



Full wwPDB/EMDataBank EM Map/Model Validation Report ⓘ

Mar 2, 2017 – 11:21 am GMT

PDB ID : 3J06
EMDB ID: : EMD-5185
Title : CryoEM Helical Reconstruction of TMV
Authors : Ge, P.; Zhou, Z.H.
Deposited on : 2011-04-26
Resolution : 3.30 Å(reported)

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

We welcome your comments at validation@mail.wwpdb.org
A user guide is available at
<http://wwpdb.org/validation/2016/EMValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

MolProbity : 4.02b-467
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : recalc29047

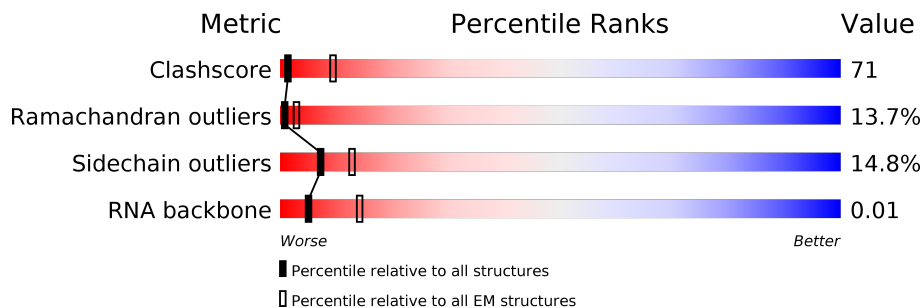
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 3.30 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)	EM structures (#Entries)
Clashscore	125131	1336
Ramachandran outliers	121729	1120
Sidechain outliers	121581	1026
RNA backbone	3398	335

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 $\leq 5\%$.

Mol	Chain	Length	Quality of chain
1	A	159	 28% 45% 21% . .
2	R	3	 100%

2 Entry composition [i](#)

There are 2 unique types of molecules in this entry. The entry contains 1284 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 Coat protein.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	155	Total	C	N	O	S	1	0
			1219	765	214	239	1		

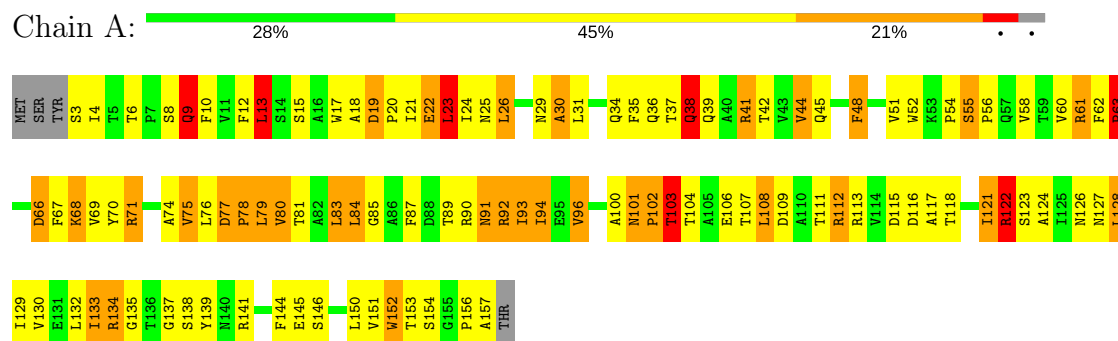
- Molecule 2 is a RNA chain called 5'-R(P*AP*UP*G)-3'.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	R	3	Total	C	N	O	P	0	0
			65	29	12	21	3		

3 Residue-property plots

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. 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: Coat protein



- Molecule 2: 5'-R(P*AP*UP*G)-3'



4 Experimental information

Property	Value	Source
Reconstruction method	HELICAL	Depositor
Imposed symmetry	POINT, Not provided	Depositor
Number of segments used	Not provided	Depositor
Resolution determination method	Truncation	Depositor
CTF correction method	Not provided	Depositor
Microscope	FEI TITAN KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	2500	Depositor
Minimum defocus (nm)	1200	Depositor
Maximum defocus (nm)	2500	Depositor
Magnification	73000	Depositor
Image detector	Kodak SO163 film	Depositor

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	$\# Z > 2$	RMSZ	$\# Z > 2$
1	A	0.54	0/1243	1.05	11/1699 (0.6%)
2	R	0.64	0/72	1.20	0/110
All	All	0.54	0/1315	1.06	11/1809 (0.6%)

There are no bond length outliers.

