



Full wwPDB X-ray Structure Validation Report ⓘ

Dec 15, 2022 – 03:28 pm GMT

PDB ID : 6S5Y
Title : Structure of tandemly arrayed consecutive Rib domains (Rib2R) from Group B Streptococcal species Streptococcus agalactiae
Authors : Whelan, F.; Griffiths, S.C.; Bateman, A.; Potts, J.R.
Deposited on : 2019-07-02
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 : 4.02b-467
Xtriage (Phenix) : 1.13
EDS : 2.31.3
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) : 2.31.3

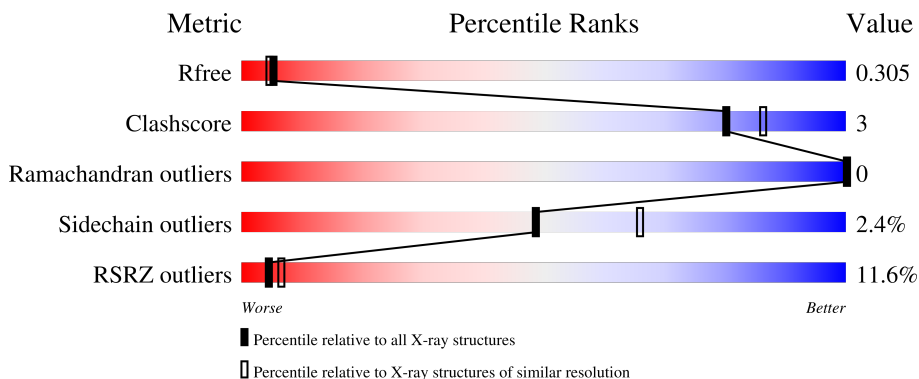
1 Overall quality at a glance

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.



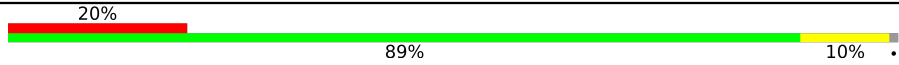
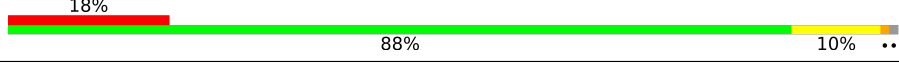
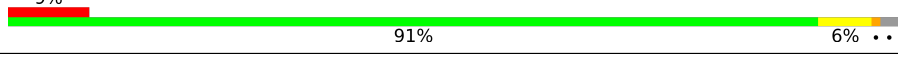
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	5042 (2.30-2.30)
Clashscore	141614	5643 (2.30-2.30)
Ramachandran outliers	138981	5575 (2.30-2.30)
Sidechain outliers	138945	5575 (2.30-2.30)
RSRZ outliers	127900	4938 (2.30-2.30)

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 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	163	 9% 93% 7%
1	B	163	 2% 90% 9% ..
1	C	163	 16% 88% 10% .
1	D	163	 9% 91% 7% .
1	E	163	 9% 90% 9% ..

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Mol	Chain	Length	Quality of chain
1	F	163	 20% 89% 10% .
1	G	163	 18% 88% 10% ..
1	H	163	 9% 91% 6% ..

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 9423 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 Group B streptococcal R4 surface protein.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
			Total	C	N	O			
1	A	163	1185	730	195	260	0	0	0
1	B	161	1173	723	193	257	0	0	0
1	C	161	1174	724	193	257	0	0	0
1	D	161	1174	724	193	257	0	0	0
1	E	161	1174	724	193	257	0	0	0
1	F	161	1174	724	193	257	0	0	0
1	G	161	1174	724	193	257	0	0	0
1	H	160	1163	718	189	256	0	0	0

