



# Full wwPDB X-ray Structure Validation Report ⓘ

Apr 19, 2021 – 08:47 AM JST

PDB ID : 7E7W  
Title : Crystal structure of RSL mutant in complex with sugar Ligand  
Authors : Li, L.; Chen, G.S.  
Deposited on : 2021-02-28  
Resolution : 4.10 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](mailto: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.

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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 : 2.18  
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) : 2.18

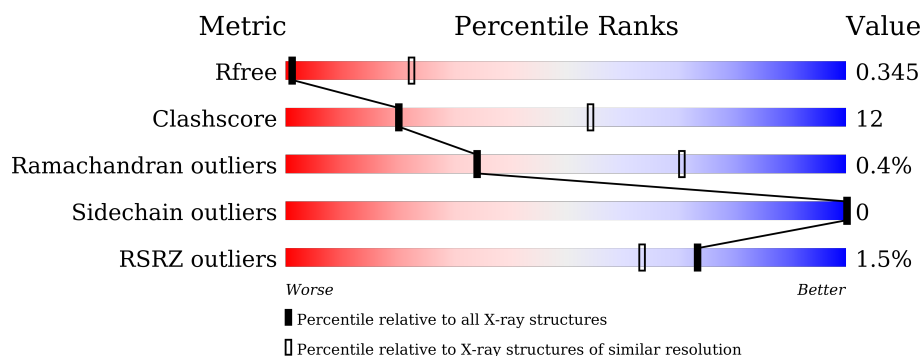
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 4.10 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	1193 (4.50-3.70)
Clashscore	141614	1003 (4.44-3.76)
Ramachandran outliers	138981	1005 (4.48-3.72)
Sidechain outliers	138945	1199 (4.50-3.70)
RSRZ outliers	127900	1034 (4.50-3.70)

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  $\geq 3$ , 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	272	<div> <div>2%</div> <div> <div></div> <div>76%</div> <div>21%</div> <div></div> </div> <div>.</div> </div>
1	B	272	<div> <div>2%</div> <div> <div></div> <div>72%</div> <div>25%</div> <div></div> </div> <div>.</div> </div>
1	C	272	<div> <div></div> <div> <div></div> <div>72%</div> <div>25%</div> <div></div> </div> <div>.</div> </div>

## 2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 6063 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 Fucose-binding lectin protein,Fucose-binding lectin protein,Fucose-binding lectin protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	264	Total	C	N	O	S	0	0	0
			2021	1271	345	399	6			
1	B	264	Total	C	N	O	S	0	0	0
			2021	1271	345	399	6			
1	C	264	Total	C	N	O	S	0	0	0
			2021	1271	345	399	6			

There are 42 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	88	SER	-	linker	UNP A0A0S4TLR1
A	89	SER	-	linker	UNP A0A0S4TLR1
A	90	THR	-	linker	UNP A0A0S4TLR1
A	91	VAL	-	linker	UNP A0A0S4TLR1
A	92	PRO	-	linker	UNP A0A0S4TLR1
A	93	GLY	-	linker	UNP A0A0S4TLR1
A	94	ASP	-	linker	UNP A0A0S4TLR1
A	179	SER	-	linker	UNP A0A0S4TLR1
A	180	SER	-	linker	UNP A0A0S4TLR1
A	181	THR	-	linker	UNP A0A0S4TLR1
A	182	VAL	-	linker	UNP A0A0S4TLR1
A	183	PRO	-	linker	UNP A0A0S4TLR1
A	184	GLY	-	linker	UNP A0A0S4TLR1
A	185	ASP	-	linker	UNP A0A0S4TLR1
B	88	SER	-	linker	UNP A0A0S4TLR1
B	89	SER	-	linker	UNP A0A0S4TLR1
B	90	THR	-	linker	UNP A0A0S4TLR1
B	91	VAL	-	linker	UNP A0A0S4TLR1
B	92	PRO	-	linker	UNP A0A0S4TLR1
B	93	GLY	-	linker	UNP A0A0S4TLR1
B	94	ASP	-	linker	UNP A0A0S4TLR1
B	179	SER	-	linker	UNP A0A0S4TLR1