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	26	LEU	CB-CA-C	9.53	128.31	110.20
1	A	41	ARG	CB-CA-C	9.20	128.81	110.40
1	A	9	GLN	CB-CA-C	-7.16	96.07	110.40
1	A	38	GLN	CB-CA-C	7.14	124.67	110.40
1	A	44	VAL	CB-CA-C	-6.92	98.26	111.40
1	A	22	GLU	CB-CA-C	6.20	122.81	110.40
1	A	80	VAL	CB-CA-C	-6.02	99.96	111.40
1	A	23	LEU	N-CA-C	5.95	127.08	111.00
1	A	42	THR	N-CA-C	5.81	126.69	111.00
1	A	30	ALA	CB-CA-C	5.74	118.72	110.10
1	A	13	LEU	N-CA-C	5.21	125.07	111.00

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	1219	0	1196	179	0
2	R	65	0	33	3	0
All	All	1284	0	1229	179	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 71.

All (179) 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:112:ARG:O	1:A:116:ASP:HB2	1.62	0.99
1:A:18:ALA:O	1:A:68:LYS:HG2	1.63	0.96
1:A:122:ARG:HH11	1:A:122:ARG:HB2	1.30	0.94
1:A:93:ILE:HG22	1:A:94:ILE:HD13	1.48	0.93
1:A:150:LEU:HD13	1:A:157:ALA:HB3	1.50	0.90
1:A:13:LEU:HD23	1:A:13:LEU:N	1.86	0.88
1:A:121:ILE:O	1:A:123:SER:N	2.10	0.85
1:A:19:ASP:HB3	1:A:22:GLU:HB3	1.59	0.84
1:A:61:ARG:HG2	1:A:62:PHE:H	1.44	0.83
1:A:61:ARG:HD2	1:A:153:THR:HG23	1.59	0.81
1:A:45:GLN:HG3	1:A:87:PHE:HB2	1.62	0.81
1:A:3:SER:HA	1:A:156:PRO:CG	2.10	0.81
1:A:151:VAL:HG23	1:A:152:TRP:H	1.45	0.81
1:A:115:ASP:O	1:A:118:THR:HG22	1.80	0.81
1:A:61:ARG:HD2	1:A:153:THR:CG2	2.12	0.80
1:A:122:ARG:NH1	1:A:122:ARG:HB2	1.95	0.80
1:A:3:SER:HA	1:A:156:PRO:HG2	1.62	0.79
1:A:78:PRO:HA	1:A:81:THR:HG22	1.65	0.78
1:A:62:PHE:HD2	1:A:141:ARG:HA	1.49	0.76
1:A:80:VAL:O	1:A:80:VAL:HG12	1.84	0.76
1:A:45:GLN:HB2	1:A:87:PHE:CD2	2.21	0.76
1:A:126:ASN:O	1:A:129:ILE:HG22	1.86	0.76
1:A:129:ILE:O	1:A:133:ILE:HG22	1.85	0.76
1:A:122:ARG:HH11	1:A:122:ARG:CB	1.99	0.75
1:A:109:ASP:O	1:A:112:ARG:HB3	1.89	0.72
1:A:115:ASP:HA	1:A:118:THR:HG22	1.71	0.72
1:A:9:GLN:HG3	1:A:150:LEU:HD21	1.71	0.71
1:A:61:ARG:HB2	1:A:153:THR:HG23	1.72	0.71
1:A:134:ARG:HH11	1:A:134:ARG:HG3	1.55	0.70
1:A:121:ILE:O	1:A:124:ALA:N	2.23	0.70
1:A:63:PRO:HG2	1:A:68:LYS:HE3	1.72	0.70
1:A:103:THR:O	1:A:107:THR:HG22	1.91	0.70