There are 40 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	225	GLY	-	expression tag	UNP P72362
A	226	PRO	-	expression tag	UNP P72362
A	227	LEU	-	expression tag	UNP P72362
A	228	GLY	-	expression tag	UNP P72362
A	229	SER	-	expression tag	UNP P72362
B	225	GLY	-	expression tag	UNP P72362
B	226	PRO	-	expression tag	UNP P72362
B	227	LEU	-	expression tag	UNP P72362
B	228	GLY	-	expression tag	UNP P72362
B	229	SER	-	expression tag	UNP P72362
C	225	GLY	-	expression tag	UNP P72362
C	226	PRO	-	expression tag	UNP P72362
C	227	LEU	-	expression tag	UNP P72362

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Chain	Residue	Modelled	Actual	Comment	Reference
C	228	GLY	-	expression tag	UNP P72362
C	229	SER	-	expression tag	UNP P72362
D	225	GLY	-	expression tag	UNP P72362
D	226	PRO	-	expression tag	UNP P72362
D	227	LEU	-	expression tag	UNP P72362
D	228	GLY	-	expression tag	UNP P72362
D	229	SER	-	expression tag	UNP P72362
E	225	GLY	-	expression tag	UNP P72362
E	226	PRO	-	expression tag	UNP P72362
E	227	LEU	-	expression tag	UNP P72362
E	228	GLY	-	expression tag	UNP P72362
E	229	SER	-	expression tag	UNP P72362
F	225	GLY	-	expression tag	UNP P72362
F	226	PRO	-	expression tag	UNP P72362
F	227	LEU	-	expression tag	UNP P72362
F	228	GLY	-	expression tag	UNP P72362
F	229	SER	-	expression tag	UNP P72362
G	225	GLY	-	expression tag	UNP P72362
G	226	PRO	-	expression tag	UNP P72362
G	227	LEU	-	expression tag	UNP P72362
G	228	GLY	-	expression tag	UNP P72362
G	229	SER	-	expression tag	UNP P72362
H	225	GLY	-	expression tag	UNP P72362
H	226	PRO	-	expression tag	UNP P72362
H	227	LEU	-	expression tag	UNP P72362
H	228	GLY	-	expression tag	UNP P72362
H	229	SER	-	expression tag	UNP P72362

- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	5	Total O 5 5	0	0
2	B	8	Total O 8 8	0	0
2	C	1	Total O 1 1	0	0
2	D	2	Total O 2 2	0	0
2	F	3	Total O 3 3	0	0
2	G	8	Total O 8 8	0	0

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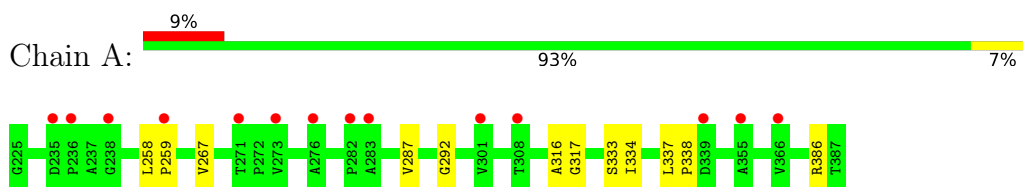
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	H	5	Total	O	0	0
			5	5		

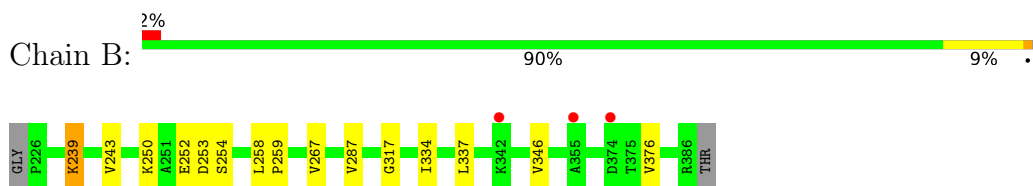
3 Residue-property plots [i](#)

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.