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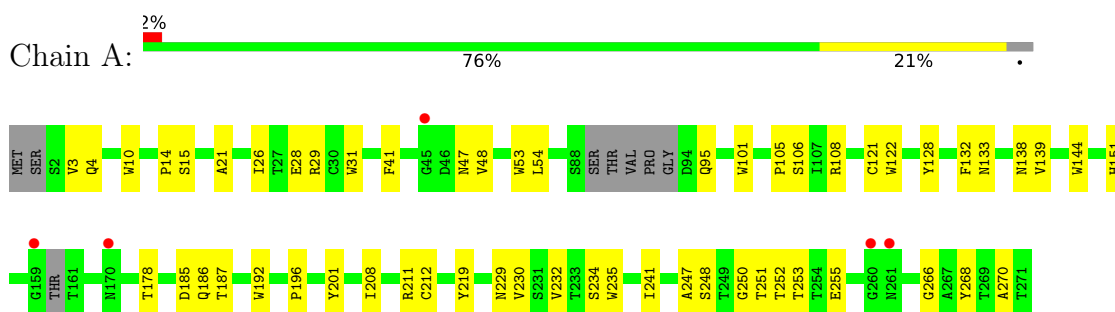
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Chain	Residue	Modelled	Actual	Comment	Reference
B	180	SER	-	linker	UNP A0A0S4TLR1
B	181	THR	-	linker	UNP A0A0S4TLR1
B	182	VAL	-	linker	UNP A0A0S4TLR1
B	183	PRO	-	linker	UNP A0A0S4TLR1
B	184	GLY	-	linker	UNP A0A0S4TLR1
B	185	ASP	-	linker	UNP A0A0S4TLR1
C	88	SER	-	linker	UNP A0A0S4TLR1
C	89	SER	-	linker	UNP A0A0S4TLR1
C	90	THR	-	linker	UNP A0A0S4TLR1
C	91	VAL	-	linker	UNP A0A0S4TLR1
C	92	PRO	-	linker	UNP A0A0S4TLR1
C	93	GLY	-	linker	UNP A0A0S4TLR1
C	94	ASP	-	linker	UNP A0A0S4TLR1
C	179	SER	-	linker	UNP A0A0S4TLR1
C	180	SER	-	linker	UNP A0A0S4TLR1
C	181	THR	-	linker	UNP A0A0S4TLR1
C	182	VAL	-	linker	UNP A0A0S4TLR1
C	183	PRO	-	linker	UNP A0A0S4TLR1
C	184	GLY	-	linker	UNP A0A0S4TLR1
C	185	ASP	-	linker	UNP A0A0S4TLR1

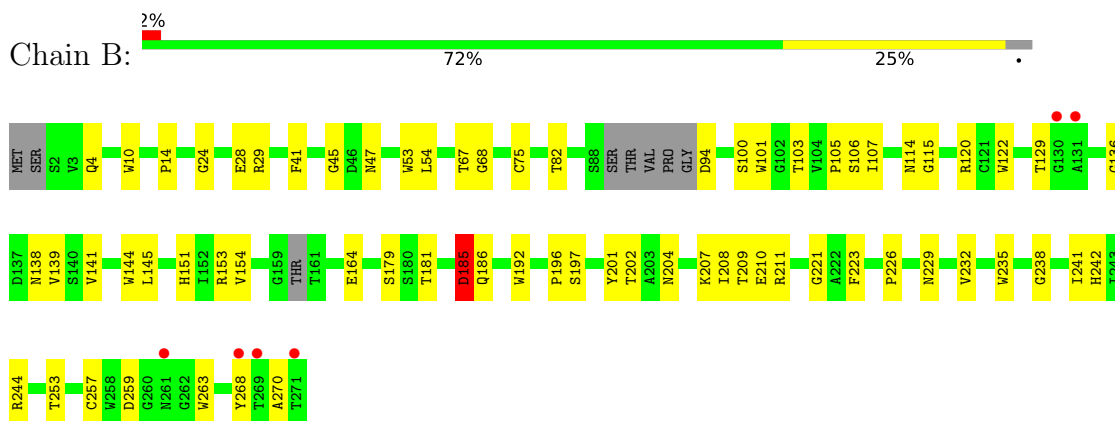
### 3 Residue-property plots [i](#)

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.