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:62:PHE:CD2	1:A:141:ARG:HA	2.27	0.69
1:A:45:GLN:HB2	1:A:87:PHE:CG	2.27	0.69
1:A:18:ALA:HB2	1:A:52:TRP:CE3	2.27	0.68
1:A:133:ILE:C	1:A:135:GLY:H	1.96	0.68
1:A:61:ARG:HG2	1:A:62:PHE:N	2.08	0.68
1:A:67:PHE:O	1:A:68:LYS:HB2	1.94	0.67
1:A:115:ASP:CA	1:A:118:THR:HG22	2.26	0.66
1:A:126:ASN:O	1:A:129:ILE:CG2	2.44	0.66
1:A:100:ALA:O	1:A:101:ASN:HB2	1.95	0.66
1:A:26:LEU:HA	1:A:29:ASN:HD21	1.61	0.65
1:A:91:ASN:HD22	1:A:91:ASN:N	1.93	0.65
1:A:91:ASN:O	1:A:93:ILE:N	2.29	0.65
1:A:126:ASN:C	1:A:129:ILE:HG22	2.16	0.64
1:A:78:PRO:HA	1:A:81:THR:CG2	2.27	0.64
1:A:152:TRP:HE3	1:A:154:SER:H	1.41	0.64
1:A:130:VAL:HA	1:A:133:ILE:CG2	2.28	0.64
1:A:52:TRP:CE2	1:A:71:ARG:HG3	2.33	0.64
1:A:35:PHE:HD1	1:A:122:ARG:HH12	1.45	0.63
1:A:76:LEU:HD23	1:A:76:LEU:O	1.98	0.63
1:A:26:LEU:HA	1:A:29:ASN:ND2	2.15	0.62
1:A:69:VAL:HG11	1:A:132:LEU:HD22	1.79	0.62
1:A:133:ILE:C	1:A:135:GLY:N	2.51	0.62
1:A:150:LEU:HD13	1:A:157:ALA:CB	2.27	0.62
1:A:115:ASP:HA	1:A:118:THR:CG2	2.31	0.61
1:A:138:SER:O	1:A:139:TYR:CG	2.54	0.61
1:A:23:LEU:HD12	1:A:69:VAL:HG21	1.82	0.60
1:A:69:VAL:HA	1:A:137:GLY:O	2.02	0.60
1:A:67:PHE:O	1:A:68:LYS:CB	2.49	0.59
1:A:19:ASP:CB	1:A:22:GLU:HB3	2.31	0.59
1:A:61:ARG:HD3	1:A:141:ARG:NH1	2.18	0.59
1:A:56:PRO:HB3	1:A:62:PHE:HD1	1.68	0.59
1:A:41:ARG:HB2	1:A:41:ARG:HH11	1.68	0.58
1:A:18:ALA:HB2	1:A:52:TRP:CZ3	2.39	0.58
1:A:121:ILE:O	1:A:122:ARG:C	2.42	0.58
1:A:107:THR:O	1:A:109:ASP:N	2.37	0.57
1:A:55:SER:HB3	1:A:56:PRO:CD	2.34	0.57
1:A:141:ARG:HH11	1:A:141:ARG:HG3	1.69	0.57
1:A:9:GLN:CG	1:A:150:LEU:HD21	2.34	0.56
1:A:12:PHE:C	1:A:13:LEU:HD23	2.26	0.56
1:A:66:ASP:OD2	1:A:68:LYS:HE2	2.06	0.56
1:A:133:ILE:O	1:A:135:GLY:N	2.38	0.56

