



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 18, 2026 – 02:30 AM UTC

PDB ID : 5J5U / pdb_00005j5u
Title : Fjoh_4561 chitin-binding protein
Authors : Koropatkin, N.M.
Deposited on : 2016-04-03
Resolution : 2.30 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

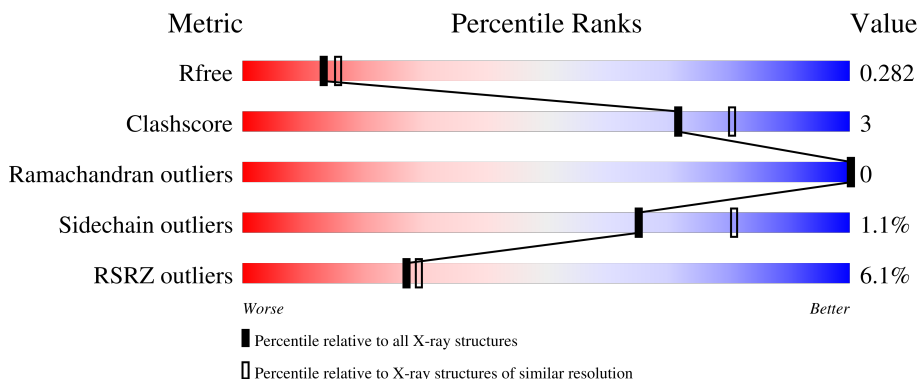
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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	6319 (2.30-2.30)
Clashscore	190562	6919 (2.30-2.30)
Ramachandran outliers	187476	6854 (2.30-2.30)
Sidechain outliers	187428	6854 (2.30-2.30)
RSRZ outliers	180081	6325 (2.30-2.30)

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	484	3% 84% 8% 7%
1	B	484	4% 85% 8% 7%
1	C	484	10% 86% 7% 7%
1	D	484	6% 86% 7% 7%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 14777 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 RagB/SusD domain protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	451	Total	C	N	O	S	0	0	0
			3536	2240	603	680	13			
1	B	452	Total	C	N	O	S	0	0	0
			3544	2246	604	681	13			
1	C	451	Total	C	N	O	S	0	0	0
			3533	2237	603	680	13			
1	D	452	Total	C	N	O	S	0	0	0
			3544	2246	604	681	13			

There are 52 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	22	HIS	-	expression tag	UNP A5FB55
A	23	HIS	-	expression tag	UNP A5FB55
A	24	HIS	-	expression tag	UNP A5FB55
A	25	HIS	-	expression tag	UNP A5FB55
A	26	HIS	-	expression tag	UNP A5FB55
A	27	HIS	-	expression tag	UNP A5FB55
A	28	GLU	-	expression tag	UNP A5FB55
A	29	ASN	-	expression tag	UNP A5FB55
A	30	LEU	-	expression tag	UNP A5FB55
A	31	TYR	-	expression tag	UNP A5FB55
A	32	PHE	-	expression tag	UNP A5FB55
A	33	GLN	-	expression tag	UNP A5FB55
A	34	GLY	-	expression tag	UNP A5FB55
B	22	HIS	-	expression tag	UNP A5FB55
B	23	HIS	-	expression tag	UNP A5FB55
B	24	HIS	-	expression tag	UNP A5FB55
B	25	HIS	-	expression tag	UNP A5FB55
B	26	HIS	-	expression tag	UNP A5FB55
B	27	HIS	-	expression tag	UNP A5FB55
B	28	GLU	-	expression tag	UNP A5FB55
B	29	ASN	-	expression tag	UNP A5FB55