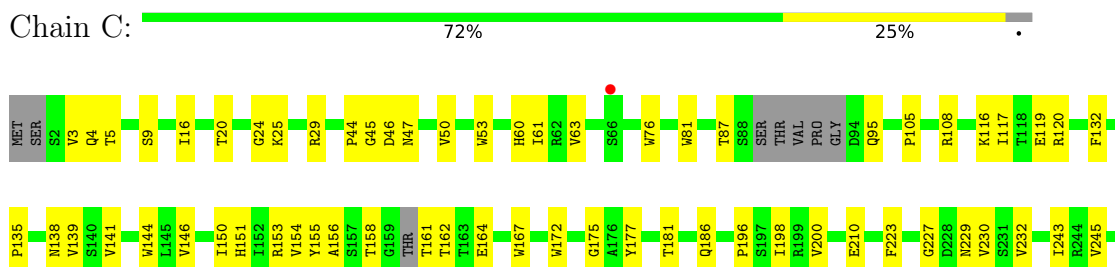
- Molecule 1: Fucose-binding lectin protein,Fucose-binding lectin protein,Fucose-binding lectin protein



- Molecule 1: Fucose-binding lectin protein,Fucose-binding lectin protein,Fucose-binding lectin protein



- Molecule 1: Fucose-binding lectin protein,Fucose-binding lectin protein,Fucose-binding lectin protein



T249	G250	E255	W258	G266	A267	Y268	T269	A270	T271
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## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	46.57Å 46.55Å 103.25Å 90.04° 89.92° 119.98°	Depositor
Resolution (Å)	26.20 – 4.10 26.20 – 4.02	Depositor EDS
% Data completeness (in resolution range)	98.4 (26.20-4.10) 96.5 (26.20-4.02)	Depositor EDS
$R_{merge}$	0.12	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.52 (at 3.97Å)	Xtriage
Refinement program	PHENIX 1.18.2_3874	Depositor
R, $R_{free}$	0.200 , 0.347 0.201 , 0.345	Depositor DCC
$R_{free}$ test set	259 reflections (4.26%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	69.8	Xtriage
Anisotropy	0.485	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.32 , 18.7	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.51$ , $\langle L^2 \rangle = 0.35$	Xtriage
Estimated twinning fraction	0.008 for h+k,-h,l 0.008 for -k,h+k,l 0.398 for k,-h-k,l 0.398 for -h-k,h,l 0.025 for -h-k,k,-l 0.028 for h,-h-k,-l 0.010 for -h,-k,l 0.017 for -k,-h,-l 0.028 for k,h,-l 0.009 for h+k,-k,-l 0.018 for -h,h+k,-l	Xtriage
$F_o, F_c$ correlation	0.91	EDS
Total number of atoms	6063	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	82.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 17.45% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>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 [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	0.27	0/2085	0.49	0/2865
1	B	0.26	0/2085	0.51	1/2865 (0.0%)
1	C	0.26	0/2085	0.49	0/2865
All	All	0.27	0/6255	0.50	1/8595 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	185	ASP	CB-CG-OD1	5.13	122.92	118.30

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	2021	0	1856	44	0
1	B	2021	0	1856	52	0
1	C	2021	0	1856	50	0
All	All	6063	0	5568	145	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 12.