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:115:ASP:C	1:A:118:THR:HG22	2.25	0.56
1:A:75:VAL:CG2	1:A:76:LEU:N	2.69	0.55
1:A:36:GLN:NE2	1:A:118:THR:HG21	2.22	0.55
1:A:107:THR:HG23	1:A:108:LEU:N	2.21	0.55
1:A:107:THR:HG23	1:A:108:LEU:H	1.72	0.55
1:A:76:LEU:O	1:A:80:VAL:HG23	2.07	0.55
1:A:102:PRO:HA	1:A:106:GLU:OE2	2.08	0.54
1:A:126:ASN:HA	1:A:129:ILE:CG2	2.37	0.54
1:A:13:LEU:CD2	1:A:13:LEU:N	2.51	0.53
1:A:141:ARG:NH1	1:A:141:ARG:HG3	2.24	0.53
1:A:80:VAL:O	1:A:80:VAL:CG1	2.55	0.53
1:A:37:THR:O	1:A:39[A]:GLN:N	2.42	0.53
1:A:79:LEU:HD21	1:A:127:ASN:HB3	1.89	0.53
1:A:6:THR:HG22	1:A:8:SER:H	1.74	0.53
1:A:15:SER:OG	1:A:54:PRO:HD3	2.10	0.52
1:A:20:PRO:HG2	1:A:21:ILE:H	1.75	0.51
1:A:126:ASN:HA	1:A:129:ILE:HG22	1.92	0.51
1:A:37:THR:O	1:A:39[B]:GLN:N	2.43	0.50
1:A:45:GLN:HG3	1:A:87:PHE:CB	2.37	0.50
1:A:89:THR:HG23	1:A:113:ARG:NH2	2.25	0.50
1:A:61:ARG:CD	1:A:153:THR:HG23	2.35	0.50
1:A:55:SER:HB3	1:A:56:PRO:HD3	1.92	0.50
1:A:4:ILE:HD13	1:A:10:PHE:CE1	2.47	0.49
1:A:38:GLN:NE2	1:A:90:ARG:HG2	2.27	0.49
1:A:38:GLN:HE22	1:A:90:ARG:HG2	1.77	0.49
1:A:130:VAL:HG12	1:A:130:VAL:O	2.12	0.49
1:A:3:SER:HA	1:A:156:PRO:CB	2.42	0.49
1:A:130:VAL:HA	1:A:133:ILE:HG21	1.95	0.49
1:A:91:ASN:O	1:A:92:ARG:C	2.50	0.49
1:A:121:ILE:C	1:A:123:SER:N	2.66	0.49
1:A:51:VAL:HG12	1:A:51:VAL:O	2.12	0.49
1:A:90:ARG:NH2	1:A:90:ARG:HB3	2.27	0.49
1:A:61:ARG:CB	1:A:153:THR:HG23	2.42	0.48
1:A:58:VAL:O	1:A:58:VAL:HG12	2.12	0.48
1:A:21:ILE:HG23	1:A:22:GLU:N	2.27	0.48
1:A:100:ALA:O	1:A:101:ASN:CB	2.62	0.48
1:A:117:ALA:HB2	2:R:6:G:O6	2.14	0.48
1:A:116:ASP:HA	2:R:4:A:O2'	2.14	0.48
1:A:44:VAL:O	1:A:48:PHE:HB2	2.14	0.48
1:A:3:SER:HA	1:A:156:PRO:CD	2.42	0.47
1:A:121:ILE:HB	1:A:122:ARG:H	1.41	0.47

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:23:LEU:HD23	1:A:23:LEU:O	2.14	0.47
1:A:107:THR:C	1:A:109:ASP:N	2.68	0.47
1:A:117:ALA:HB2	2:R:6:G:C6	2.50	0.47
1:A:25:ASN:O	1:A:29:ASN:ND2	2.48	0.47
1:A:38:GLN:OE1	1:A:38:GLN:O	2.33	0.47
1:A:24:ILE:CG1	1:A:25:ASN:N	2.77	0.47
1:A:69:VAL:HG11	1:A:132:LEU:CD2	2.44	0.46
1:A:90:ARG:CZ	1:A:90:ARG:HB3	2.45	0.46
1:A:126:ASN:CA	1:A:129:ILE:HG22	2.46	0.46
1:A:144:PHE:CD1	1:A:145:GLU:HG3	2.50	0.46
1:A:83:LEU:O	1:A:84:LEU:C	2.53	0.46
1:A:17:TRP:CD1	1:A:56:PRO:HD2	2.50	0.46
1:A:55:SER:CB	1:A:56:PRO:CD	2.93	0.46
1:A:151:VAL:HG23	1:A:152:TRP:N	2.22	0.46
1:A:62:PHE:O	1:A:63:PRO:C	2.53	0.46
1:A:126:ASN:O	1:A:130:VAL:HG23	2.16	0.46
1:A:75:VAL:HG23	1:A:76:LEU:N	2.30	0.45
1:A:36:GLN:HE22	1:A:115:ASP:CB	2.30	0.45
1:A:138:SER:O	1:A:139:TYR:CD1	2.70	0.45
1:A:63:PRO:O	1:A:141:ARG:HB2	2.16	0.44
1:A:58:VAL:O	1:A:58:VAL:CG1	2.66	0.44
1:A:107:THR:C	1:A:109:ASP:H	2.20	0.44
1:A:3:SER:HA	1:A:156:PRO:HB2	2.00	0.44
1:A:29:ASN:OD1	1:A:30:ALA:N	2.50	0.44
1:A:80:VAL:HG22	1:A:128:LEU:HD12	1.99	0.44
1:A:79:LEU:HD21	1:A:127:ASN:CB	2.48	0.44
1:A:4:ILE:HG21	1:A:10:PHE:CE1	2.53	0.44
1:A:38:GLN:OE1	1:A:38:GLN:HA	2.19	0.43
1:A:17:TRP:CD1	1:A:17:TRP:N	2.81	0.43
1:A:9:GLN:O	1:A:12:PHE:HB2	2.18	0.43
1:A:10:PHE:C	1:A:12:PHE:H	2.22	0.43
1:A:31:LEU:HD23	1:A:31:LEU:HA	1.80	0.43
1:A:83:LEU:HD11	1:A:87:PHE:CE2	2.53	0.43
1:A:91:ASN:ND2	1:A:91:ASN:N	2.63	0.43
1:A:152:TRP:HE3	1:A:154:SER:N	2.13	0.43
1:A:76:LEU:HD23	1:A:76:LEU:C	2.37	0.43
1:A:112:ARG:HD3	1:A:116:ASP:OD1	2.18	0.43
1:A:36:GLN:HE21	1:A:118:THR:HG21	1.82	0.42
1:A:3:SER:HA	1:A:156:PRO:HD2	2.00	0.42
1:A:93:ILE:HG22	1:A:94:ILE:N	2.33	0.42
1:A:10:PHE:O	1:A:12:PHE:N	2.52	0.42

