



Full wwPDB EM Validation Report ⓘ

Mar 27, 2026 – 02:09 PM UTC

PDB ID : 2QU4 / pdb_00002qu4
Title : Model for Bacterial ParM Filament
Authors : Orlova, A.; Garner, E.C.; Galkin, V.E.; Heuser, J.; Mullins, R.D.; Egelman, E.H.
Deposited on : 2007-08-03
Resolution : 16.00 Å (reported)
Based on initial models : 1MWM, 1MWK

This is a Full wwPDB EM Validation Report for a publicly released PDB/EMDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/EMValidationReportHelp>

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
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

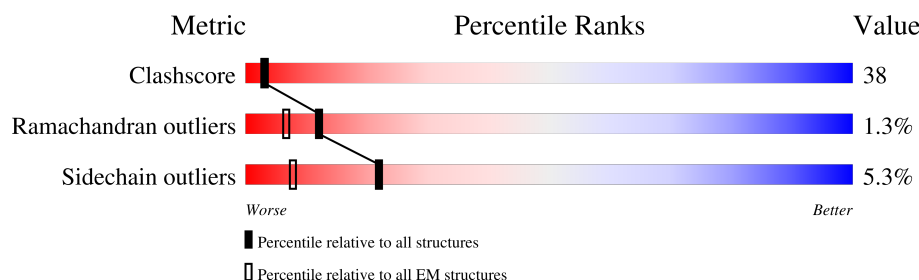
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 16.00 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)	EM structures (#Entries)
Clashscore	229148	23984
Ramachandran outliers	224038	23583
Sidechain outliers	223484	23102

The table below summarises the geometric issues observed across the polymeric chains and their fit to the map. The red, orange, yellow and green segments of the 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 $\leq 5\%$.

Mol	Chain	Length	Quality of chain
1	A	320	70% 27% .

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 2517 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, 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.

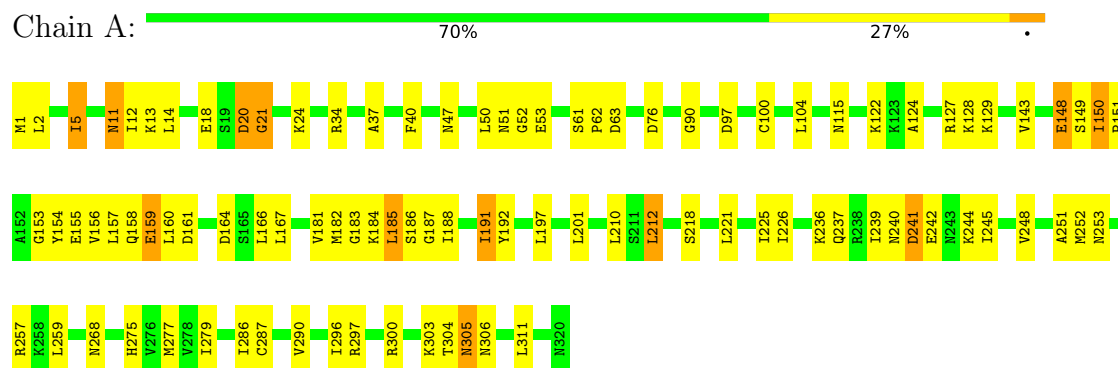
- Molecule 1 is a protein called Plasmid segregation protein parM.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	320	Total	C	N	O	S	0	0
			2517	1586	425	498	8		

3 Residue-property plots

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. 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. 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: Plasmid segregation protein parM



4 Data and refinement statistics

Xtriage (Phenix) and EDS failed to run properly - this section is therefore incomplete.

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	1.00Å 1.00Å 1.00Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	(Not available) – 16.00	Depositor
% Data completeness (in resolution range)	(Not available) ((Not available)-16.00)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	unknown	Depositor
R, R_{free}	(Not available) , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	2517	wwPDB-VP
Average B, all atoms (Å ²)	30.0	wwPDB-VP

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	0.44	0/2556	0.91	8/3457 (0.2%)

There are no bond length outliers.

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	149	SER	N-CA-C	7.98	119.98	111.28
1	A	5	ILE	N-CA-C	6.89	117.96	107.77
1	A	90	GLY	N-CA-C	-6.66	106.55	114.48
1	A	40	PHE	N-CA-C	-6.14	105.65	113.02
1	A	148	GLU	N-CA-C	5.59	122.70	110.80
1	A	150	ILE	N-CA-C	5.32	120.38	108.88
1	A	251	ALA	N-CA-C	-5.29	105.59	111.36
1	A	97	ASP	N-CA-C	-5.05	102.03	110.17

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	2517	0	2499	190	0
All	All	2517	0	2499	190	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 38.