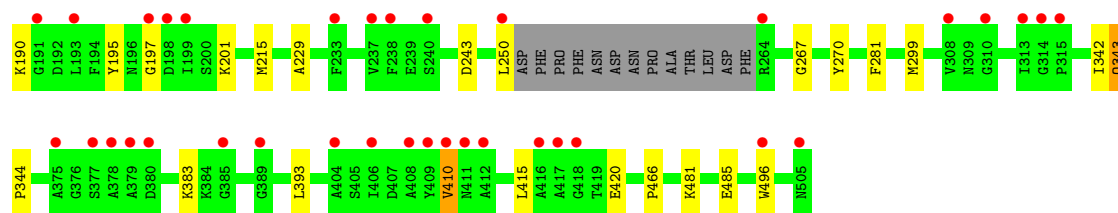
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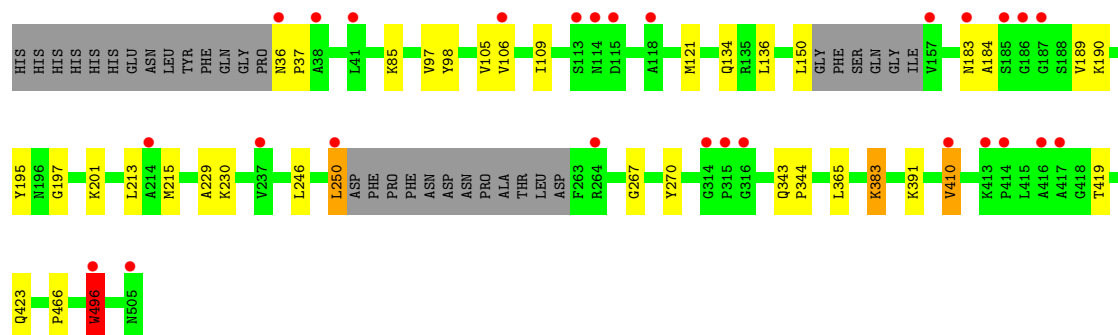
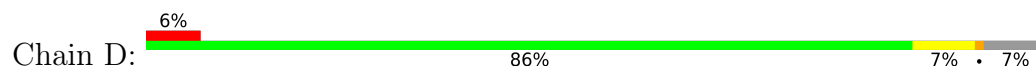
Chain	Residue	Modelled	Actual	Comment	Reference
B	30	LEU	-	expression tag	UNP A5FB55
B	31	TYR	-	expression tag	UNP A5FB55
B	32	PHE	-	expression tag	UNP A5FB55
B	33	GLN	-	expression tag	UNP A5FB55
B	34	GLY	-	expression tag	UNP A5FB55
C	22	HIS	-	expression tag	UNP A5FB55
C	23	HIS	-	expression tag	UNP A5FB55
C	24	HIS	-	expression tag	UNP A5FB55
C	25	HIS	-	expression tag	UNP A5FB55
C	26	HIS	-	expression tag	UNP A5FB55
C	27	HIS	-	expression tag	UNP A5FB55
C	28	GLU	-	expression tag	UNP A5FB55
C	29	ASN	-	expression tag	UNP A5FB55
C	30	LEU	-	expression tag	UNP A5FB55
C	31	TYR	-	expression tag	UNP A5FB55
C	32	PHE	-	expression tag	UNP A5FB55
C	33	GLN	-	expression tag	UNP A5FB55
C	34	GLY	-	expression tag	UNP A5FB55
D	22	HIS	-	expression tag	UNP A5FB55
D	23	HIS	-	expression tag	UNP A5FB55
D	24	HIS	-	expression tag	UNP A5FB55
D	25	HIS	-	expression tag	UNP A5FB55
D	26	HIS	-	expression tag	UNP A5FB55
D	27	HIS	-	expression tag	UNP A5FB55
D	28	GLU	-	expression tag	UNP A5FB55
D	29	ASN	-	expression tag	UNP A5FB55
D	30	LEU	-	expression tag	UNP A5FB55
D	31	TYR	-	expression tag	UNP A5FB55
D	32	PHE	-	expression tag	UNP A5FB55
D	33	GLN	-	expression tag	UNP A5FB55
D	34	GLY	-	expression tag	UNP A5FB55

- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	163	Total O 163 163	0	0
2	B	184	Total O 184 184	0	0
2	C	134	Total O 134 134	0	0
2	D	139	Total O 139 139	0	0



