



Full wwPDB EM Validation Report ⓘ

Mar 7, 2026 – 03:03 AM UTC

PDB ID : 1SOR / pdb_00001sor
Title : Aquaporin-0 membrane junctions reveal the structure of a closed water pore
Authors : Gonen, T.; Sliz, P.; Kistler, J.; Cheng, Y.; Walz, T.
Deposited on : 2004-03-15
Resolution : 3.00 Å(reported)

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)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

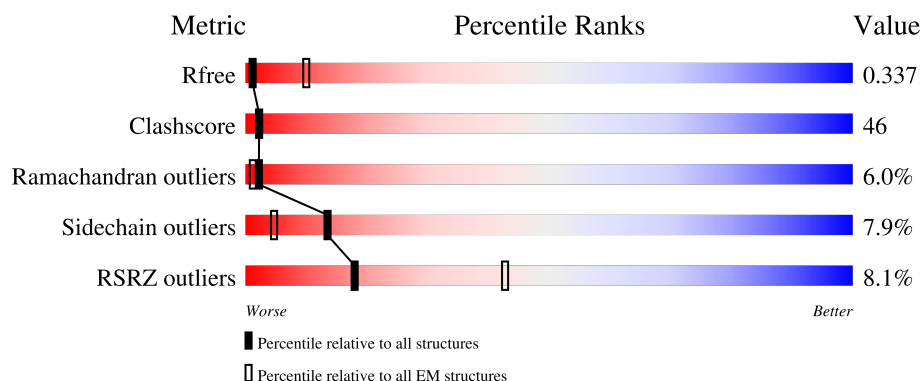
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON CRYSTALLOGRAPHY

The reported resolution of this entry is 3.00 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)
R_{free}	180332	208
Clashscore	229148	23984
Ramachandran outliers	224038	23583
Sidechain outliers	223484	23102
RSRZ outliers	180361	209

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	235	8% 38% 53% 8%

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 1778 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.

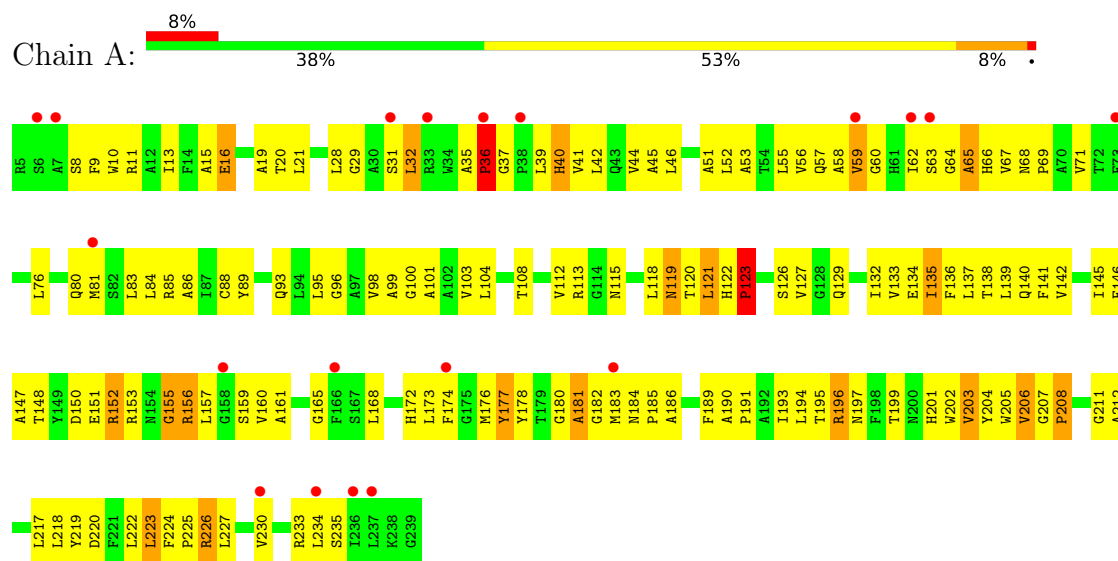
- Molecule 1 is a protein called Aquaporin-0.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
1	A	235	1778	1179	299	295	5	0	0

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. 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: Aquaporin-0



4 Data and refinement statistics

Property	Value	Source
Space group	P 4 2 2	Depositor
Cell constants a, b, c, α , β , γ	65.50Å 65.50Å 160.00Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	22.25 – 3.00 22.25 – 3.00	Depositor EDS
% Data completeness (in resolution range)	88.2 (22.25-3.00) 87.8 (22.25-3.00)	Depositor EDS
R_{merge}	0.16	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.03 (at 2.89Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.299 , 0.338 0.298 , 0.337	Depositor DCC
R_{free} test set	730 reflections (10.53%)	wwPDB-VP
Wilson B-factor (Å ²)	81.5	Xtriage
Anisotropy	0.456	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 84.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.41$, $\langle L^2 \rangle = 0.24$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.89	EDS
Total number of atoms	1778	wwPDB-VP
Average B, all atoms (Å ²)	81.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 13.26% 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	0.36	0/1828	0.90	5/2495 (0.2%)

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	59	VAL	N-CA-C	9.12	119.10	110.53
1	A	178	TYR	N-CA-C	5.69	117.29	111.14
1	A	115	ASN	N-CA-C	-5.64	105.08	112.41
1	A	177	TYR	N-CA-C	5.35	116.92	111.14
1	A	206	VAL	N-CA-C	5.31	115.52	110.42

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	1778	0	1807	165	0
All	All	1778	0	1807	165	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 46.