All (190) 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:159:GLU:HG2	1:A:183:GLY:CA	1.31	1.59
1:A:159:GLU:CG	1:A:183:GLY:HA3	1.10	1.54
1:A:156:VAL:CG2	1:A:166:LEU:HB3	1.42	1.50
1:A:159:GLU:CG	1:A:183:GLY:CA	1.82	1.47
1:A:160:LEU:HA	1:A:182:MET:CE	1.22	1.43
1:A:159:GLU:HB2	1:A:164:ASP:N	1.31	1.40
1:A:153:GLY:O	1:A:181:VAL:CG1	1.68	1.38
1:A:279:ILE:HG21	1:A:306:ASN:N	1.05	1.36
1:A:159:GLU:HG3	1:A:183:GLY:N	1.42	1.34
1:A:279:ILE:HG21	1:A:306:ASN:CA	1.55	1.34
1:A:160:LEU:CA	1:A:182:MET:HE2	1.26	1.34
1:A:154:TYR:CD1	1:A:185:LEU:O	1.81	1.33
1:A:157:LEU:CA	1:A:182:MET:HB2	1.61	1.28
1:A:160:LEU:CA	1:A:182:MET:CE	1.79	1.25
1:A:279:ILE:CG2	1:A:306:ASN:N	2.00	1.24
1:A:154:TYR:HD1	1:A:185:LEU:O	1.11	1.21
1:A:156:VAL:HG22	1:A:166:LEU:CB	1.70	1.20
1:A:279:ILE:CG2	1:A:306:ASN:HA	1.73	1.18
1:A:279:ILE:CG2	1:A:306:ASN:CA	2.23	1.16
1:A:157:LEU:HA	1:A:182:MET:CB	1.75	1.14
1:A:159:GLU:CB	1:A:164:ASP:N	2.12	1.12
1:A:153:GLY:C	1:A:185:LEU:HD23	1.71	1.11
1:A:153:GLY:C	1:A:185:LEU:CD2	1.95	1.11
1:A:153:GLY:C	1:A:181:VAL:HG11	1.75	1.10
1:A:153:GLY:O	1:A:181:VAL:HG11	1.40	1.07
1:A:160:LEU:HD13	1:A:182:MET:HA	1.19	1.07
1:A:158:GLN:HA	1:A:186:SER:HB3	1.12	1.07
1:A:279:ILE:HD13	1:A:306:ASN:N	1.69	1.07
1:A:154:TYR:CG	1:A:184:LYS:O	2.08	1.06
1:A:159:GLU:HG3	1:A:183:GLY:CA	1.64	1.06
1:A:156:VAL:HB	1:A:181:VAL:HG12	1.37	1.04
1:A:157:LEU:HD12	1:A:185:LEU:CD2	1.88	1.04
1:A:158:GLN:HA	1:A:186:SER:CB	1.87	1.03
1:A:154:TYR:CD2	1:A:184:LYS:O	2.11	1.02
1:A:158:GLN:N	1:A:186:SER:N	2.08	1.02
1:A:279:ILE:HG22	1:A:306:ASN:HA	1.38	1.01
1:A:160:LEU:CA	1:A:182:MET:HE3	1.68	1.00
1:A:160:LEU:HD13	1:A:182:MET:CA	1.75	1.00
1:A:156:VAL:HG22	1:A:166:LEU:HB3	1.05	1.00
1:A:156:VAL:HB	1:A:181:VAL:CG1	1.69	0.96

