wwPDB EM Map/Model Validation Report (i)

Apr 10, 2016 – 01:35 PM BST

PDB ID	:	2QU4
EMDB ID:	:	unknown
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)

This is a wwPDB EM Map/Model Validation Report for a publicly released PDB/EMDB entry.
For rigid body fitted models, validation errors reported here could
stem from errors in the original structure(s) used in the fitting.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at
http://wwpdb.org/validation/2016/EMValidationReportHelp

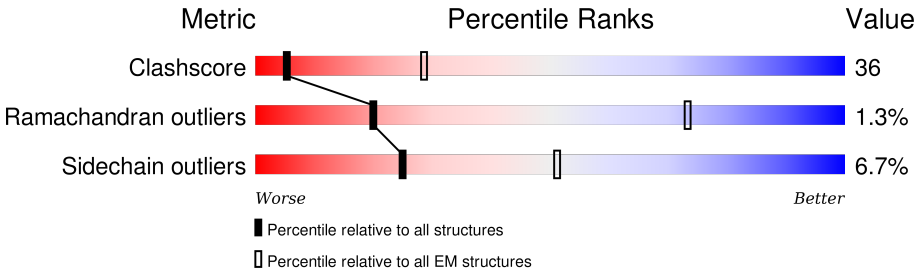
MolProbity	:	4.02b-467
Mogul	:	unknown
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP)	:	trunk27241

1 Overall quality at a glance (i)

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	114402	924
Ramachandran outliers	111179	726
Sidechain outliers	111093	686

The table below summarises the geometric issues observed across the polymeric chains. The red, orange, yellow and green segments on the bar indicate the fraction of residues that contain outliers for >=3, 2, 1 and 0 types of geometric quality criteria. 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 <=5%

Mol	Chain	Length	Quality of chain
1	A	320	71% 26% .

2 Entry composition (i)

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.

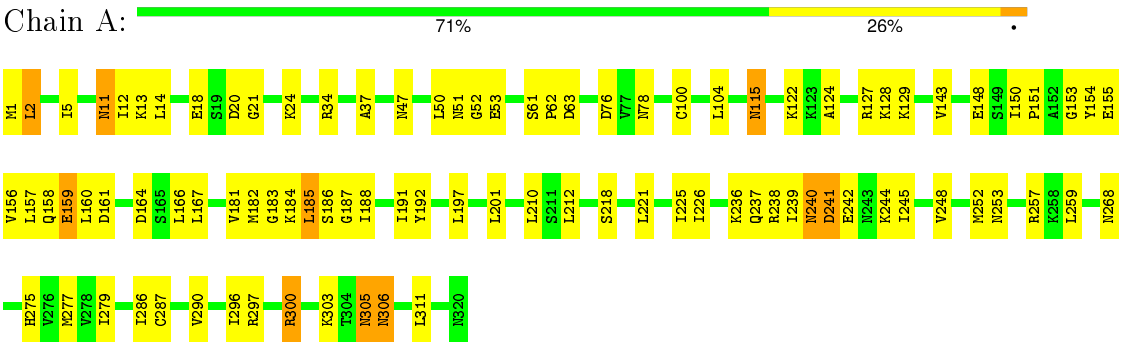
Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
1	A	320	2517	1586	425	498	8	0	0

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 errors 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

Chain A:



4 Experimental information (i)

Property	Value	Source
Reconstruction method	HELICAL	Depositor
Imposed symmetry	POINT, Not provided	Depositor
Number of images	Not provided	Depositor
Resolution determination method	Not provided	Depositor
CTF correction method	Not provided	Depositor
Microscope	Not provided	Depositor
Voltage (kV)	200	Depositor
Electron dose ($e^-/Å^2$)	Not provided	Depositor
Minimum defocus (nm)	Not provided	Depositor
Maximum defocus (nm)	Not provided	Depositor
Magnification	Not provided	Depositor
Image detector	Not provided	Depositor

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	$\#\|Z\| > 2$	RMSZ	$\#\|Z\| > 2$
1	A	0.35	0/2556	0.61	0/3457

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

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	182	0
All	All	2517	0	2499	182	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 36.