All (165) 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:19:ALA:HB2	1:A:59:VAL:HG21	1.44	1.00
1:A:98:VAL:HG13	1:A:194:LEU:HD21	1.49	0.95
1:A:204:TYR:O	1:A:208:PRO:HD2	1.72	0.89
1:A:135:ILE:HD12	1:A:207:GLY:HA2	1.55	0.86
1:A:226:ARG:HH11	1:A:233:ARG:HH12	1.27	0.82
1:A:182:GLY:O	1:A:183:MET:HG2	1.81	0.79
1:A:152:ARG:HB2	1:A:152:ARG:NH1	1.99	0.78
1:A:81:MET:HE3	1:A:86:ALA:HB2	1.65	0.77
1:A:28:LEU:HD21	1:A:191:PRO:HG3	1.65	0.77
1:A:172:HIS:HA	1:A:183:MET:HE1	1.67	0.75
1:A:226:ARG:NH1	1:A:233:ARG:HH12	1.84	0.75
1:A:196:ARG:H	1:A:196:ARG:HD2	1.51	0.75
1:A:57:GLN:HE21	1:A:161:ALA:HB3	1.52	0.74
1:A:51:ALA:O	1:A:55:LEU:HD23	1.88	0.74
1:A:84:LEU:HD12	1:A:85:ARG:N	2.04	0.73
1:A:41:VAL:HG13	1:A:42:LEU:HD12	1.71	0.72
1:A:101:ALA:HB2	1:A:190:ALA:HB1	1.74	0.70
1:A:220:ASP:HA	1:A:226:ARG:NH2	2.07	0.70
1:A:226:ARG:HH11	1:A:226:ARG:HB2	1.57	0.70
1:A:155:GLY:C	1:A:156:ARG:HD3	2.18	0.68
1:A:98:VAL:HG13	1:A:194:LEU:CD2	2.23	0.68
1:A:225:PRO:HA	1:A:226:ARG:NH2	2.09	0.67
1:A:57:GLN:HE21	1:A:161:ALA:CB	2.06	0.67
1:A:13:ILE:HD12	1:A:88:CYS:HB3	1.76	0.67
1:A:52:LEU:O	1:A:56:VAL:HG23	1.95	0.66
1:A:220:ASP:HA	1:A:226:ARG:HH22	1.60	0.66
1:A:138:THR:HG21	1:A:208:PRO:HA	1.77	0.66
1:A:135:ILE:CD1	1:A:207:GLY:HA2	2.25	0.66
1:A:40:HIS:O	1:A:44:VAL:HG12	1.95	0.66
1:A:152:ARG:HB2	1:A:152:ARG:HH11	1.60	0.65
1:A:151:GLU:HG3	1:A:152:ARG:N	2.12	0.65
1:A:157:LEU:HD12	1:A:157:LEU:H	1.61	0.65
1:A:134:GLU:HB3	1:A:208:PRO:HD3	1.78	0.64
1:A:53:ALA:HB2	1:A:165:GLY:HA3	1.80	0.64
1:A:35:ALA:O	1:A:37:GLY:N	2.31	0.64
1:A:148:THR:HG23	1:A:160:VAL:HB	1.80	0.63
1:A:225:PRO:HA	1:A:226:ARG:HH22	1.62	0.63
1:A:19:ALA:CB	1:A:59:VAL:HG21	2.23	0.63
1:A:226:ARG:HH11	1:A:233:ARG:NH1	1.94	0.63
1:A:8:SER:O	1:A:11:ARG:HG2	1.99	0.63
1:A:104:LEU:O	1:A:108:THR:HG23	1.99	0.62
1:A:156:ARG:HD3	1:A:156:ARG:N	2.14	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:35:ALA:N	1:A:36:PRO:CD	2.62	0.62
1:A:195:THR:HG22	1:A:195:THR:O	2.00	0.61
1:A:113:ARG:HD2	1:A:113:ARG:C	2.26	0.61
1:A:122:HIS:CD2	1:A:123:PRO:HD2	2.34	0.61
1:A:135:ILE:HD12	1:A:207:GLY:CA	2.29	0.61
1:A:204:TYR:O	1:A:208:PRO:CD	2.47	0.60
1:A:135:ILE:HG22	1:A:136:PHE:HD1	1.66	0.60
1:A:148:THR:CG2	1:A:160:VAL:HB	2.32	0.60
1:A:226:ARG:HD2	1:A:233:ARG:NH1	2.16	0.60
