



Full wwPDB X-ray Structure Validation Report ⓘ

May 29, 2020 – 03:43 am BST

PDB ID : 3T1R
Title : MglB with tetrameric arrangement
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Deposited on : 2011-07-22
Resolution : 2.00 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

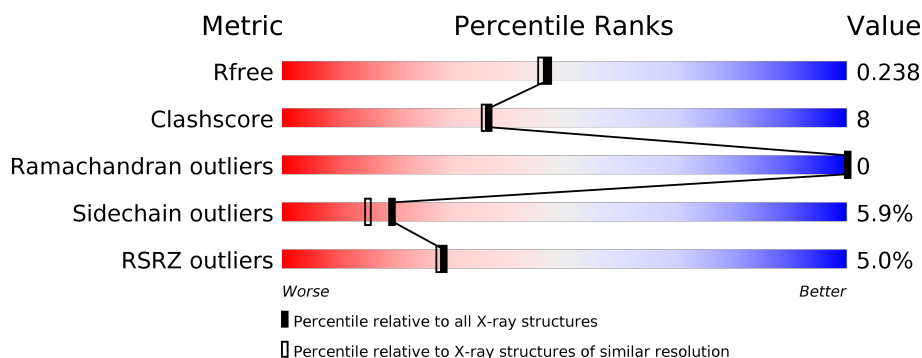
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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	8085 (2.00-2.00)
Clashscore	141614	9178 (2.00-2.00)
Ramachandran outliers	138981	9054 (2.00-2.00)
Sidechain outliers	138945	9053 (2.00-2.00)
RSRZ outliers	127900	7900 (2.00-2.00)

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 $\leq 5\%$. 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	165	<div> <div>3%</div> <div> <div></div> <div>64%</div> <div>15%</div> <div>•</div> <div>19%</div> </div> </div>
1	B	165	<div> <div></div> <div> <div>66%</div> <div>13%</div> <div>•</div> <div>19%</div> </div> </div>
1	C	165	<div> <div>7%</div> <div> <div></div> <div>70%</div> <div>9%</div> <div>••</div> <div>18%</div> </div> </div>
1	D	165	<div> <div>7%</div> <div> <div></div> <div>69%</div> <div>9%</div> <div>•</div> <div>19%</div> </div> </div>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 4485 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 Gliding protein MglB.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	133	Total	C	N	O	S	0	8	0
			1066	684	181	200	1			
1	B	134	Total	C	N	O	S	0	4	0
			1054	674	183	196	1			
1	C	136	Total	C	N	O	S	0	3	0
			1061	681	183	196	1			
1	D	134	Total	C	N	O	S	0	6	0
			1065	682	187	195	1			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	GLY	-	EXPRESSION TAG	UNP Q5SJ83
A	0	SER	-	EXPRESSION TAG	UNP Q5SJ83
A	65	SER	GLY	VARIANT	UNP Q5SJ83
B	-1	GLY	-	EXPRESSION TAG	UNP Q5SJ83
B	0	SER	-	EXPRESSION TAG	UNP Q5SJ83
B	65	SER	GLY	ENGINEERED MUTATION	UNP Q5SJ83
C	-1	GLY	-	EXPRESSION TAG	UNP Q5SJ83
C	0	SER	-	EXPRESSION TAG	UNP Q5SJ83
C	65	SER	GLY	VARIANT	UNP Q5SJ83
D	-1	GLY	-	EXPRESSION TAG	UNP Q5SJ83
D	0	SER	-	EXPRESSION TAG	UNP Q5SJ83
D	65	SER	GLY	VARIANT	UNP Q5SJ83

- Molecule 2 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	1	Total	Mg	0	0
			1	1		
2	A	1	Total	Mg	0	0
			1	1		

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	D	2	Total 2	Mg 2	0	0
2	C	1	Total 1	Mg 1	0	0