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:158:GLN:CA	1:A:186:SER:HB3	1.96	0.96
1:A:153:GLY:O	1:A:181:VAL:HG12	1.65	0.96
1:A:153:GLY:CA	1:A:181:VAL:HG11	1.94	0.95
1:A:156:VAL:CG2	1:A:166:LEU:CB	2.35	0.94
1:A:156:VAL:CG2	1:A:181:VAL:HB	1.92	0.94
1:A:160:LEU:N	1:A:182:MET:HE3	1.82	0.94
1:A:159:GLU:H	1:A:183:GLY:N	1.66	0.92
1:A:37:ALA:H	1:A:47:ASN:HD21	1.18	0.91
1:A:159:GLU:CG	1:A:183:GLY:N	2.18	0.91
1:A:157:LEU:HD12	1:A:185:LEU:HD22	1.50	0.91
1:A:159:GLU:H	1:A:183:GLY:CA	1.84	0.91
1:A:156:VAL:HG21	1:A:166:LEU:HB3	1.53	0.90
1:A:158:GLN:HG3	1:A:184:LYS:O	1.30	0.90
1:A:153:GLY:O	1:A:185:LEU:HA	1.70	0.89
1:A:157:LEU:CA	1:A:182:MET:CB	2.43	0.88
1:A:159:GLU:HG3	1:A:183:GLY:H	1.08	0.88
1:A:155:GLU:HA	1:A:184:LYS:CA	2.03	0.88
1:A:157:LEU:HA	1:A:182:MET:HB2	0.88	0.87
1:A:153:GLY:O	1:A:185:LEU:HD23	1.76	0.85
1:A:157:LEU:HD12	1:A:185:LEU:HD23	1.57	0.84
1:A:154:TYR:CD1	1:A:184:LYS:O	2.28	0.84
1:A:226:ILE:HG12	1:A:252:MET:HE2	1.57	0.83
1:A:279:ILE:CD1	1:A:306:ASN:N	2.41	0.82
1:A:154:TYR:CD1	1:A:185:LEU:C	2.56	0.82
1:A:158:GLN:CA	1:A:186:SER:CB	2.50	0.81
1:A:225:ILE:HG22	1:A:252:MET:HE1	1.64	0.80
1:A:155:GLU:HA	1:A:184:LYS:HA	1.63	0.80
1:A:210:LEU:HB3	1:A:212:LEU:HD13	1.64	0.79
1:A:305:ASN:HA	1:A:306:ASN:O	1.83	0.78
1:A:157:LEU:C	1:A:182:MET:HB2	2.08	0.78
1:A:154:TYR:CE2	1:A:184:LYS:O	2.36	0.78
1:A:159:GLU:CG	1:A:183:GLY:H	1.89	0.78
1:A:277:MET:CE	1:A:306:ASN:HB2	2.15	0.77
1:A:158:GLN:CG	1:A:184:LYS:O	2.21	0.76
1:A:159:GLU:HB2	1:A:164:ASP:CA	2.15	0.76
1:A:160:LEU:CD1	1:A:182:MET:HA	2.05	0.76
1:A:160:LEU:N	1:A:182:MET:HB3	1.99	0.76
1:A:160:LEU:H	1:A:182:MET:HB3	1.53	0.74
1:A:156:VAL:HG21	1:A:181:VAL:HB	1.70	0.72
1:A:159:GLU:N	1:A:183:GLY:CA	2.53	0.72
1:A:156:VAL:CB	1:A:181:VAL:HG12	2.18	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:210:LEU:CB	1:A:212:LEU:HD13	2.20	0.71
1:A:150:ILE:O	1:A:185:LEU:HD21	1.91	0.70
1:A:156:VAL:HG11	1:A:181:VAL:N	1.67	0.69
1:A:11:ASN:C	1:A:11:ASN:HD22	1.99	0.68
1:A:159:GLU:C	1:A:182:MET:HE3	2.18	0.68
1:A:244:LYS:O	1:A:248:VAL:HG23	1.94	0.68
1:A:287:CYS:SG	1:A:303:LYS:HD2	2.34	0.68
1:A:277:MET:HE2	1:A:306:ASN:HB2	1.75	0.68
1:A:153:GLY:O	1:A:181:VAL:HG13	1.87	0.67
1:A:160:LEU:HA	1:A:182:MET:HE2	0.81	0.67
1:A:37:ALA:H	1:A:47:ASN:ND2	1.91	0.66
1:A:153:GLY:C	1:A:185:LEU:HD21	2.17	0.65
1:A:157:LEU:C	1:A:182:MET:CB	2.69	0.65
1:A:155:GLU:C	1:A:185:LEU:N	2.52	0.64
1:A:157:LEU:HD11	1:A:188:ILE:CD1	2.29	0.62
1:A:253:ASN:O	1:A:257:ARG:HG3	1.99	0.62
1:A:154:TYR:CE1	1:A:184:LYS:O	2.52	0.61
1:A:157:LEU:CD1	1:A:185:LEU:HD23	2.30	0.61
1:A:61:SER:C	1:A:63:ASP:H	2.08	0.61
1:A:187:GLY:C	1:A:188:ILE:HD12	2.26	0.60
1:A:157:LEU:CD1	1:A:185:LEU:CD2	2.73	0.60
1:A:154:TYR:CZ	1:A:184:LYS:O	2.55	0.60
1:A:156:VAL:HG23	1:A:166:LEU:HB3	1.70	0.59
1:A:153:GLY:C	1:A:181:VAL:CG1	2.50	0.59
1:A:275:HIS:HA	1:A:296:ILE:HD11	1.83	0.59
1:A:157:LEU:CA	1:A:182:MET:CA	2.77	0.59
1:A:34:ARG:HH11	1:A:53:GLU:HG3	1.67	0.59
1:A:156:VAL:HG13	1:A:166:LEU:H	1.68	0.58
1:A:37:ALA:N	1:A:47:ASN:HD21	1.96	0.58
1:A:212:LEU:HB3	1:A:218:SER:HB2	1.84	0.58
1:A:226:ILE:CG1	1:A:252:MET:HE2	2.31	0.58
1:A:157:LEU:H	1:A:181:VAL:HG12	1.69	0.58
1:A:275:HIS:CD2	1:A:300:ARG:HG3	2.39	0.57
1:A:159:GLU:CD	1:A:183:GLY:HA3	2.17	0.57
1:A:153:GLY:HA2	1:A:181:VAL:HG11	1.85	0.57
