



Full wwPDB X-ray Structure Validation Report ⓘ

May 23, 2020 – 10:52 am BST

PDB ID : 5J5U
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 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

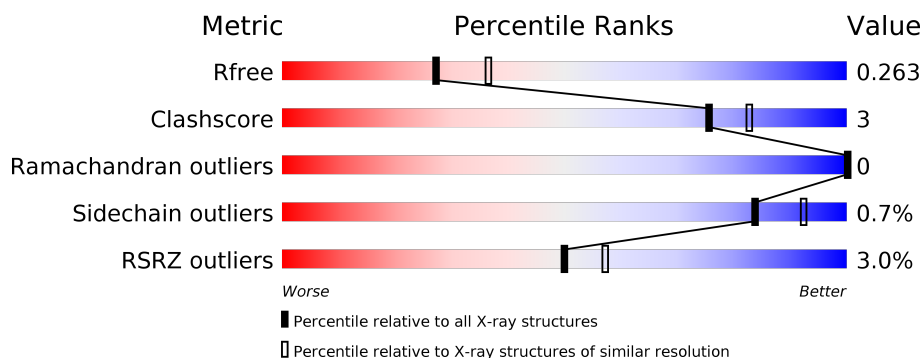
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 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}	130704	5042 (2.30-2.30)
Clashscore	141614	5643 (2.30-2.30)
Ramachandran outliers	138981	5575 (2.30-2.30)
Sidechain outliers	138945	5575 (2.30-2.30)
RSRZ outliers	127900	4938 (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 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 $\leq 5\%$. 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	<div> <div>2%</div> <div> <div></div> <div>86%</div> <div>7%</div> <div>7%</div> </div> </div>
1	B	484	<div> <div>2%</div> <div> <div></div> <div>86%</div> <div>7%</div> <div>7%</div> </div> </div>
1	C	484	<div> <div>4%</div> <div> <div></div> <div>87%</div> <div>6%</div> <div>7%</div> </div> </div>
1	D	484	<div> <div>3%</div> <div> <div></div> <div>86%</div> <div>6%</div> <div>7%</div> </div> </div>

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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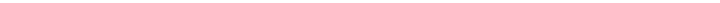
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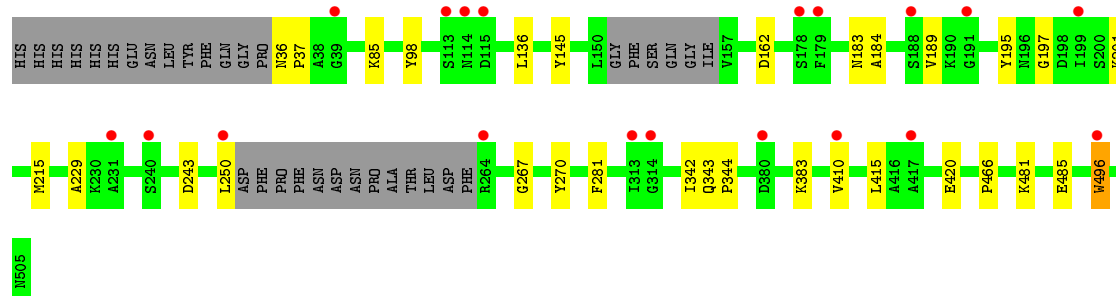
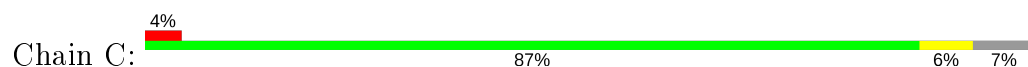
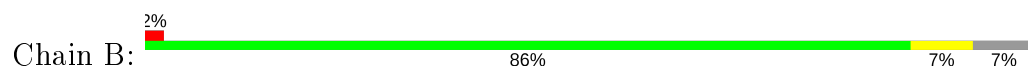
Chain	Residue	Modelled	Actual	Comment	Reference
B	30	LEU	-	expression tag	UNP A5FB55
B	31	TYR	-	expression tag	UNP A5FB55
B	32	PHE	-	expression tag	UNP A5FB55
B	33	GLN	-	expression tag	UNP A5FB55
B	34	GLY	-	expression tag	UNP A5FB55
C	22	HIS	-	expression tag	UNP A5FB55
C	23	HIS	-	expression tag	UNP A5FB55
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.

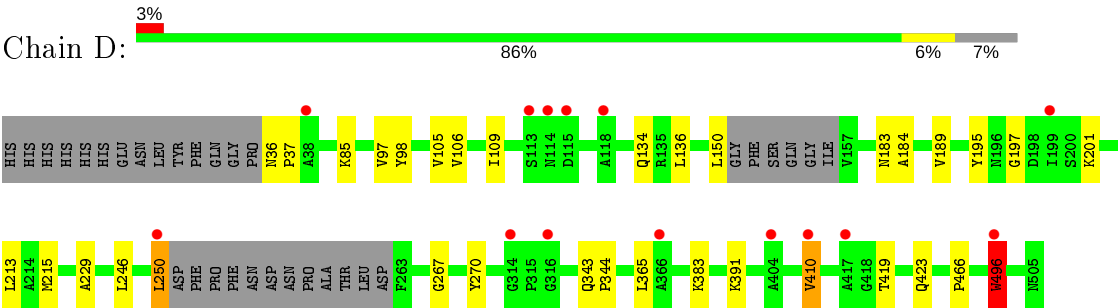
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	163	Total O 163 163	0	0
2	B	184	Total O 184 184	0	0
2	C	134	Total O 134 134	0	0
2	D	139	Total O 139 139	0	0

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 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.