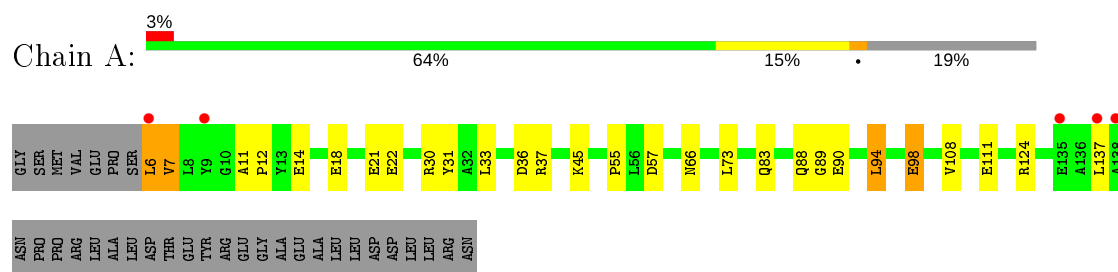
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	83	Total 83	O 83	0	0
3	B	67	Total 67	O 67	0	0
3	C	43	Total 43	O 43	0	0
3	D	41	Total 41	O 41	0	0

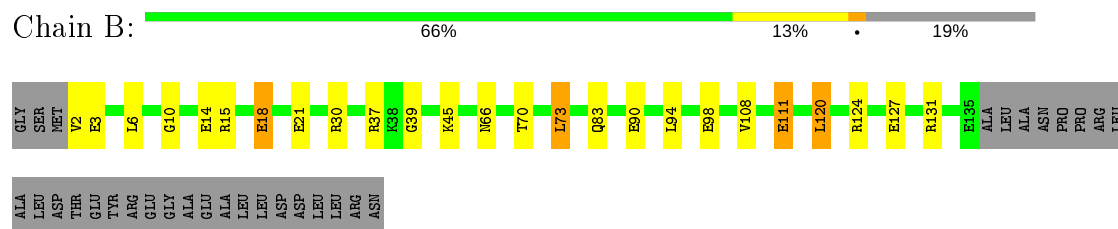
3 Residue-property plots [i](#)

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.

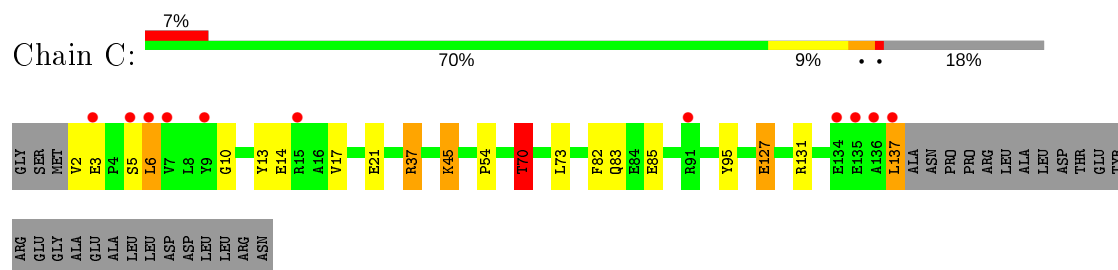
• Molecule 1: Gliding protein MglB



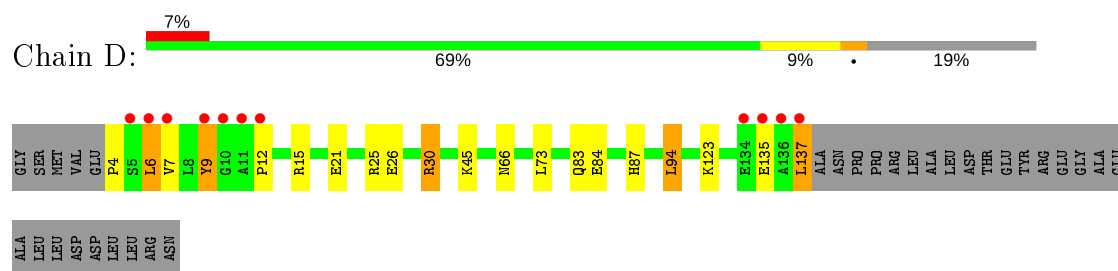
• Molecule 1: Gliding protein MglB



• Molecule 1: Gliding protein MglB



• Molecule 1: Gliding protein MglB



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	99.20Å 118.00Å 111.70Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	45.33 – 2.00 45.33 – 2.00	Depositor EDS
% Data completeness (in resolution range)	100.0 (45.33-2.00) 98.2 (45.33-2.00)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.22 (at 2.00Å)	Xtriage
Refinement program	REFMAC 5.5.0110	Depositor
R, R_{free}	0.182 , 0.236 0.183 , 0.238	Depositor DCC
R_{free} test set	2189 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	27.1	Xtriage
Anisotropy	0.229	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.40 , 52.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	4485	wwPDB-VP
Average B, all atoms (Å ²)	29.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.62% 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](#)