1:A:159:GLU:N	1:A:183:GLY:C	2.60	0.57
1:A:277:MET:HE2	1:A:304:THR:OG1	2.05	0.56
1:A:296:ILE:HD11	1:A:300:ARG:HB3	1.87	0.56
1:A:157:LEU:HD21	1:A:188:ILE:HD12	1.88	0.56
1:A:239:ILE:HB	1:A:245:ILE:CD1	2.36	0.56
1:A:1:MET:HE3	1:A:1:MET:HA	1.88	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:156:VAL:HA	1:A:182:MET:C	2.22	0.55
1:A:241:ASP:HB3	1:A:244:LYS:HB2	1.88	0.55
1:A:156:VAL:CG1	1:A:182:MET:CA	2.57	0.55
1:A:153:GLY:CA	1:A:185:LEU:CD2	2.80	0.55
1:A:11:ASN:C	1:A:12:ILE:HD12	2.32	0.54
1:A:155:GLU:CA	1:A:184:LYS:CA	2.80	0.54
1:A:157:LEU:O	1:A:182:MET:CB	2.57	0.53
1:A:159:GLU:N	1:A:183:GLY:N	2.49	0.53
1:A:157:LEU:HA	1:A:182:MET:CA	2.37	0.53
1:A:241:ASP:O	1:A:245:ILE:HG12	2.10	0.52
1:A:157:LEU:N	1:A:181:VAL:HG12	2.26	0.51
1:A:159:GLU:HB3	1:A:164:ASP:N	2.20	0.49
1:A:201:LEU:C	1:A:201:LEU:HD23	2.38	0.49
1:A:153:GLY:CA	1:A:185:LEU:HD21	2.43	0.49
1:A:51:ASN:HD21	1:A:128:LYS:CD	2.26	0.49
1:A:11:ASN:C	1:A:11:ASN:ND2	2.70	0.48
1:A:157:LEU:H	1:A:185:LEU:HA	1.77	0.48
1:A:158:GLN:C	1:A:186:SER:OG	2.54	0.48
1:A:236:LYS:HE2	1:A:242:GLU:OE1	2.13	0.48
1:A:239:ILE:HB	1:A:245:ILE:HD11	1.96	0.48
1:A:115:ASN:C	1:A:115:ASN:HD22	2.22	0.47
1:A:296:ILE:CD1	1:A:300:ARG:HB3	2.44	0.47
1:A:159:GLU:HB2	1:A:164:ASP:O	2.14	0.47
1:A:157:LEU:O	1:A:182:MET:HB3	2.14	0.47
1:A:157:LEU:O	1:A:182:MET:HB2	2.13	0.47
1:A:279:ILE:CB	1:A:306:ASN:N	2.74	0.47
1:A:157:LEU:HD11	1:A:188:ILE:HD12	1.96	0.47
1:A:240:ASN:O	1:A:241:ASP:HB2	2.14	0.47
1:A:156:VAL:HG22	1:A:166:LEU:CA	2.41	0.47
1:A:157:LEU:N	1:A:182:MET:N	2.48	0.47
1:A:129:LYS:H	1:A:129:LYS:HG2	1.54	0.46
1:A:11:ASN:HD21	1:A:13:LYS:NZ	2.14	0.46
1:A:155:GLU:CA	1:A:184:LYS:C	2.87	0.46
1:A:5:ILE:HG12	1:A:14:LEU:HD22	1.97	0.46
1:A:188:ILE:HG21	1:A:191:ILE:HD13	1.98	0.46
1:A:239:ILE:HB	1:A:245:ILE:HD13	1.98	0.46
1:A:159:GLU:CG	1:A:183:GLY:HA2	2.23	0.45
1:A:192:TYR:CD2	1:A:192:TYR:C	2.95	0.45
1:A:20:ASP:CG	1:A:21:GLY:H	2.25	0.45
1:A:61:SER:C	1:A:63:ASP:N	2.71	0.45
1:A:127:ARG:HH22	1:A:143:VAL:HB	1.81	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:156:VAL:HG22	1:A:166:LEU:HB2	1.84	0.44
1:A:157:LEU:CD1	1:A:185:LEU:HD22	2.36	0.44
1:A:157:LEU:CG	1:A:188:ILE:HD12	2.48	0.44
1:A:18:GLU:HG3	1:A:24:LYS:HE3	1.99	0.44
1:A:50:LEU:HD23	1:A:76:ASP:HB3	1.99	0.44
1:A:156:VAL:CA	1:A:182:MET:C	2.01	0.44
1:A:277:MET:HE2	1:A:306:ASN:CB	2.45	0.44
1:A:51:ASN:HD21	1:A:128:LYS:HD3	1.82	0.44
1:A:153:GLY:CA	1:A:181:VAL:CG1	2.82	0.44
1:A:210:LEU:HB2	1:A:212:LEU:HD13	1.97	0.43
1:A:100:CYS:SG	1:A:122:LYS:HE3	2.59	0.43
1:A:268:ASN:HD22	1:A:268:ASN:HA	1.66	0.43
1:A:154:TYR:HB2	1:A:185:LEU:HD13	1.15	0.43
1:A:157:LEU:C	1:A:182:MET:C	2.79	0.42
1:A:297:ARG:HD3	1:A:300:ARG:HD3	2.02	0.42
1:A:148:GLU:O	1:A:151:PRO:HD2	2.20	0.42
1:A:155:GLU:HA	1:A:184:LYS:C	2.43	0.42
1:A:158:GLN:HB3	1:A:184:LYS:HD2	1.86	0.41
1:A:210:LEU:HD21	1:A:248:VAL:HG21	2.02	0.41
1:A:210:LEU:HB3	1:A:212:LEU:CD1	2.40	0.41
1:A:286:ILE:O	1:A:290:VAL:HG13	2.21	0.41
1:A:124:ALA:O	1:A:127:ARG:HB2	2.21	0.41
1:A:159:GLU:HB2	1:A:164:ASP:C	2.46	0.41
1:A:12:ILE:HD12	1:A:12:ILE:N	2.36	0.40
1:A:150:ILE:O	1:A:185:LEU:CD2	2.66	0.40
1:A:61:SER:HA	1:A:62:PRO:HD3	1.92	0.40
1:A:305:ASN:CA	1:A:306:ASN:O	2.63	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 PDB entries followed by that with respect to all EM entries.