● Molecule 1: RagB/SusD domain protein



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	82.71Å 82.89Å 93.91Å 66.44° 78.33° 67.56°	Depositor
Resolution (Å)	44.49 – 2.30 44.49 – 2.30	Depositor EDS
% Data completeness (in resolution range)	96.8 (44.49-2.30) 95.0 (44.49-2.30)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.90 (at 2.29Å)	Xtriage
Refinement program	PHENIX 1.9_1692	Depositor
R, R_{free}	0.214 , 0.263 (Not available) , 0.282	Depositor DCC
R_{free} test set	1995 reflections (2.19%)	wwPDB-VP
Wilson B-factor (Å ²)	21.2	Xtriage
Anisotropy	0.233	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 45.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.90	EDS
Total number of atoms	14777	wwPDB-VP
Average B, all atoms (Å ²)	26.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 22.32 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 5.8368e-03. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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	0.62	0/3622	0.82	6/4909 (0.1%)
1	B	0.60	0/3630	0.84	2/4920 (0.0%)
1	C	0.58	0/3618	0.85	2/4904 (0.0%)
1	D	0.60	0/3630	0.84	1/4920 (0.0%)
All	All	0.60	0/14500	0.84	11/19653 (0.1%)

There are no bond length outliers.

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	496	TRP	CA-CB-CG	6.48	125.90	113.60
1	A	496	TRP	CA-CB-CG	6.30	125.58	113.60
1	D	496	TRP	CA-CB-CG	5.86	124.73	113.60
1	B	417	ALA	N-CA-C	5.55	117.01	111.07
1	A	221	SER	CA-C-N	5.53	125.63	119.32
1	A	221	SER	C-N-CA	5.53	125.63	119.32
1	A	417	ALA	N-CA-C	5.49	116.94	111.07
1	A	343	GLN	CA-C-N	5.35	124.99	119.05
1	A	343	GLN	C-N-CA	5.35	124.99	119.05
1	C	343	GLN	CA-C-N	5.16	124.95	119.28
1	C	343	GLN	C-N-CA	5.16	124.95	119.28

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	3536	0	3397	26	0
1	B	3544	0	3408	24	0
1	C	3533	0	3399	22	0
1	D	3544	0	3408	22	0
2	A	163	0	0	2	0
2	B	184	0	0	1	0
2	C	134	0	0	0	0
2	D	139	0	0	0	0
All	All	14777	0	13612	92	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 3.