All (145) 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:158:THR:HG1	1:C:161:THR:N	1.63	0.97
1:B:181:THR:HG21	1:B:202:THR:HG21	1.58	0.85
1:A:47:ASN:HD22	1:A:95:GLN:HA	1.40	0.84
1:A:138:ASN:HD22	1:A:186:GLN:HA	1.42	0.83
1:A:4:GLN:HA	1:A:229:ASN:HD22	1.50	0.75
1:B:244:ARG:HH11	1:B:257:CYS:HB3	1.52	0.74
1:B:138:ASN:HD22	1:B:186:GLN:HA	1.53	0.74
1:C:116:LYS:HA	1:C:135:PRO:HA	1.70	0.74
1:C:158:THR:OG1	1:C:161:THR:N	2.22	0.72
1:B:138:ASN:ND2	1:B:186:GLN:HA	2.07	0.70
1:C:5:THR:HG22	1:C:20:THR:HG23	1.73	0.69
1:A:10:TRP:HB2	1:A:54:LEU:HD11	1.74	0.69
1:A:47:ASN:ND2	1:A:95:GLN:HA	2.09	0.68
1:B:29:ARG:HE	1:B:268:TYR:HE1	1.40	0.67
1:C:87:THR:OG1	1:C:120:ARG:NH2	2.27	0.67
1:A:138:ASN:OD1	1:A:139:VAL:N	2.27	0.67
1:A:252:THR:HG23	1:A:270:ALA:HB1	1.78	0.66
1:A:29:ARG:HE	1:A:268:TYR:HE1	1.44	0.65
1:B:120:ARG:NH1	1:B:129:THR:OG1	2.28	0.65
1:B:4:GLN:HG2	1:B:229:ASN:HD21	1.61	0.65
1:C:24:GLY:HA2	1:C:46:ASP:HA	1.80	0.64
1:C:47:ASN:HD21	1:C:95:GLN:HG2	1.63	0.63
1:B:103:THR:HG21	1:C:24:GLY:HA3	1.81	0.62
1:A:4:GLN:HA	1:A:229:ASN:ND2	2.15	0.62
1:A:138:ASN:ND2	1:A:186:GLN:HG2	2.15	0.61
1:C:164:GLU:HB3	1:C:175:GLY:HA3	1.83	0.61
1:B:14:PRO:HG3	1:B:235:TRP:CE2	2.37	0.59
1:A:253:THR:N	1:A:270:ALA:O	2.35	0.58
1:B:138:ASN:ND2	1:B:186:GLN:OE1	2.36	0.58
1:B:138:ASN:OD1	1:B:139:VAL:N	2.36	0.58
1:B:210:GLU:HG2	1:B:221:GLY:HA3	1.85	0.57
1:C:162:THR:H	1:C:181:THR:HG21	1.68	0.57
1:C:138:ASN:ND2	1:C:186:GLN:HG2	2.18	0.57
1:B:144:TRP:CE2	1:B:196:PRO:HG3	2.40	0.56
1:B:53:TRP:CZ2	1:B:105:PRO:HG3	2.40	0.56
1:B:144:TRP:CZ2	1:B:196:PRO:HG3	2.41	0.56
1:A:229:ASN:HB3	1:A:248:SER:HB2	1.87	0.56
1:B:101:TRP:CD1	1:B:145:LEU:HD21	2.40	0.56
1:C:25:LYS:HA	1:C:44:PRO:HA	1.87	0.55
1:C:138:ASN:OD1	1:C:139:VAL:N	2.40	0.55
1:C:138:ASN:HD22	1:C:186:GLN:HA	1.72	0.55
1:A:47:ASN:OD1	1:A:48:VAL:N	2.40	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:24:GLY:O	1:C:45:GLY:N	2.40	0.54
1:A:14:PRO:HG3	1:A:235:TRP:CE2	2.43	0.54
1:A:229:ASN:OD1	1:A:230:VAL:N	2.40	0.54
1:A:26:ILE:HD13	1:A:48:VAL:HB	1.90	0.54
1:C:255:GLU:HB2	1:C:268:TYR:HB2	1.90	0.54
1:C:164:GLU:HB2	1:C:177:TYR:HB3	1.89	0.54
1:B:202:THR:OG1	1:B:211:ARG:NH1	2.36	0.53
1:B:24:GLY:O	1:B:45:GLY:N	2.33	0.52
1:C:232:VAL:HG23	1:C:245:VAL:HG22	1.91	0.52
1:A:144:TRP:CE2	1:A:196:PRO:HG3	2.43	0.52
1:B:144:TRP:CE2	1:B:151:HIS:HB2	2.45	0.52
1:C:229:ASN:OD1	1:C:230:VAL:N	2.43	0.52
1:C:132:PHE:HB2	1:C:172:TRP:NE1	2.24	0.52
1:C:144:TRP:CD1	1:C:146:VAL:HG23	2.45	0.52
1:A:21:ALA:HB2	1:A:26:ILE:HG12	1.92	0.51
1:B:28:GLU:HB2	1:B:41:PHE:HB3	1.92	0.51
1:B:186:GLN:HG2	1:B:229:ASN:OD1	2.10	0.51
1:C:255:GLU:HB3	1:C:266:GLY:HA3	1.92	0.51
1:B:45:GLY:HA2	1:B:67:THR:HG23	1.91	0.51
1:B:235:TRP:O	1:B:241:ILE:HG13	2.11	0.51
1:A:132:PHE:O	1:A:133:ASN:ND2	2.44	0.50
1:B:192:TRP:CD2	1:B:241:ILE:HD13	2.47	0.50
1:C:117:ILE:HD13	1:C:139:VAL:HB	1.93	0.50
1:B:4:GLN:OE1	1:B:47:ASN:ND2	2.46	0.49
1:B:181:THR:OG1	1:B:185:ASP:OD2	2.24	0.49
1:B:75:CYS:HB2	1:B:82:THR:OG1	2.12	0.49
1:A:178:THR:O	1:A:211:ARG:NE	2.46	0.49
1:C:53:TRP:NE1	1:C:60:HIS:HB2	2.28	0.49
1:C:53:TRP:CE2	1:C:60:HIS:HB2	2.48	0.49
1:B:209:THR:OG1	1:B:211:ARG:NH2	2.46	0.48
1:C:141:VAL:HG23	1:C:154:VAL:HG22	1.94	0.48
1:B:68:GLY:O	1:B:94:ASP:HB2	2.12	0.48
1:A:3:VAL:HG13	1:A:250:GLY:O	2.14	0.48