- Chain A:  29% 86% 7% 7%



● 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 85.93 – 2.30	Depositor EDS
% Data completeness (in resolution range)	96.8 (44.49-2.30) 95.0 (85.93-2.30)	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 0.218 , 0.263	Depositor DCC
R_{free} test set	1992 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)	0.36 , 49.9	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	0.91	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.*

¹ Intensities estimated from amplitudes.

² Theoretical 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	0.49	0/3622	0.58	1/4909 (0.0%)
1	B	0.48	0/3630	0.58	1/4920 (0.0%)
1	C	0.47	0/3618	0.57	0/4904
1	D	0.47	0/3630	0.59	2/4920 (0.0%)
All	All	0.48	0/14500	0.58	4/19653 (0.0%)

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	496	TRP	CA-CB-CG	6.42	125.90	113.70
1	A	496	TRP	CA-CB-CG	6.25	125.58	113.70
1	D	496	TRP	CA-CB-CG	5.81	124.73	113.70
1	D	246	LEU	CA-CB-CG	5.20	127.25	115.30

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	24	0
1	B	3544	0	3408	24	0
1	C	3533	0	3399	20	0
1	D	3544	0	3408	20	0
2	A	163	0	0	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
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	86	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 (86) 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:N	1:A:505:ASN:HD22	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: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: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
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: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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:A:215:MET:HE1	1:A:229:ALA:HB2	1.93	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:215:MET:HE1	1:B:229:ALA:HB2	1.95	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: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:B:383:LYS:HZ1	1:B:415:LEU:H	1.61	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: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
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:77:THR:HB	1:B:328:PHE:CE1	2.54	0.43
1:B:53:TYR:O	1:B:57:ARG:HD2	2.19	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:201:LYS:HZ2	1:A:243:ASP:HA	1.83	0.43
1:B:55:GLN:O	1:B:59:ASN:HB3	2.18	0.42
1:A:267:GLY:O	1:A:270:TYR:HB3	2.18	0.42
1:A:193:LEU:HD21	1:C:250:LEU:HD22	2.01	0.42
1:A:216:ARG:HA	1:A:445:ARG:HD3	2.00	0.42
1:A:36:ASN:HB2	1:A:37:PRO:HD3	1.99	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:C:215:MET:HE1	1:C:229:ALA:HB2	2.01	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
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 [i](#)

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%)	367 (99%)	3 (1%)	81	91
1	B	371/399 (93%)	369 (100%)	2 (0%)	88	95
1	C	370/399 (93%)	369 (100%)	1 (0%)	92	97
1	D	371/399 (93%)	367 (99%)	4 (1%)	73	86
All	All	1482/1596 (93%)	1472 (99%)	10 (1%)	84	92

All (10) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	185	SER
1	A	496	TRP
1	A	505	ASN
1	B	349	ARG
1	B	496	TRP
1	C	496	TRP
1	D	250	LEU
1	D	391	LYS
1	D	410	VAL
1	D	496	TRP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	505	ASN
1	B	505	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 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 ⓘ

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%)	0.15	9 (1%) 65 71	9, 23, 40, 59	0
1	B	452/484 (93%)	0.24	12 (2%) 54 62	9, 23, 43, 57	0
1	C	451/484 (93%)	0.39	19 (4%) 36 43	9, 28, 51, 63	0
1	D	452/484 (93%)	0.30	14 (3%) 49 56	9, 26, 48, 60	0
All	All	1806/1936 (93%)	0.27	54 (2%) 50 57	9, 24, 47, 63	0

All (54) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	186	GLY	4.7
1	A	113	SER	4.1
1	C	496	TRP	3.8
1	C	115	ASP	3.7
1	D	114	ASN	3.6
1	D	496	TRP	3.6
1	C	114	ASN	3.4
1	A	496	TRP	3.4
1	B	496	TRP	3.3
1	A	114	ASN	3.3
1	C	39	GLY	3.2
1	A	186	GLY	3.1
1	C	264	ARG	3.1
1	D	410	VAL	3.1
1	C	380	ASP	2.9
1	C	314	GLY	2.9
1	D	113	SER	2.8
1	D	250	LEU	2.8
1	C	199	ILE	2.8
1	A	157	VAL	2.7
1	C	188	SER	2.7

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Mol	Chain	Res	Type	RSRZ
1	B	114	ASN	2.7
1	C	178	SER	2.7
1	B	263	PHE	2.7
1	C	231	ALA	2.6
1	D	38	ALA	2.6
1	D	366	ALA	2.6
1	D	115	ASP	2.5
1	D	316	GLY	2.5
1	A	263	PHE	2.5
1	A	118	ALA	2.5
1	C	240	SER	2.5
1	D	199	ILE	2.5
1	D	118	ALA	2.5
1	D	417	ALA	2.4
1	A	250	LEU	2.4
1	C	179	PHE	2.4
1	A	505	ASN	2.4
1	C	250	LEU	2.3
1	C	113	SER	2.3
1	B	199	ILE	2.3
1	C	191	GLY	2.2
1	B	191	GLY	2.2
1	C	417	ALA	2.2
1	B	193	LEU	2.2
1	C	313	ILE	2.1
1	B	113	SER	2.1
1	B	115	ASP	2.1
1	D	404	ALA	2.1
1	B	39	GLY	2.1
1	C	410	VAL	2.0
1	B	505	ASN	2.0
1	B	250	LEU	2.0
1	D	314	GLY	2.0

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

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

6.4 Ligands

There are no ligands in this entry.

6.5 Other polymers

There are no such residues in this entry.