All (92) 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:505:ASN:HD22	1:A:505:ASN:H	1.21	0.88
1:A:505:ASN:HD22	1:A:505:ASN:N	1.78	0.81
1:B:383:LYS:NZ	1:B:415:LEU:H	1.86	0.72
1:B:383:LYS:HG3	1:B:410:VAL:HG13	1.72	0.71
1:C:466:PRO:HB3	1:C:496:TRP:HD1	1.59	0.67
1:B:121:MET:HE1	1:B:190:LYS:H	1.60	0.67
1:D:121:MET:HE1	1:D:190:LYS:H	1.62	0.64
1:A:505:ASN:N	1:A:505:ASN:ND2	2.46	0.63
1:D:383:LYS:HE3	1:D:410:VAL:HG13	1.80	0.63
1:D:183:ASN:OD1	1:D:184:ALA:N	2.31	0.62
1:C:466:PRO:HB3	1:C:496:TRP:CD1	2.34	0.61
1:A:201:LYS:NZ	1:A:243:ASP:OD1	2.35	0.59
1:C:195:TYR:CD2	1:C:201:LYS:HD3	2.36	0.59
1:B:134:GLN:HG3	1:B:213:LEU:HD21	1.85	0.59
1:C:189:VAL:HB	1:C:197:GLY:HA2	1.84	0.59
1:C:85:LYS:HD3	1:C:496:TRP:CE3	2.38	0.59
1:D:419:THR:OG1	1:D:423:GLN:OE1	2.19	0.58
1:A:183:ASN:OD1	1:A:184:ALA:N	2.36	0.58
1:B:379:ALA:O	1:B:383:LYS:HD3	2.04	0.57
1:A:419:THR:OG1	1:A:423:GLN:OE1	2.22	0.57
1:B:383:LYS:HZ2	1:B:415:LEU:H	1.52	0.57
1:C:201:LYS:NZ	1:C:243:ASP:OD1	2.38	0.56
1:A:121:MET:HE1	1:A:190:LYS:H	1.70	0.56
1:A:195:TYR:CD2	1:A:201:LYS:HD3	2.41	0.55
1:D:466:PRO:HB3	1:D:496:TRP:CD1	2.41	0.55
1:B:195:TYR:CD2	1:B:201:LYS:HD3	2.41	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:421:LYS:NZ	2:A:610:HOH:O	2.38	0.55
1:B:419:THR:OG1	1:B:423:GLN:OE1	2.23	0.54
1:A:98:TYR:CZ	1:A:136:LEU:HD21	2.42	0.54
1:D:383:LYS:HG3	1:D:410:VAL:HG13	1.90	0.54
1:D:195:TYR:CD2	1:D:201:LYS:HD3	2.42	0.54
1:A:134:GLN:HG3	1:A:213:LEU:HD21	1.90	0.53
1:C:299:MET:HE2	1:C:393:LEU:HB2	1.92	0.52
1:A:201:LYS:HZ2	1:A:243:ASP:HA	1.74	0.51
1:C:183:ASN:OD1	1:C:184:ALA:N	2.43	0.51
1:A:215:MET:CE	1:A:229:ALA:HB2	2.41	0.51
1:B:215:MET:CE	1:B:229:ALA:HB2	2.41	0.51
1:D:466:PRO:HB3	1:D:496:TRP:HD1	1.75	0.50
1:D:134:GLN:HG3	1:D:213:LEU:HD21	1.94	0.50
1:B:191:GLY:HA2	1:D:250:LEU:HD11	1.93	0.50
1:C:98:TYR:CZ	1:C:136:LEU:HD21	2.48	0.49
1:B:466:PRO:HB3	1:B:496:TRP:CD1	2.47	0.48
1:C:267:GLY:O	1:C:270:TYR:HB3	2.13	0.48
1:C:383:LYS:HG2	1:C:410:VAL:HG13	1.95	0.48
1:D:215:MET:HE3	1:D:229:ALA:HB2	1.96	0.48
1:B:36:ASN:HB2	1:B:37:PRO:HD3	1.95	0.48
1:B:481:LYS:NZ	1:B:485:GLU:OE2	2.44	0.48
1:C:481:LYS:NZ	1:C:485:GLU:OE2	2.47	0.48
1:D:85:LYS:HD3	1:D:496:TRP:CE3	2.48	0.48
1:C:121:MET:HE1	1:C:190:LYS:H	1.78	0.47
1:B:383:LYS:HZ1	1:B:415:LEU:H	1.61	0.47
1:C:215:MET:HE3	1:C:229:ALA:HB2	1.96	0.47