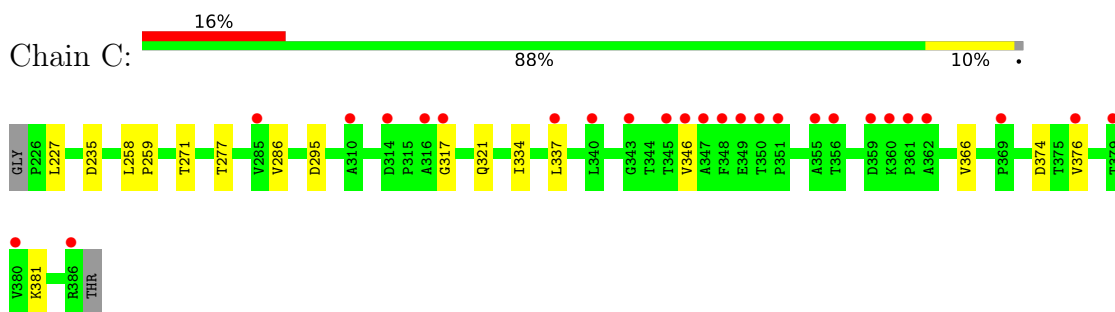
- Molecule 1: Group B streptococcal R4 surface protein



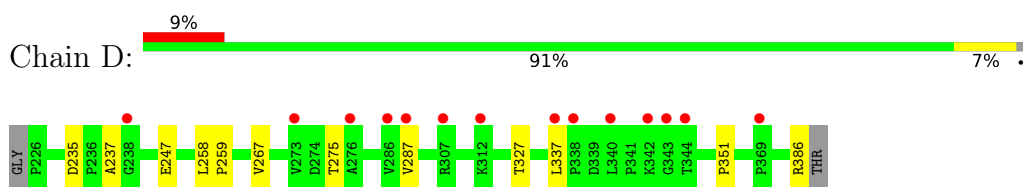
- Molecule 1: Group B streptococcal R4 surface protein



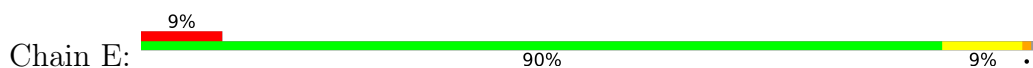
- Molecule 1: Group B streptococcal R4 surface protein

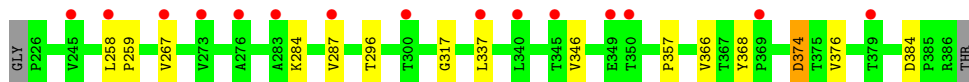


- Molecule 1: Group B streptococcal R4 surface protein

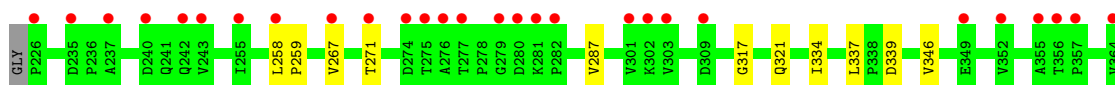
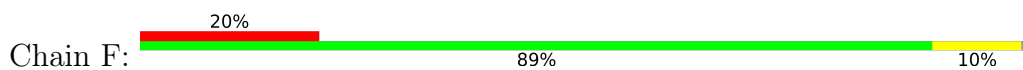


- Molecule 1: Group B streptococcal R4 surface protein

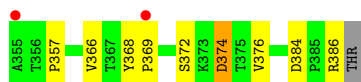
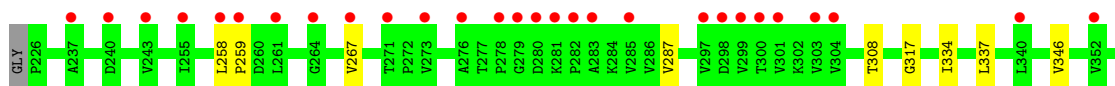
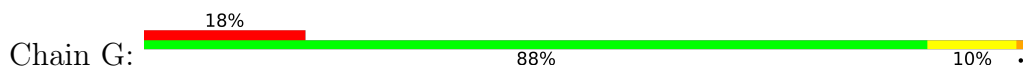