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:62:PHE:O	1:A:63:PRO:O	2.37	0.42
1:A:69:VAL:O	1:A:70:TYR:C	2.58	0.42
1:A:20:PRO:HG2	1:A:21:ILE:N	2.34	0.42
1:A:78:PRO:CA	1:A:81:THR:HG22	2.41	0.42
1:A:10:PHE:C	1:A:12:PHE:N	2.72	0.42
1:A:134:ARG:HH11	1:A:134:ARG:CG	2.26	0.42
1:A:62:PHE:CD1	1:A:63:PRO:HD2	2.55	0.42
1:A:77:ASP:N	1:A:78:PRO:HD2	2.35	0.42
1:A:79:LEU:O	1:A:79:LEU:HD23	2.18	0.42
1:A:60:VAL:O	1:A:61:ARG:O	2.38	0.42
1:A:103:THR:CG2	1:A:104:THR:N	2.82	0.42
1:A:92:ARG:CZ	1:A:96:VAL:HG21	2.50	0.42
1:A:62:PHE:HE2	1:A:139:TYR:O	2.03	0.41
1:A:45:GLN:CB	1:A:87:PHE:CD2	2.97	0.41
1:A:21:ILE:HG23	1:A:22:GLU:H	1.84	0.41
1:A:77:ASP:CG	1:A:78:PRO:N	2.71	0.41
1:A:3:SER:CA	1:A:156:PRO:HG2	2.42	0.41
1:A:67:PHE:N	1:A:67:PHE:CD1	2.89	0.41
1:A:18:ALA:CB	1:A:52:TRP:CZ3	3.03	0.41
1:A:77:ASP:CB	1:A:78:PRO:CD	2.98	0.41
1:A:108:LEU:HA	1:A:111:THR:HG22	2.01	0.41
1:A:4:ILE:H	1:A:156:PRO:HG2	1.85	0.41
1:A:36:GLN:HE22	1:A:115:ASP:HB3	1.85	0.41
1:A:134:ARG:CG	1:A:134:ARG:NH1	2.85	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	154/159 (97%)	101 (66%)	32 (21%)	21 (14%)	0 2

All (21) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	66	ASP
1	A	68	LYS
1	A	74	ALA
1	A	83	LEU
1	A	92	ARG
1	A	101	ASN
1	A	122	ARG
1	A	38	GLN
1	A	55	SER
1	A	61	ARG
1	A	108	LEU
1	A	63	PRO
1	A	102	PRO
1	A	103	THR
1	A	84	LEU
1	A	121	ILE
1	A	34	GLN
1	A	85	GLY
1	A	146	SER
1	A	96	VAL
1	A	134	ARG

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	136/139 (98%)	116 (85%)	20 (15%)	3 17

All (20) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	9	GLN
1	A	13	LEU
1	A	19	ASP
1	A	23	LEU

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	48	PHE
1	A	63	PRO
1	A	71	ARG
1	A	75	VAL
1	A	77	ASP
1	A	78	PRO
1	A	79	LEU
1	A	91	ASN
1	A	93	ILE
1	A	94	ILE
1	A	103	THR
1	A	112	ARG
1	A	122	ARG
1	A	128	LEU
1	A	133	ILE
1	A	152	TRP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	36	GLN
1	A	57	GLN
1	A	73	ASN
1	A	91	ASN

5.3.3 RNA ⓘ

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
2	R	2/3 (66%)	1 (50%)	0

All (1) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
2	R	5	U

There are no RNA pucker outliers to report.

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.