZ	2.31	0.78
1:B:414:VAL:CG1	1:B:414:VAL:O	2.30	0.77
1:B:454:GLY:O	1:B:455:TYR:CB	2.32	0.77
1:B:612:GLY:HA2	1:B:618:SER:HA	1.65	0.77
1:B:414:VAL:HG13	1:B:414:VAL:O	1.83	0.76
1:B:629:ARG:N	1:B:629:ARG:HD3	2.00	0.75
1:A:583:ASN:C	1:A:583:ASN:ND2	2.30	0.75
1:B:450:SER:O	1:B:451:ARG:CB	2.33	0.74
1:B:586:MET:CE	1:B:612:GLY:O	2.37	0.73
1:A:592:VAL:HG22	1:A:593:TYR:H	1.54	0.73
1:B:628:TYR:HD2	1:B:629:ARG:N	1.80	0.73
1:B:657:LEU:O	1:B:658:ALA:HB3	1.89	0.72
1:B:403:LEU:CB	1:B:404:PRO:CD	2.62	0.72
1:B:210:ILE:HG13	1:B:211:PRO:HD3	1.71	0.71
1:A:521:LEU:HD23	1:A:522:ASN:H	1.56	0.71
1:B:379:ASN:OD1	1:B:379:ASN:C	2.29	0.71
1:A:640:SER:HB3	1:A:647:GLN:HB2	1.71	0.70
1:B:628:TYR:CE2	1:B:630:GLY:CA	2.75	0.70
1:B:586:MET:HE1	1:B:612:GLY:O	1.92	0.69
1:A:521:LEU:HD22	1:A:522:ASN:N	2.04	0.69
1:A:499:TYR:CE2	1:A:520:GLY:HA3	2.28	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:159:SER:O	1:B:160:HIS:CD2	2.47	0.67
1:B:356:LEU:HG	1:B:366:TYR:HB3	1.78	0.66
1:B:628:TYR:OH	1:B:630:GLY:HA3	1.96	0.66
1:B:601:ASN:HA	1:B:629:ARG:HE	1.61	0.65
1:B:529:ASN:HB3	1:B:554:ASN:HB2	1.78	0.65
1:B:627:ASN:OD1	1:B:628:TYR:N	2.30	0.65
1:B:453:ASN:OD1	1:B:454:GLY:N	2.30	0.65
1:A:346:GLU:OE1	1:A:347:LYS:N	2.29	0.65
1:B:379:ASN:OD1	1:B:380:PHE:N	2.30	0.64
1:A:309:SER:O	1:A:310:THR:CG2	2.44	0.64
1:A:287:ILE:HD12	1:A:288:ASN:H	1.63	0.63
1:B:629:ARG:HD3	1:B:629:ARG:H	1.62	0.63
1:B:142:LEU:HD12	1:B:143:ASN:CA	2.25	0.63
1:A:261:ARG:HB2	1:A:305:GLU:OE2	1.99	0.63
1:A:264:ALA:HB2	1:A:305:GLU:OE1	1.98	0.62
1:B:586:MET:HG2	1:B:613:GLY:HA3	1.81	0.62
1:B:614:GLY:O	1:B:616:GLY:N	2.33	0.61
1:A:306:ALA:O	1:A:307:ASP:CB	2.48	0.61
1:A:583:ASN:O	1:A:583:ASN:ND2	2.30	0.61
1:A:483:LYS:NZ	1:A:485:GLN:HG2	2.17	0.60
1:A:261:ARG:NE	1:A:583:ASN:OD1	2.35	0.60
1:A:483:LYS:HZ2	1:A:485:GLN:HG3	1.66	0.59
1:B:657:LEU:O	1:B:658:ALA:CB	2.49	0.59
1:B:628:TYR:CZ	1:B:630:GLY:N	2.61	0.59
1:B:629:ARG:N	1:B:629:ARG:CD	2.66	0.58
1:B:579:SER:C	1:B:580:HIS:ND1	2.57	0.58
1:A:529:ASN:HB3	1:A:554:ASN:HB2	1.85	0.58
1:B:287:ILE:HD12	1:B:288:ASN:H	1.67	0.58
1:A:306:ALA:O	1:A:307:ASP:HB2	2.03	0.58
1:B:581:ASP:HB3	1:B:588:ASN:HB3	1.86	0.57
1:B:379:ASN:HB2	1:B:397:THR:HG23	1.86	0.57
1:B:587:THR:N	1:B:613:GLY:HA2	2.09	0.57
1:B:517:PHE:HB3	1:B:536:LEU:HB3	1.87	0.57
1:B:295:ASN:HB2	1:B:317:TYR:CZ	2.39	0.57
1:A:483:LYS:NZ	1:A:485:GLN:CG	2.67	0.57
1:A:430:GLN:HG2	1:A:488:VAL:HG12	1.86	0.56
1:B:403:LEU:HB3	1:B:404:PRO:HD3	1.81	0.56
1:A:483:LYS:HZ1	1:A:485:GLN:HG2	1.70	0.55
1:A:384:LYS:HG2	1:A:385:ASN:N	2.22	0.55
1:B:404:PRO:HB3	1:B:450:SER:HA	1.87	0.55
1:B:628:TYR:CZ	1:B:630:GLY:CA	2.90	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:322:LEU:HD23	1:B:356:LEU:HD21	1.89	0.54
1:B:403:LEU:HD22	1:B:476:LEU:CD2	2.37	0.54
1:B:168:SER:HB3	1:B:179:ASP:HB3	1.89	0.54
1:A:517:PHE:HB3	1:A:536:LEU:HB3	1.90	0.54
1:B:629:ARG:HA	1:B:634:ASN:HB3	1.91	0.52
1:B:257:HIS:HB2	1:B:518:GLN:HG2	1.92	0.52
1:B:628:TYR:OH	1:B:630:GLY:CA	2.58	0.52
1:B:145:ASN:HB3	1:B:165:ASN:HB2	1.92	0.51
1:B:397:THR:O	1:B:412:GLN:HA	2.10	0.51
1:B:586:MET:HE2	1:B:612:GLY:O	2.07	0.51
1:B:170:LEU:HB3	1:B:177:LEU:HB3	1.93	0.51
1:B:343:ALA:O	1:B:344:GLN:HG2	2.10	0.51