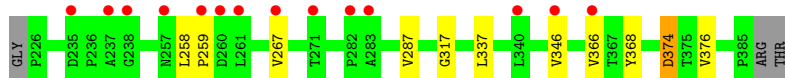
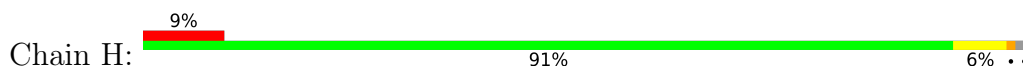
- Molecule 1: Group B streptococcal R4 surface protein



- Molecule 1: Group B streptococcal R4 surface protein



- Molecule 1: Group B streptococcal R4 surface protein



4 Data and refinement statistics i

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	37.83Å 380.85Å 37.85Å 90.00° 91.60° 90.00°	Depositor
Resolution (Å)	38.09 – 2.30 47.61 – 2.30	Depositor EDS
% Data completeness (in resolution range)	99.8 (38.09-2.30) 99.8 (47.61-2.30)	Depositor EDS
R_{merge}	0.15	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.75 (at 2.29Å)	Xtrriage
Refinement program	BUSTER 2.10.3	Depositor
R, R_{free}	0.234 , 0.274 0.262 , 0.305	Depositor DCC
R_{free} test set	2349 reflections (4.99%)	wwPDB-VP
Wilson B-factor (Å ²)	48.3	Xtrriage
Anisotropy	0.689	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 70.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	0.003 for l,k,-h 0.096 for h,-k,-l 0.056 for l,-k,h	Xtrriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	9423	wwPDB-VP
Average B, all atoms (Å ²)	91.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 9.87% 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	0.43	0/1207	0.61	0/1658
1	B	0.55	0/1195	0.68	0/1641
1	C	0.45	0/1196	0.61	0/1642
1	D	0.43	0/1196	0.60	0/1642
1	E	0.40	0/1196	0.59	0/1642
1	F	0.40	0/1196	0.58	0/1642
1	G	0.42	0/1196	0.61	0/1642
1	H	0.41	0/1185	0.61	0/1628
All	All	0.44	0/9567	0.61	0/13137

There are no bond length outliers.

There are no bond angle outliers.