Bond lengths and bond angles in the following residue types are not validated in this section: MG

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	1.15	2/1105 (0.2%)	1.10	7/1499 (0.5%)
1	B	1.16	2/1085 (0.2%)	1.22	8/1472 (0.5%)
1	C	1.12	2/1089 (0.2%)	1.03	4/1478 (0.3%)
1	D	1.05	0/1102	1.03	3/1493 (0.2%)
All	All	1.12	6/4381 (0.1%)	1.10	22/5942 (0.4%)

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	18	GLU	CB-CG	6.84	1.65	1.52
1	A	98	GLU	CG-CD	6.06	1.61	1.51
1	C	95	TYR	CD1-CE1	5.82	1.48	1.39
1	A	31	TYR	CD1-CE1	5.53	1.47	1.39
1	C	70	THR	CB-CG2	-5.31	1.34	1.52
1	B	108	VAL	CB-CG1	5.09	1.63	1.52

All (22) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	37	ARG	NE-CZ-NH2	-16.86	111.87	120.30
1	B	37	ARG	NE-CZ-NH1	15.84	128.22	120.30
1	D	30	ARG	NE-CZ-NH2	-13.56	113.52	120.30
1	A	37	ARG	NE-CZ-NH2	-10.95	114.83	120.30
1	A	37	ARG	NE-CZ-NH1	10.67	125.64	120.30
1	B	30	ARG	NE-CZ-NH2	-10.29	115.15	120.30
1	C	37	ARG	NE-CZ-NH1	7.50	124.05	120.30
1	D	30	ARG	NE-CZ-NH1	7.43	124.02	120.30
1	B	94	LEU	CB-CG-CD2	-6.92	99.23	111.00
1	C	37	ARG	NE-CZ-NH2	-6.34	117.13	120.30
1	A	30	ARG	NE-CZ-NH2	-6.30	117.15	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	94	LEU	CB-CG-CD2	-5.83	101.08	111.00
1	C	6	LEU	CA-CB-CG	5.76	128.54	115.30
1	B	37	ARG	CD-NE-CZ	5.64	131.50	123.60
1	B	131	ARG	NE-CZ-NH1	-5.57	117.52	120.30
1	C	70	THR	N-CA-CB	-5.56	99.74	110.30
1	B	37	ARG	CG-CD-NE	-5.52	100.22	111.80
1	A	36	ASP	CB-CG-OD2	5.20	122.98	118.30
1	A	30	ARG	NE-CZ-NH1	5.18	122.89	120.30
1	B	73	LEU	CA-CB-CG	5.14	127.12	115.30
1	A	37	ARG	CG-CD-NE	-5.04	101.22	111.80
1	A	33	LEU	CB-CG-CD2	-5.02	102.47	111.00

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	1066	0	1095	23	1
1	B	1054	0	1080	17	1
1	C	1061	0	1093	22	0
1	D	1065	0	1103	18	1
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	2	0	0	0	0
3	A	83	0	0	9	0
3	B	67	0	0	4	1
3	C	43	0	0	3	0
3	D	41	0	0	5	0
All	All	4485	0	4371	68	2

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 8.