1:B:580:HIS:N	1:B:580:HIS:ND1	2.58	0.51
1:B:628:TYR:CD2	1:B:629:ARG:CA	2.93	0.51
1:B:519:ALA:HB3	1:B:534:TYR:HB3	1.92	0.51
1:B:590:ALA:HA	1:B:609:GLY:HA3	1.94	0.50
1:B:160:HIS:N	1:B:160:HIS:CD2	2.79	0.50
1:B:492:LEU:HB2	1:B:496:SER:HB2	1.94	0.50
1:B:338:TYR:HB3	1:B:348:PRO:HD2	1.93	0.50
1:B:628:TYR:CZ	1:B:630:GLY:HA3	2.45	0.50
1:B:453:ASN:OD1	1:B:454:GLY:O	2.30	0.49
1:B:603:SER:OG	1:B:629:ARG:NH1	2.45	0.49
1:A:402:THR:HG22	1:A:408:GLN:HG2	1.94	0.49
1:B:141:LEU:O	1:B:141:LEU:HD12	2.02	0.49
1:A:500:LEU:HD12	1:A:519:ALA:HB2	1.94	0.49
1:B:656:VAL:CG1	1:B:657:LEU:N	2.72	0.49
1:A:445:ALA:O	1:A:448:THR:HG23	2.12	0.48
1:A:581:ASP:HB2	1:A:588:ASN:HB3	1.95	0.48
1:B:151:VAL:HG12	1:B:647:GLN:HG2	1.96	0.48
1:B:232:ASN:HD22	1:B:271:ASN:HD22	1.62	0.48
1:B:413:SER:HB2	1:B:441:TYR:CZ	2.48	0.48
1:A:154:ARG:O	1:A:155:ILE:HG13	2.14	0.48
1:B:579:SER:O	1:B:579:SER:OG	2.30	0.48
1:A:448:THR:O	1:A:448:THR:CG2	2.38	0.47
1:A:515:GLU:HB2	1:A:538:LYS:HB3	1.95	0.47
1:A:262:GLY:N	1:A:305:GLU:OE2	2.41	0.47
1:B:656:VAL:HG13	1:B:657:LEU:N	2.28	0.47
1:A:261:ARG:CZ	1:A:583:ASN:OD1	2.62	0.47
1:A:388:ALA:O	1:A:389:LEU:HB2	2.13	0.47
1:A:533:SER:HB2	1:A:550:ALA:HB3	1.95	0.47
1:A:610:TYR:HA	1:A:620:SER:HA	1.96	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:420:LYS:O	1:B:421:SER:OG	2.30	0.47
1:B:628:TYR:O	1:B:634:ASN:CB	2.57	0.47
1:A:415:ARG:HD2	1:A:435:ARG:HH21	1.78	0.47
1:A:583:ASN:O	1:A:584:GLY:C	2.50	0.46
1:A:352:GLN:HG3	1:A:370:GLN:HB3	1.97	0.46
1:B:628:TYR:CD2	1:B:630:GLY:N	2.61	0.46
1:A:420:LYS:HB3	1:A:430:GLN:HB3	1.98	0.46
1:A:196:ASN:HD21	1:B:196:ASN:HB3	1.82	0.45
1:B:495:THR:HG22	1:B:525:PHE:HB3	1.98	0.45
1:B:162:ALA:HB3	1:B:185:TYR:HB3	1.99	0.45
1:A:322:LEU:HD23	1:A:356:LEU:HD21	1.99	0.44
1:A:445:ALA:O	1:A:448:THR:CG2	2.65	0.44
1:A:262:GLY:HA2	1:A:282:PRO:HB3	2.00	0.43
1:B:302:THR:HG22	1:B:312:ILE:HG12	1.99	0.43
1:B:597:LEU:C	1:B:599:ASP:H	2.21	0.43
1:A:516:GLN:HE21	1:A:518:GLN:HG2	1.83	0.43
1:B:455:TYR:HE2	1:B:472:ASP:CB	2.09	0.43
1:A:403:LEU:HB3	1:A:404:PRO:HD2	2.00	0.43
1:A:256:ILE:H	1:A:287:ILE:HD11	1.84	0.43
1:B:601:ASN:HA	1:B:629:ARG:NE	2.31	0.42
1:A:525:PHE:CE2	1:A:527:ASP:HB2	2.54	0.42
1:B:164:LEU:HB3	1:B:183:TRP:HB3	2.02	0.42
1:B:451:ARG:CA	1:B:451:ARG:NE	2.57	0.42
1:A:592:VAL:HG13	1:A:593:TYR:N	2.35	0.42
1:A:145:ASN:HB3	1:A:165:ASN:HB2	2.01	0.42
1:B:287:ILE:H	1:B:287:ILE:HG13	1.63	0.42
1:A:531:THR:HG23	1:A:552:ASN:HB2	2.01	0.41
1:A:234:ARG:HE	1:A:339:ARG:HD3	1.85	0.41
1:B:402:THR:OG1	1:B:402:THR:O	2.30	0.41
1:A:214:SER:HA	1:A:240:SER:HA	2.03	0.41
1:B:230:GLY:HA3	1:B:296:SER:HB2	2.02	0.41
1:B:505:GLN:HB3	1:B:514:ASP:HB3	2.03	0.41
1:A:175:TRP:CE3	1:A:205:LEU:HG	2.55	0.41
1:B:146:PHE:HD1	1:B:164:LEU:HD13	1.86	0.41
1:B:525:PHE:O	1:B:526:GLU:HG2	2.20	0.41
1:A:302:THR:HB	1:A:312:ILE:HG23	2.03	0.41
1:B:510:THR:HG22	1:B:512:ASN:H	1.86	0.41
1:B:203:THR:H	1:B:220:ASP:HB2	1.86	0.40
1:B:604:TYR:HB3	1:B:626:LEU:HD23	2.02	0.40
1:B:141:LEU:CD1	1:B:142:LEU:CA	2.86	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	456/558 (82%)	422 (92%)	29 (6%)	5 (1%)	14	48
1	B	449/558 (80%)	409 (91%)	34 (8%)	6 (1%)	12	43
All	All	905/1116 (81%)	831 (92%)	63 (7%)	11 (1%)	13	46