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	1185	0	1162	7	0
1	B	1173	0	1148	7	0
1	C	1174	0	1153	7	0
1	D	1174	0	1153	7	0
1	E	1174	0	1153	7	0
1	F	1174	0	1153	7	0
1	G	1174	0	1153	10	0
1	H	1163	0	1140	6	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	A	5	0	0	0	0
2	B	8	0	0	0	0
2	C	1	0	0	0	0
2	D	2	0	0	0	0
2	F	3	0	0	0	0
2	G	8	0	0	0	0
2	H	5	0	0	0	0
All	All	9423	0	9215	51	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 (51) 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:E:357:PRO:HG3	1:E:384:ASP:HB2	1.68	0.74
1:B:239:LYS:O	1:B:254:SER:HA	1.91	0.69
1:C:227:LEU:HB3	1:D:327:THR:HB	1.75	0.69
1:F:321:GLN:HG2	1:F:381:LYS:HB3	1.77	0.66
1:E:284:LYS:HZ3	1:E:296:THR:HG21	1.65	0.61
1:A:292:GLY:HA2	1:D:351:PRO:HG2	1.84	0.59
1:G:317:GLY:HA3	1:G:376:VAL:HG11	1.85	0.57
1:C:321:GLN:HG3	1:C:381:LYS:HB3	1.88	0.55
1:E:317:GLY:HA3	1:E:376:VAL:HG11	1.86	0.55
1:H:317:GLY:HA3	1:H:376:VAL:HG11	1.90	0.53
1:C:317:GLY:HA3	1:C:376:VAL:HG11	1.90	0.53
1:B:317:GLY:HA3	1:B:376:VAL:HG11	1.91	0.52
1:F:317:GLY:HA3	1:F:376:VAL:HG11	1.92	0.52
1:B:252:GLU:HG3	1:B:267:VAL:CG2	2.40	0.51
1:C:277:THR:HG21	1:F:271:THR:HG23	1.93	0.50
1:H:267:VAL:HG12	1:H:287:VAL:HG22	1.95	0.49
1:D:247:GLU:O	1:D:275:THR:HG21	2.13	0.48
1:H:346:VAL:HG12	1:H:366:VAL:HG22	1.96	0.47
1:H:368:TYR:HE2	1:H:374:ASP:HB2	1.79	0.47
1:B:267:VAL:HG12	1:B:287:VAL:HG22	1.98	0.46
1:G:308:THR:OG1	1:G:372:SER:HB3	2.16	0.46
1:F:346:VAL:HG12	1:F:366:VAL:HG22	1.99	0.45
1:G:384:ASP:C	1:G:386:ARG:H	2.19	0.45
1:G:267:VAL:HG12	1:G:287:VAL:HG22	1.99	0.44
1:E:267:VAL:HG12	1:E:287:VAL:HG22	2.00	0.44
1:F:267:VAL:HG12	1:F:287:VAL:HG22	2.00	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:317:GLY:O	1:D:235:ASP:HB3	2.18	0.43
1:B:334:ILE:HG13	1:B:346:VAL:HG11	2.00	0.43
1:E:368:TYR:HE2	1:E:374:ASP:HB2	1.83	0.43
1:F:334:ILE:HG13	1:F:346:VAL:HG11	2.00	0.43
1:G:258:LEU:N	1:G:259:PRO:HD2	2.34	0.43
1:G:368:TYR:HE2	1:G:374:ASP:HB2	1.84	0.43
1:E:346:VAL:HG12	1:E:366:VAL:HG22	2.01	0.43
1:G:334:ILE:HG13	1:G:346:VAL:HG11	2.01	0.43
1:F:258:LEU:N	1:F:259:PRO:HD2	2.34	0.42
1:G:346:VAL:HG12	1:G:366:VAL:HG22	2.01	0.42
1:A:267:VAL:HG12	1:A:287:VAL:HG22	1.99	0.42
1:C:346:VAL:HG12	1:C:366:VAL:HG22	2.01	0.42
1:A:258:LEU:N	1:A:259:PRO:HD2	2.35	0.42
1:B:252:GLU:HG3	1:B:267:VAL:HG21	2.02	0.41
1:D:267:VAL:HG12	1:D:287:VAL:HG22	2.02	0.41
1:H:258:LEU:N	1:H:259:PRO:HD2	2.35	0.41
1:A:338:PRO:HG2	1:G:357:PRO:HB2	2.02	0.41
1:C:258:LEU:N	1:C:259:PRO:HD2	2.36	0.41
1:B:258:LEU:N	1:B:259:PRO:HD2	2.35	0.41
1:E:258:LEU:N	1:E:259:PRO:HD2	2.36	0.41
1:A:317:GLY:HA2	1:A:334:ILE:HD13	2.03	0.41
1:D:258:LEU:N	1:D:259:PRO:HD2	2.36	0.40
1:A:316:ALA:HB2	1:D:237:ALA:HB2	2.04	0.40
1:G:369:PRO:O	1:H:259:PRO:HD3	2.21	0.40
1:C:334:ILE:HG13	1:C:346:VAL:HG11	2.04	0.40

There are no symmetry-related clashes.