1:B:153:ARG:HD2	1:B:164:GLU:OE2	2.14	0.48
1:C:144:TRP:CE2	1:C:196:PRO:HG3	2.49	0.48
1:A:192:TRP:CD2	1:A:241:ILE:HG12	2.49	0.47
1:B:115:GLY:O	1:B:136:GLY:N	2.29	0.47
1:A:28:GLU:HB2	1:A:41:PHE:HB3	1.95	0.47
1:C:76:TRP:HE3	1:C:81:TRP:CE2	2.32	0.47
1:A:108:ARG:NH1	1:A:121:CYS:HB3	2.30	0.47
1:C:50:VAL:HG23	1:C:63:VAL:HG22	1.95	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:192:TRP:CE3	1:B:241:ILE:HD13	2.50	0.47
1:A:138:ASN:ND2	1:A:187:THR:H	2.11	0.47
1:C:9:SER:HB3	1:C:16:ILE:HG12	1.97	0.46
1:A:15:SER:HA	1:A:31:TRP:O	2.15	0.46
1:A:106:SER:HA	1:A:122:TRP:O	2.15	0.46
1:C:153:ARG:NH1	1:C:167:TRP:O	2.48	0.46
1:A:144:TRP:CE2	1:A:151:HIS:HB2	2.50	0.46
1:A:255:GLU:HG2	1:A:266:GLY:HA3	1.98	0.46
1:B:223:PHE:HB2	1:B:263:TRP:CE2	2.51	0.45
1:B:101:TRP:CE2	1:B:106:SER:HB2	2.51	0.45
1:C:4:GLN:HG2	1:C:229:ASN:HD21	1.81	0.45
1:B:179:SER:HB2	1:B:211:ARG:NE	2.32	0.45
1:B:201:TYR:HE2	1:B:232:VAL:HG21	1.82	0.45
1:A:101:TRP:CE2	1:A:106:SER:HB2	2.52	0.44
1:A:247:ALA:O	1:A:253:THR:HA	2.16	0.44
1:C:144:TRP:O	1:C:150:ILE:HD12	2.17	0.44
1:C:61:ILE:HB	1:C:76:TRP:HB3	1.99	0.44
1:A:121:CYS:HB2	1:A:128:TYR:CZ	2.53	0.44
1:C:144:TRP:CE2	1:C:151:HIS:HB2	2.52	0.44
1:B:235:TRP:CZ3	1:B:244:ARG:HG2	2.53	0.43
1:B:253:THR:H	1:B:270:ALA:HB2	1.83	0.43
1:C:155:TYR:CE1	1:C:164:GLU:HG3	2.53	0.43
1:B:181:THR:HB	1:B:204:ASN:HD21	1.84	0.43
1:B:202:THR:O	1:B:208:ILE:HA	2.18	0.43
1:A:3:VAL:HG13	1:A:251:THR:HA	2.01	0.43
1:B:242:HIS:ND1	1:B:259:ASP:HA	2.34	0.43
1:B:244:ARG:NH1	1:B:257:CYS:HB3	2.25	0.43
1:C:146:VAL:HB	1:C:151:HIS:CD2	2.53	0.43
1:A:14:PRO:HG3	1:A:235:TRP:CZ2	2.53	0.42
1:B:235:TRP:HH2	1:B:244:ARG:NE	2.17	0.42
1:C:155:TYR:CE1	1:C:200:VAL:HG21	2.54	0.42
1:B:192:TRP:CE2	1:B:197:SER:HB2	2.54	0.42
1:A:47:ASN:ND2	1:A:95:GLN:HG2	2.34	0.42
1:C:53:TRP:CZ2	1:C:105:PRO:HG3	2.54	0.42
1:C:138:ASN:HD21	1:C:186:GLN:HG2	1.80	0.42
1:C:4:GLN:HA	1:C:229:ASN:ND2	2.34	0.42
1:A:208:ILE:HD13	1:A:230:VAL:HB	2.00	0.42
1:A:255:GLU:O	1:A:266:GLY:N	2.52	0.42
1:B:192:TRP:CE2	1:B:241:ILE:HG21	2.55	0.42
1:C:243:ILE:HB	1:C:258:TRP:HB3	2.00	0.42
1:A:185:ASP:N	1:A:185:ASP:OD1	2.52	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:234:SER:HB2	1:A:241:ILE:HD11	2.01	0.42
1:A:235:TRP:O	1:A:241:ILE:HD12	2.19	0.42
1:B:10:TRP:HB2	1:B:54:LEU:HD11	2.01	0.41
1:B:207:LYS:HA	1:B:226:PRO:HA	2.03	0.41
1:C:29:ARG:NH2	1:C:269:THR:OG1	2.54	0.41
1:A:53:TRP:CZ2	1:A:105:PRO:HG3	2.56	0.41
1:B:100:SER:CB	1:B:107:ILE:HG12	2.50	0.41
1:B:107:ILE:HB	1:B:122:TRP:HB3	2.03	0.41
1:B:235:TRP:HZ3	1:B:244:ARG:HG2	1.85	0.41
1:C:156:ALA:O	1:C:162:THR:HA	2.20	0.41
1:C:164:GLU:HB2	1:C:177:TYR:CB	2.51	0.41
1:C:227:GLY:HA2	1:C:249:THR:OG1	2.21	0.41
1:A:201:TYR:CE2	1:A:232:VAL:HG21	2.55	0.41
1:C:210:GLU:HB2	1:C:223:PHE:HB3	2.03	0.40
1:B:106:SER:HA	1:B:122:TRP:O	2.21	0.40
1:C:108:ARG:HD2	1:C:119:GLU:OE2	2.21	0.40
1:C:3:VAL:HG13	1:C:250:GLY:O	2.22	0.40
1:A:138:ASN:HD21	1:A:186:GLN:HG2	1.86	0.40
1:A:212:CYS:HB2	1:A:219:TYR:CE1	2.57	0.40
1:B:141:VAL:HG23	1:B:154:VAL:HG22	2.04	0.40
1:C:144:TRP:CZ3	1:C:198:ILE:HD11	2.57	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	258/272 (95%)	241 (93%)	17 (7%)	0	100	100
1	B	258/272 (95%)	240 (93%)	15 (6%)	3 (1%)	13	48
1	C	258/272 (95%)	237 (92%)	21 (8%)	0	100	100
All	All	774/816 (95%)	718 (93%)	53 (7%)	3 (0%)	34	71