All (11) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	155	ILE
1	A	592	VAL
1	B	451	ARG
1	A	348	PRO
1	A	541	TRP
1	B	592	VAL
1	B	615	ASP
1	B	325	ARG
1	B	405	ASP
1	B	526	GLU
1	A	612	GLY

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	381/457 (83%)	360 (94%)	21 (6%)	21	55
1	B	378/457 (83%)	345 (91%)	33 (9%)	10	35

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
All	All	759/914 (83%)	705 (93%)	54 (7%)	14	44

All (54) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	155	ILE
1	A	208	ASP
1	A	220	ASP
1	A	229	ASP
1	A	242	ASP
1	A	287	ILE
1	A	307	ASP
1	A	310	THR
1	A	346	GLU
1	A	347	LYS
1	A	448	THR
1	A	449	TYR
1	A	514	ASP
1	A	515	GLU
1	A	521	LEU
1	A	545	ARG
1	A	546	ASP
1	A	583	ASN
1	A	596	LEU
1	A	610	TYR
1	A	642	SER
1	B	141	LEU
1	B	142	LEU
1	B	200	HIS
1	B	205	LEU
1	B	208	ASP
1	B	229	ASP
1	B	256	ILE
1	B	287	ILE
1	B	309	SER
1	B	317	TYR
1	B	349	ARG
1	B	356	LEU
1	B	359	LEU
1	B	364	THR
1	B	365	ILE
1	B	386	MET

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Mol	Chain	Res	Type
1	B	397	THR
1	B	410	ASP
1	B	415	ARG
1	B	420	LYS
1	B	430	GLN
1	B	435	ARG
1	B	451	ARG
1	B	452	MET
1	B	455	TYR
1	B	480	LYS
1	B	580	HIS
1	B	586	MET
1	B	596	LEU
1	B	615	ASP
1	B	628	TYR
1	B	629	ARG
1	B	656	VAL

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (19) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	196	ASN
1	A	199	GLN
1	A	379	ASN
1	A	479	ASN
1	A	491	GLN
1	A	505	GLN
1	A	580	HIS
1	A	627	ASN
1	B	145	ASN
1	B	160	HIS
1	B	165	ASN
1	B	199	GLN
1	B	271	ASN
1	B	479	ASN
1	B	485	GLN
1	B	491	GLN
1	B	504	HIS
1	B	505	GLN
1	B	518	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 carbohydrates 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

Unable to reproduce the depositors R factor - this section is therefore empty.

6.2 Non-standard residues in protein, DNA, RNA chains

Unable to reproduce the depositors R factor - this section is therefore empty.

6.3 Carbohydrates

Unable to reproduce the depositors R factor - this section is therefore empty.

6.4 Ligands

Unable to reproduce the depositors R factor - this section is therefore empty.

6.5 Other polymers

Unable to reproduce the depositors R factor - this section is therefore empty.