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	161/163 (99%)	158 (98%)	3 (2%)	0	100 100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	159/163 (98%)	156 (98%)	3 (2%)	0	100	100
1	C	159/163 (98%)	155 (98%)	4 (2%)	0	100	100
1	D	159/163 (98%)	157 (99%)	2 (1%)	0	100	100
1	E	159/163 (98%)	157 (99%)	2 (1%)	0	100	100
1	F	159/163 (98%)	157 (99%)	2 (1%)	0	100	100
1	G	159/163 (98%)	156 (98%)	3 (2%)	0	100	100
1	H	158/163 (97%)	157 (99%)	1 (1%)	0	100	100
All	All	1273/1304 (98%)	1253 (98%)	20 (2%)	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	137/137 (100%)	134 (98%)	3 (2%)	52	69
1	B	135/137 (98%)	130 (96%)	5 (4%)	34	48
1	C	136/137 (99%)	130 (96%)	6 (4%)	28	39
1	D	136/137 (99%)	134 (98%)	2 (2%)	65	79
1	E	136/137 (99%)	134 (98%)	2 (2%)	65	79
1	F	136/137 (99%)	132 (97%)	4 (3%)	42	58
1	G	136/137 (99%)	134 (98%)	2 (2%)	65	79
1	H	135/137 (98%)	133 (98%)	2 (2%)	65	79
All	All	1087/1096 (99%)	1061 (98%)	26 (2%)	49	66

All (26) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	333	SER
1	A	337	LEU

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Mol	Chain	Res	Type
1	A	386	ARG
1	B	239	LYS
1	B	243	VAL
1	B	250	LYS
1	B	253	ASP
1	B	337	LEU
1	C	235	ASP
1	C	271	THR
1	C	286	VAL
1	C	295	ASP
1	C	337	LEU
1	C	374	ASP
1	D	337	LEU
1	D	386	ARG
1	E	337	LEU
1	E	374	ASP
1	F	337	LEU
1	F	339	ASP
1	F	374	ASP
1	F	386	ARG
1	G	337	LEU
1	G	374	ASP
1	H	337	LEU
1	H	374	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	242	GLN
1	C	242	GLN
1	C	321	GLN
1	D	313	ASN
1	E	242	GLN
1	H	242	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 monosaccharides 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	163/163 (100%)	0.76	14 (8%) 10 14	61, 86, 112, 118	0
1	B	161/163 (98%)	0.68	3 (1%) 66 73	33, 83, 112, 138	0
1	C	161/163 (98%)	0.94	26 (16%) 1 2	46, 93, 134, 164	0
1	D	161/163 (98%)	0.55	14 (8%) 10 14	59, 81, 99, 117	0
1	E	161/163 (98%)	0.87	15 (9%) 8 11	73, 91, 114, 124	0
1	F	161/163 (98%)	1.17	33 (20%) 1 1	60, 99, 130, 165	0
1	G	161/163 (98%)	1.23	30 (18%) 1 1	76, 94, 126, 139	0
1	H	160/163 (98%)	0.73	14 (8%) 10 13	71, 90, 109, 118	0
All	All	1289/1304 (98%)	0.87	149 (11%) 4 6	33, 89, 121, 165	0

All (149) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	G	282	PRO	9.3
1	F	355	ALA	6.7
1	F	356	THR	6.3
1	G	258	LEU	6.0
1	G	283	ALA	5.6
1	C	349	GLU	5.5
1	F	349	GLU	5.4
1	F	279	GLY	5.0
1	A	238	GLY	4.9
1	G	280	ASP	4.8
1	C	350	THR	4.8
1	C	386	ARG	4.8
1	G	279	GLY	4.8
1	F	366	VAL	4.5
1	G	281	LYS	4.4
1	H	282	PRO	4.3

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Mol	Chain	Res	Type	RSRZ
1	C	361	PRO	4.3
1	G	259	PRO	4.3
1	C	379	THR	4.2
1	G	271	THR	4.2
1	C	369	PRO	4.1
1	G	276	ALA	4.0
1	H	346	VAL	4.0
1	A	273	VAL	4.0
1	D	238	GLY	4.0
1	A	366	VAL	4.0
1	F	281	LYS	4.0
1	G	237	ALA	3.8
1	F	258	LEU	3.8
1	C	317	GLY	3.8
1	F	369	PRO	3.8
1	C	360	LYS	3.7
1	H	237	ALA	3.7
1	A	235	ASP	3.6
1	F	226	PRO	3.6
1	G	273	VAL	3.6
1	E	283	ALA	3.6
1	D	369	PRO	3.6
1	G	240	ASP	3.6
1	G	255	ILE	3.6
1	C	362	ALA	3.4
1	C	380	VAL	3.4
1	C	343	GLY	3.4
1	E	258	LEU	3.4
1	F	357	PRO	3.3
1	H	235	ASP	3.3
1	A	283	ALA	3.3
1	A	259	PRO	3.3
1	C	314	ASP	3.2
1	G	303	VAL	3.2
1	F	240	ASP	3.2
1	G	352	VAL	3.1
1	E	340	LEU	3.1
1	H	283	ALA	3.1
1	E	245	VAL	3.1
1	B	374	ASP	3.0
1	A	339	ASP	3.0
1	A	282	PRO	3.0