1:C:281:PHE:CE2	1:C:342:ILE:HG12	2.49	0.46
1:D:215:MET:CE	1:D:229:ALA:HB2	2.45	0.46
1:B:189:VAL:HB	1:B:197:GLY:HA2	1.96	0.46
1:C:215:MET:CE	1:C:229:ALA:HB2	2.45	0.46
1:A:198:ASP:OD1	1:A:200:SER:OG	2.32	0.46
1:A:215:MET:HE3	1:A:229:ALA:HB2	1.96	0.46
1:D:189:VAL:HB	1:D:197:GLY:HA2	1.96	0.46
1:D:106:VAL:HG11	1:D:150:LEU:HD23	1.97	0.46
1:B:216:ARG:HA	1:B:445:ARG:HD3	1.98	0.46
1:C:145:TYR:HB3	1:C:162:ASP:OD2	2.15	0.46
1:D:36:ASN:HB2	1:D:37:PRO:HD3	1.97	0.45
1:B:499:LYS:NZ	1:B:505:ASN:HD22	2.14	0.45
1:B:85:LYS:HD3	1:B:496:TRP:CE3	2.52	0.45
1:D:98:TYR:CZ	1:D:136:LEU:HD21	2.50	0.45
1:A:403:GLN:NE2	1:A:407:ASP:OD1	2.47	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:466:PRO:HB3	1:A:496:TRP:CD1	2.52	0.45
1:D:267:GLY:O	1:D:270:TYR:HB3	2.17	0.44
1:C:343:GLN:HA	1:C:344:PRO:HD3	1.81	0.44
1:D:343:GLN:HA	1:D:344:PRO:HD3	1.86	0.44
1:D:365:LEU:HD23	1:D:365:LEU:HA	1.86	0.44
1:D:97:VAL:HG11	1:D:136:LEU:HD22	1.99	0.44
1:C:36:ASN:HB2	1:C:37:PRO:HD3	1.99	0.43
1:A:55:GLN:O	1:A:59:ASN:HB3	2.19	0.43
1:B:51:ASP:OD2	1:B:96:ASN:ND2	2.41	0.43
1:A:208:THR:HG23	1:A:361:SER:HB3	2.00	0.43
1:B:53:TYR:O	1:B:57:ARG:HD2	2.19	0.43
1:B:77:THR:HB	1:B:328:PHE:CE1	2.54	0.43
1:B:55:GLN:O	1:B:59:ASN:HB3	2.18	0.42
1:A:193:LEU:HD21	1:C:250:LEU:HD22	2.01	0.42
1:A:267:GLY:O	1:A:270:TYR:HB3	2.18	0.42
1:A:36:ASN:HB2	1:A:37:PRO:HD3	1.99	0.42
1:A:216:ARG:HA	1:A:445:ARG:HD3	2.00	0.42
1:A:246:LEU:HD13	1:A:354:LEU:HA	2.00	0.42
1:C:415:LEU:HG	1:C:420:GLU:HG3	2.02	0.42
1:A:365:LEU:HD23	1:A:365:LEU:HA	1.89	0.42
1:B:191:GLY:N	2:B:626:HOH:O	2.52	0.41
1:B:281:PHE:CE2	1:B:342:ILE:HG12	2.55	0.41
1:C:85:LYS:HD3	1:C:496:TRP:CD2	2.55	0.41
1:A:392:GLN:NE2	2:A:609:HOH:O	2.36	0.40
1:D:105:VAL:O	1:D:109:ILE:HG13	2.22	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	445/484 (92%)	431 (97%)	14 (3%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	446/484 (92%)	433 (97%)	13 (3%)	0	100	100
1	C	445/484 (92%)	431 (97%)	14 (3%)	0	100	100
1	D	446/484 (92%)	433 (97%)	13 (3%)	0	100	100
All	All	1782/1936 (92%)	1728 (97%)	54 (3%)	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	370/399 (93%)	364 (98%)	6 (2%)	55	73
1	B	371/399 (93%)	369 (100%)	2 (0%)	81	90
1	C	370/399 (93%)	369 (100%)	1 (0%)	86	93
1	D	371/399 (93%)	364 (98%)	7 (2%)	50	69
All	All	1482/1596 (93%)	1466 (99%)	16 (1%)	65	81