All (68) 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:127:GLU:OE2	1:C:131[A]:ARG:NE	2.04	0.91
1:D:83[A]:GLN:HG3	1:D:84[A]:GLU:CD	1.91	0.90
1:A:66[B]:ASN:HD21	1:B:70:THR:HG22	1.37	0.90
1:D:21:GLU:HG2	1:D:45:LYS:HE2	1.57	0.83
1:D:83[A]:GLN:HG3	1:D:84[A]:GLU:OE2	1.77	0.82
1:D:123:LYS:HD2	3:D:288:HOH:O	1.77	0.82
1:B:66[A]:ASN:OD1	3:B:168:HOH:O	1.98	0.79
1:B:111:GLU:HG3	3:B:205:HOH:O	1.82	0.77
1:C:73[A]:LEU:HD12	1:D:66[A]:ASN:ND2	2.01	0.74
1:D:66[B]:ASN:ND2	3:D:303:HOH:O	2.21	0.73
1:D:66[B]:ASN:OD1	3:D:167:HOH:O	2.07	0.72
1:B:124:ARG:NH1	3:B:180:HOH:O	1.84	0.71
1:A:111[A]:GLU:HG2	3:A:211:HOH:O	1.90	0.70
1:D:25[B]:ARG:NH1	3:D:175:HOH:O	2.09	0.70
1:C:127:GLU:CD	1:C:131[A]:ARG:HE	1.94	0.70
1:A:66[B]:ASN:HD21	1:B:70:THR:CG2	2.05	0.68
1:C:70:THR:HG21	1:C:85:GLU:OE2	1.94	0.67
1:A:57[B]:ASP:OD2	3:A:239:HOH:O	2.11	0.67
1:C:73[A]:LEU:HD12	1:D:66[A]:ASN:HD22	1.60	0.67
1:C:131[B]:ARG:HH11	1:C:131[B]:ARG:HG3	1.58	0.67
1:A:14[A]:GLU:OE2	3:A:225:HOH:O	2.12	0.66
1:C:127:GLU:OE2	1:C:131[B]:ARG:NH1	2.30	0.65
1:A:90:GLU:HG2	1:B:83[A]:GLN:OE1	1.96	0.65
1:C:73[B]:LEU:HD21	1:D:94:LEU:CD2	2.28	0.64
1:C:21:GLU:HG2	1:C:45:LYS:HE2	1.79	0.64
1:A:6:LEU:HD13	1:A:137:LEU:HD13	1.78	0.64
1:A:90:GLU:O	3:A:230:HOH:O	2.15	0.64
1:A:111[B]:GLU:CD	3:A:214:HOH:O	2.36	0.64
1:B:39:GLY:O	3:B:235:HOH:O	2.16	0.63
1:D:83[A]:GLN:CG	1:D:84[A]:GLU:OE2	2.47	0.62
1:A:21:GLU:HG2	1:A:45:LYS:HE2	1.81	0.62
1:D:6:LEU:HD12	1:D:137:LEU:HB3	1.80	0.62
1:A:18:GLU:O	1:A:22[B]:GLU:HG2	1.99	0.62
1:C:73[A]:LEU:CD1	1:D:66[A]:ASN:HD22	2.14	0.61
1:A:66[B]:ASN:ND2	1:B:70:THR:HG22	2.11	0.59
1:A:6:LEU:N	3:A:243:HOH:O	2.35	0.58
1:C:10:GLY:O	1:C:14[A]:GLU:HG3	2.03	0.58
1:A:88[A]:GLN:HG3	1:A:89:GLY:O	2.03	0.57
1:A:11:ALA:HB3	1:A:12:PRO:HD3	1.86	0.57
1:C:131[B]:ARG:NH1	1:C:131[B]:ARG:HG3	2.20	0.56
1:A:83[B]:GLN:NE2	1:B:90:GLU:OE1	2.38	0.55
1:D:9:TYR:O	1:D:12:PRO:HD2	2.09	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:94[B]:LEU:CD2	1:A:108:VAL:HG22	2.40	0.52
1:C:73[B]:LEU:HD21	1:D:94:LEU:HD21	1.91	0.51
1:C:13:TYR:O	1:C:17:VAL:HG23	2.11	0.51
1:B:83[B]:GLN:OE1	1:B:98:GLU:OE1	2.30	0.49
1:D:21:GLU:CG	1:D:45:LYS:HE2	2.38	0.48
1:A:14[A]:GLU:CG	3:A:225:HOH:O	2.61	0.48
1:A:111[B]:GLU:HG2	3:A:214:HOH:O	2.14	0.48
1:C:6:LEU:HD12	1:C:137:LEU:HD12	1.94	0.48
1:A:83[A]:GLN:NE2	1:A:98:GLU:OE1	2.46	0.47
1:B:120:LEU:C	1:B:120:LEU:HD12	2.35	0.47
1:B:10:GLY:O	1:B:14[B]:GLU:HG2	2.15	0.47
1:A:111[B]:GLU:CG	3:A:214:HOH:O	2.64	0.45
1:C:45:LYS:HD2	1:C:45:LYS:HA	1.78	0.44
1:C:21:GLU:CG	1:C:45:LYS:HE2	2.47	0.44
1:A:90:GLU:HG2	1:B:83[A]:GLN:CD	2.39	0.43
1:B:21:GLU:OE1	1:B:45:LYS:HE3	2.19	0.43
1:B:45:LYS:HA	1:B:45:LYS:HD2	1.74	0.43
1:C:83:GLN:HG3	3:C:284:HOH:O	2.19	0.42
1:C:70:THR:HG23	1:C:82:PHE:HD2	1.84	0.41
1:B:2:VAL:HG11	1:C:54:PRO:HB3	2.03	0.41
1:C:37:ARG:NH2	3:C:186:HOH:O	2.45	0.41
1:C:70:THR:CG2	3:C:171:HOH:O	2.69	0.41
1:B:15[A]:ARG:HH21	1:B:18:GLU:CD	2.24	0.41
1:D:4:PRO:HB2	1:D:7:VAL:HG22	2.02	0.41
1:D:87:HIS:HE1	3:D:179:HOH:O	2.04	0.40
1:A:6:LEU:O	1:A:7:VAL:HG13	2.21	0.40