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Mol	Chain	Res	Type	RSRZ
1	C	351	PRO	3.0
1	D	343	GLY	3.0
1	F	282	PRO	3.0
1	F	309	ASP	3.0
1	G	355	ALA	3.0
1	C	345	THR	2.9
1	C	346	VAL	2.9
1	F	380	VAL	2.9
1	E	369	PRO	2.9
1	G	300	THR	2.9
1	C	310	ALA	2.9
1	H	259	PRO	2.9
1	C	316	ALA	2.8
1	D	287	VAL	2.8
1	F	242	GLN	2.8
1	D	342	LYS	2.8
1	F	267	VAL	2.8
1	E	300	THR	2.8
1	G	369	PRO	2.7
1	C	376	VAL	2.7
1	G	267	VAL	2.7
1	F	302	LYS	2.7
1	H	340	LEU	2.7
1	F	276	ALA	2.7
1	C	347	ALA	2.7
1	D	338	PRO	2.6
1	E	276	ALA	2.6
1	C	340	LEU	2.6
1	E	350	THR	2.6
1	H	271	THR	2.6
1	F	385	PRO	2.6
1	D	340	LEU	2.6
1	B	355	ALA	2.6
1	E	349	GLU	2.6
1	E	379	THR	2.6
1	G	298	ASP	2.6
1	F	301	VAL	2.6
1	D	344	THR	2.6
1	G	301	VAL	2.6
1	D	337	LEU	2.5
1	G	304	VAL	2.5
1	F	364	VAL	2.5

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Mol	Chain	Res	Type	RSRZ
1	F	271	THR	2.5
1	G	340	LEU	2.5
1	C	355	ALA	2.5
1	E	267	VAL	2.5
1	F	370	ASP	2.5
1	C	285	VAL	2.4
1	F	255	ILE	2.4
1	G	297	VAL	2.4
1	A	276	ALA	2.4
1	A	308	THR	2.4
1	D	286	VAL	2.4
1	G	278	PRO	2.4
1	G	299	VAL	2.4
1	F	352	VAL	2.4
1	H	260	ASP	2.3
1	G	264	GLY	2.3
1	E	345	THR	2.3
1	F	277	THR	2.3
1	F	303	VAL	2.3
1	D	312	LYS	2.3
1	A	236	PRO	2.3
1	E	273	VAL	2.3
1	D	307	ARG	2.3
1	F	274	ASP	2.3
1	C	337	LEU	2.2
1	H	261	LEU	2.2
1	F	280	ASP	2.2
1	G	243	VAL	2.2
1	C	348	PHE	2.2
1	D	276	ALA	2.2
1	C	359	ASP	2.2
1	F	243	VAL	2.2
1	A	271	THR	2.2
1	G	261	LEU	2.2
1	H	238	GLY	2.2
1	B	342	LYS	2.1
1	H	257	ASN	2.1
1	E	287	VAL	2.1
1	A	301	VAL	2.1
1	D	273	VAL	2.1
1	G	285	VAL	2.1
1	A	355	ALA	2.1

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Mol	Chain	Res	Type	RSRZ
1	H	267	VAL	2.1
1	F	237	ALA	2.1
1	F	275	THR	2.1
1	H	366	VAL	2.1
1	F	235	ASP	2.0
1	E	337	LEU	2.0
1	C	356	THR	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.