All (16) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	185	SER
1	A	246	LEU
1	A	410	VAL
1	A	413	LYS
1	A	496	TRP
1	A	505	ASN
1	B	410	VAL
1	B	496	TRP
1	C	410	VAL
1	D	230	LYS
1	D	246	LEU
1	D	250	LEU
1	D	383	LYS

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Mol	Chain	Res	Type
1	D	391	LYS
1	D	410	VAL
1	D	496	TRP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	505	ASN
1	B	196	ASN
1	B	505	ASN
1	C	473	ASN
1	D	40	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

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	451/484 (93%)	0.22	13 (2%) 53 56	9, 23, 40, 59	0
1	B	452/484 (93%)	0.35	20 (4%) 39 41	9, 23, 43, 57	0
1	C	451/484 (93%)	0.66	50 (11%) 10 11	9, 28, 51, 63	0
1	D	452/484 (93%)	0.46	27 (5%) 27 29	9, 26, 48, 60	0
All	All	1806/1936 (93%)	0.42	110 (6%) 27 29	9, 24, 47, 63	0

All (110) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	250	LEU	5.2
1	D	496	TRP	5.1
1	C	496	TRP	4.8
1	A	250	LEU	4.8
1	A	496	TRP	4.4
1	C	115	ASP	4.2
1	D	113	SER	4.2
1	B	496	TRP	4.0
1	C	380	ASP	4.0
1	A	113	SER	3.9
1	A	505	ASN	3.6
1	C	264	ARG	3.5
1	C	505	ASN	3.3
1	C	199	ILE	3.2
1	C	250	LEU	3.2
1	D	118	ALA	3.1
1	C	406	ILE	3.1
1	B	114	ASN	3.1
1	B	505	ASN	3.1
1	C	193	LEU	3.1
1	C	197	GLY	3.1

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Mol	Chain	Res	Type	RSRZ
1	D	316	GLY	3.0
1	B	186	GLY	3.0
1	C	188	SER	3.0
1	C	114	ASN	3.0
1	D	505	ASN	3.0
1	C	157	VAL	2.9
1	C	314	GLY	2.9
1	C	416	ALA	2.8
1	C	183	ASN	2.8
1	D	115	ASP	2.8
1	D	157	VAL	2.8
1	A	157	VAL	2.8
1	A	118	ALA	2.7
1	C	379	ALA	2.7
1	C	410	VAL	2.7
1	B	115	ASP	2.7
1	B	118	ALA	2.6
1	C	117	ALA	2.6
1	B	410	VAL	2.6
1	A	263	PHE	2.5
1	A	410	VAL	2.5
1	C	237	VAL	2.5
1	D	410	VAL	2.5
1	A	41	LEU	2.5
1	C	375	ALA	2.5
1	D	417	ALA	2.5
1	C	103	LYS	2.5
1	B	157	VAL	2.5
1	C	417	ALA	2.5
1	B	263	PHE	2.5
1	C	310	GLY	2.5
1	D	186	GLY	2.5
1	C	377	SER	2.4
1	B	117	ALA	2.4
1	C	191	GLY	2.4
1	B	36	ASN	2.4
1	C	385	GLY	2.4
1	D	416	ALA	2.4
1	D	36	ASN	2.3
1	C	389	GLY	2.3
1	C	408	ALA	2.3
1	C	240	SER	2.3

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Mol	Chain	Res	Type	RSRZ
1	D	106	VAL	2.3
1	D	114	ASN	2.3
1	D	315	PRO	2.3
1	B	416	ALA	2.3
1	D	237	VAL	2.3
1	C	198	ASP	2.3
1	C	315	PRO	2.3
1	B	385	GLY	2.3
1	C	187	GLY	2.3
1	C	418	GLY	2.3
1	C	409	TYR	2.3
1	D	264	ARG	2.3
1	B	377	SER	2.3
1	C	113	SER	2.3
1	C	313	ILE	2.3
1	B	150	LEU	2.2
1	C	179	PHE	2.2
1	B	250	LEU	2.2
1	C	308	VAL	2.2
1	A	413	LYS	2.2
1	A	417	ALA	2.2
1	D	214	ALA	2.2
1	B	349	ARG	2.2
1	D	183	ASN	2.2
1	B	237	VAL	2.2
1	D	41	LEU	2.1
1	A	36	ASN	2.1
1	B	179	PHE	2.1
1	C	238	PHE	2.1
1	B	197	GLY	2.1
1	D	187	GLY	2.1
1	C	404	ALA	2.1
1	D	185	SER	2.1
1	C	36	ASN	2.1
1	C	39	GLY	2.1
1	D	314	GLY	2.1
1	C	411	ASN	2.1
1	A	187	GLY	2.1
1	C	233	PHE	2.1
1	C	378	ALA	2.0
1	C	412	ALA	2.0
1	C	107	ASP	2.0

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Mol	Chain	Res	Type	RSRZ
1	C	120	ASN	2.0
1	D	414	PRO	2.0
1	C	106	VAL	2.0
1	D	413	LYS	2.0
1	D	38	ALA	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 [i](#)

There are no such residues in this entry.