All (182) 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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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: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:160:LEU:CA	1:A:182:MET:HE2	1.38	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:160:LEU:CA	1:A:182:MET:HE3	1.56	1.09
1:A:153:GLY:C	1:A:185:LEU:CD2	1.95	1.09
1:A:153:GLY:C	1:A:185:LEU:HD23	1.71	1.08
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:153:GLY:C	1:A:181:VAL:HG11	1.75	1.05
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:160:LEU:N	1:A:182:MET:HE3	1.73	1.01
1:A:279:ILE:HG22	1:A:306:ASN:HA	1.38	1.01
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
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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:159:GLU:H	1:A:183:GLY:N	1.66	0.92
1:A:159:GLU:CG	1:A:183:GLY:N	2.18	0.91
1:A:37:ALA:H	1:A:47:ASN:HD21	1.18	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:279:ILE:CD1	1:A:306:ASN:N	2.41	0.82
1:A:158:GLN:CA	1:A:186:SER:CB	2.50	0.81
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:160:LEU:HA	1:A:182:MET:HE2	0.86	0.79
1:A:305:ASN:HA	1:A:306:ASN:O	1.83	0.78
1:A:154:TYR:CD1	1:A:185:LEU:C	2.56	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:277:MET:HE1	1:A:306:ASN:HB2	1.68	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:157:LEU:C	1:A:182:MET:HB2	2.08	0.73
1:A:159:GLU:N	1:A:183:GLY:CA	2.53	0.72
1:A:156:VAL:HG21	1:A:181:VAL:HB	1.70	0.72
1:A:156:VAL:CB	1:A:181:VAL:HG12	2.18	0.71
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: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:153:GLY:O	1:A:181:VAL:HG13	1.87	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:37:ALA:H	1:A:47:ASN:ND2	1.91	0.66
1:A:159:GLU:C	1:A:182:MET:HE3	2.17	0.64
1:A:11:ASN:C	1:A:11:ASN:HD22	1.99	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:155:GLU:C	1:A:185:LEU:N	2.52	0.61
1:A:157:LEU:CD1	1:A:185:LEU:HD23	2.30	0.61
1:A:157:LEU:CD1	1:A:185:LEU:CD2	2.73	0.60
1:A:157:LEU:C	1:A:182:MET:CB	2.69	0.60
1:A:153:GLY:CA	1:A:185:LEU:CD2	2.80	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: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: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:61:SER:C	1:A:63:ASP:H	2.08	0.57
1:A:153:GLY:HA2	1:A:181:VAL:HG11	1.85	0.57
1:A:226:ILE:HG12	1:A:252:MET:HE2	1.86	0.57
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:187:GLY:C	1:A:188:ILE:HD12	2.26	0.56
1:A:239:ILE:HB	1:A:245:ILE:CD1	2.36	0.56
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:C	1:A:181:VAL:CG1	2.50	0.55
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:156:VAL:CA	1:A:182:MET:C	2.01	0.51
1:A:157:LEU:N	1:A:181:VAL:HG12	2.26	0.51
1:A:225:ILE:HG22	1:A:252:MET:CE	2.41	0.50
1:A:1:MET:HE2	1:A:2:LEU:H	1.77	0.50
1:A:11:ASN:C	1:A:12:ILE:HD12	2.32	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:159:GLU:HB3	1:A:164:ASP:N	2.20	0.49
1:A:153:GLY:CA	1:A:185:LEU:HD21	2.43	0.49
1:A:226:ILE:HA	1:A:252:MET:HE1	1.94	0.49
1:A:51:ASN:HD21	1:A:128:LYS:CD	2.26	0.49
1:A:157:LEU:H	1:A:185:LEU:HA	1.77	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: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:HG2	1:A:129:LYS:H	1.54	0.46
1:A:11:ASN:HD21	1:A:13:LYS:NZ	2.14	0.46
1:A:188:ILE:HG21	1:A:191:ILE:HD13	1.98	0.46
1:A:5:ILE:HG12	1:A:14:LEU:HD22	1.97	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:201:LEU:C	1:A:201:LEU:HD23	2.38	0.45
1:A:127:ARG:HH22	1:A:143:VAL:HB	1.81	0.44
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: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:115:ASN:C	1:A:115:ASN:HD22	2.22	0.43
1:A:158:GLN:C	1:A:186:SER:OG	2.54	0.43
1:A:61:SER:C	1:A:63:ASP:N	2.71	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: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:CA	1:A:184:LYS:C	2.87	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:225:ILE:HG22	1:A:252:MET:HE3	2.02	0.41
1:A:124:ALA:O	1:A:127:ARG:HB2	2.21	0.41
1:A:192:TYR:C	1:A:192:TYR:CD2	2.95	0.41
1:A:12:ILE:N	1:A:12:ILE:HD12	2.36	0.40
1:A:226:ILE:CA	1:A:252:MET:HE1	2.50	0.40
1:A:150:ILE:O	1:A:185:LEU:CD2	2.66	0.40
1:A:305:ASN:CA	1:A:306:ASN:O	2.63	0.40
1:A:11:ASN:ND2	1:A:11:ASN:C	2.70	0.40
1:A:61:SER:HA	1:A:62:PRO:HD3	1.92	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%)	15 60

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 (i)

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%)	265 (93%)	19 (7%)	20 57

All (19) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	2	LEU
1	A	11	ASN
1	A	78	ASN
1	A	104	LEU
1	A	115	ASN
1	A	159	GLU
1	A	161	ASP
1	A	167	LEU
1	A	185	LEU
1	A	197	LEU
1	A	221	LEU
1	A	237	GLN
1	A	238	ARG
1	A	240	ASN
1	A	259	LEU
1	A	300	ARG
1	A	305	ASN
1	A	306	ASN
1	A	311	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) 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

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Mol	Chain	Res	Type
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 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.