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	114	ASN
1	B	185	ASP
1	B	238	GLY

### 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	208/215 (97%)	208 (100%)	0	100	100
1	B	208/215 (97%)	208 (100%)	0	100	100
1	C	208/215 (97%)	208 (100%)	0	100	100
All	All	624/645 (97%)	624 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

### 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 monosaccharides in this entry.

## 5.6 Ligand geometry

There are no ligands in this entry.

## 5.7 Other polymers

There are no such residues in this entry.

## 5.8 Polymer linkage issues

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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	264/272 (97%)	0.01	5 (1%) 66 58	44, 77, 114, 153	0
1	B	264/272 (97%)	0.05	6 (2%) 60 51	44, 83, 113, 142	0
1	C	264/272 (97%)	0.01	1 (0%) 92 87	50, 82, 117, 144	0
All	All	792/816 (97%)	0.02	12 (1%) 73 63	44, 81, 116, 153	0

All (12) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	269	THR	4.3
1	A	159	GLY	3.7
1	B	268	TYR	3.4
1	A	170	ASN	2.5
1	A	260	GLY	2.5
1	B	130	GLY	2.5
1	C	66	SER	2.4
1	A	261	ASN	2.3
1	B	271	THR	2.1
1	B	131	ALA	2.1
1	B	261	ASN	2.0
1	A	45	GLY	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 monosaccharides 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.