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	314/320 (98%)	299 (95%)	11 (4%)	4 (1%)	9	42

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	20	ASP
1	A	52	GLY
1	A	241	ASP
1	A	21	GLY

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 PDB entries followed by that with respect to all EM entries.

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	284/284 (100%)	269 (95%)	15 (5%)	20	41

All (15) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	2	LEU
1	A	11	ASN
1	A	104	LEU
1	A	159	GLU
1	A	161	ASP
1	A	167	LEU
1	A	185	LEU
1	A	191	ILE
1	A	197	LEU
1	A	212	LEU
1	A	221	LEU
1	A	237	GLN
1	A	259	LEU
1	A	305	ASN
1	A	311	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (17) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	11	ASN
1	A	25	GLN
1	A	47	ASN
1	A	51	ASN
1	A	54	GLN
1	A	73	GLN
1	A	83	HIS
1	A	84	HIS
1	A	111	ASN
1	A	115	ASN
1	A	118	ASN
1	A	228	HIS
1	A	237	GLN
1	A	240	ASN
1	A	253	ASN
1	A	268	ASN
1	A	308	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

The following chains have linkage breaks:

Mol	Chain	Number of breaks
1	A	2

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	A	163:LEU	C	164:ASP	N	12.40
1	A	305:ASN	C	306:ASN	N	7.02