All (2) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:127:GLU:OE1	1:D:30:ARG:NH2[8_545]	2.02	0.18
1:A:124:ARG:NH2	3:B:184:HOH:O[4_545]	2.17	0.03

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	139/165 (84%)	137 (99%)	2 (1%)	0	100	100
1	B	136/165 (82%)	135 (99%)	1 (1%)	0	100	100
1	C	137/165 (83%)	135 (98%)	2 (2%)	0	100	100
1	D	138/165 (84%)	132 (96%)	6 (4%)	0	100	100
All	All	550/660 (83%)	539 (98%)	11 (2%)	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	107/126 (85%)	101 (94%)	6 (6%)	21	17
1	B	106/126 (84%)	101 (95%)	5 (5%)	26	22
1	C	106/126 (84%)	99 (93%)	7 (7%)	16	12
1	D	107/126 (85%)	99 (92%)	8 (8%)	13	9
All	All	426/504 (84%)	400 (94%)	26 (6%)	19	14

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

Mol	Chain	Res	Type
1	A	6	LEU
1	A	7	VAL
1	A	55	PRO
1	A	73	LEU
1	A	94[A]	LEU
1	A	94[B]	LEU
1	B	3	GLU
1	B	6	LEU

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Mol	Chain	Res	Type
1	B	73	LEU
1	B	111	GLU
1	B	120	LEU
1	C	2	VAL
1	C	3	GLU
1	C	5	SER
1	C	45	LYS
1	C	70	THR
1	C	127	GLU
1	C	137	LEU
1	D	6	LEU
1	D	9	TYR
1	D	15[A]	ARG
1	D	15[B]	ARG
1	D	26	GLU
1	D	73	LEU
1	D	135	GLU
1	D	137	LEU

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

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

Of 5 ligands modelled in this entry, 5 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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 ⓘ

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	133/165 (80%)	-0.15	5 (3%) 40 39	15, 23, 41, 63	0
1	B	134/165 (81%)	-0.29	0 100 100	16, 24, 40, 65	0
1	C	136/165 (82%)	0.05	11 (8%) 12 11	19, 29, 55, 67	0
1	D	134/165 (81%)	0.17	11 (8%) 11 11	20, 30, 65, 72	0
All	All	537/660 (81%)	-0.05	27 (5%) 28 28	15, 26, 54, 72	0

All (27) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	9	TYR	7.2
1	A	138	ALA	5.3
1	D	137	LEU	5.1
1	D	11	ALA	5.0
1	C	136	ALA	4.1
1	C	9	TYR	4.0
1	C	91	ARG	3.9
1	A	137	LEU	3.6
1	C	137	LEU	3.5
1	C	5	SER	3.3
1	D	7	VAL	3.2
1	C	7	VAL	3.1
1	D	6	LEU	3.0
1	D	136	ALA	2.9
1	D	10	GLY	2.9
1	C	3	GLU	2.8
1	C	135	GLU	2.8
1	D	5	SER	2.7
1	A	135	GLU	2.6
1	C	15	ARG	2.3
1	D	134	GLU	2.3

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Mol	Chain	Res	Type	RSRZ
1	D	135	GLU	2.2
1	C	6	LEU	2.1
1	A	9	TYR	2.1
1	C	134	GLU	2.1
1	D	12	PRO	2.1
1	A	6	LEU	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 carbohydrates in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q<0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	MG	D	164	1/1	0.78	0.07	47,47,47,47	0
2	MG	C	164	1/1	0.93	0.05	29,29,29,29	0
2	MG	D	1305	1/1	0.96	0.10	53,53,53,53	1
2	MG	B	164	1/1	0.99	0.08	29,29,29,29	0
2	MG	A	164	1/1	0.99	0.03	21,21,21,21	0

6.5 Other polymers [i](#)

There are no such residues in this entry.