1:A:64:GLY:O	1:A:65:ALA:HB3	2.02	0.59
1:A:153:ARG:HG2	1:A:157:LEU:HD11	1.84	0.59
1:A:135:ILE:HG22	1:A:136:PHE:CD1	2.37	0.59
1:A:129:GLN:O	1:A:133:VAL:HG23	2.03	0.58
1:A:133:VAL:O	1:A:137:LEU:HG	2.04	0.58
1:A:190:ALA:HB3	1:A:191:PRO:HD3	1.85	0.58
1:A:226:ARG:NH1	1:A:226:ARG:HB2	2.18	0.58
1:A:195:THR:O	1:A:196:ARG:C	2.47	0.57
1:A:230:VAL:O	1:A:233:ARG:HG2	2.05	0.57
1:A:52:LEU:HD22	1:A:168:LEU:HG	1.87	0.57
1:A:60:GLY:N	1:A:65:ALA:HB2	2.21	0.56
1:A:126:SER:OG	1:A:129:GLN:HG3	2.06	0.56
1:A:185:PRO:HB3	1:A:205:TRP:CE2	2.41	0.56
1:A:101:ALA:CB	1:A:190:ALA:HB1	2.36	0.55
1:A:118:LEU:O	1:A:120:THR:HG23	2.06	0.55
1:A:15:ALA:HB1	1:A:59:VAL:HG13	1.89	0.55
1:A:141:PHE:CZ	1:A:145:ILE:HD11	2.42	0.54
1:A:233:ARG:C	1:A:235:SER:H	2.15	0.54
1:A:21:LEU:HA	1:A:96:GLY:O	2.07	0.54
1:A:83:LEU:O	1:A:83:LEU:HD23	2.08	0.54
1:A:65:ALA:O	1:A:67:VAL:HG23	2.08	0.54
1:A:168:LEU:HD11	1:A:172:HIS:CE1	2.43	0.53
1:A:152:ARG:HG2	1:A:224:PHE:CZ	2.44	0.53
1:A:31:SER:O	1:A:32:LEU:C	2.51	0.53
1:A:157:LEU:HD12	1:A:157:LEU:N	2.24	0.53
1:A:122:HIS:HD2	1:A:123:PRO:HD2	1.73	0.52
1:A:100:GLY:O	1:A:103:VAL:HG22	2.09	0.52
1:A:202:TRP:O	1:A:206:VAL:HG23	2.09	0.52
1:A:176:MET:HA	1:A:181:ALA:N	2.24	0.52
1:A:62:ILE:HD12	1:A:62:ILE:N	2.25	0.52
1:A:35:ALA:CB	1:A:40:HIS:HB3	2.40	0.52
1:A:151:GLU:HG3	1:A:152:ARG:H	1.74	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:157:LEU:H	1:A:157:LEU:CD1	2.23	0.52
1:A:173:LEU:C	1:A:173:LEU:HD23	2.35	0.51
1:A:138:THR:HG21	1:A:208:PRO:CA	2.41	0.50
1:A:152:ARG:HG2	1:A:224:PHE:HZ	1.76	0.50
1:A:227:LEU:HD12	1:A:227:LEU:N	2.26	0.50
1:A:60:GLY:CA	1:A:65:ALA:HB2	2.42	0.50
1:A:184:ASN:OD1	1:A:186:ALA:HB3	2.12	0.50
1:A:199:THR:C	1:A:201:HIS:H	2.19	0.50
1:A:189:PHE:O	1:A:193:ILE:HG22	2.12	0.49
1:A:217:LEU:C	1:A:217:LEU:HD23	2.38	0.49
1:A:46:LEU:C	1:A:46:LEU:HD13	2.37	0.49
1:A:135:ILE:HG22	1:A:136:PHE:N	2.27	0.49
1:A:137:LEU:O	1:A:140:GLN:HG3	2.12	0.49
1:A:122:HIS:HD2	1:A:123:PRO:CD	2.26	0.49
1:A:196:ARG:HD2	1:A:196:ARG:N	2.25	0.49
1:A:137:LEU:HD21	1:A:174:PHE:HD2	1.78	0.48
1:A:32:LEU:HD12	1:A:112:VAL:HG12	1.96	0.48
1:A:64:GLY:O	1:A:65:ALA:CB	2.61	0.48
1:A:202:TRP:CZ2	1:A:203:VAL:HG12	2.49	0.48
1:A:222:LEU:O	1:A:225:PRO:HD3	2.14	0.48
1:A:29:GLY:HA2	1:A:32:LEU:HD23	1.95	0.47
1:A:29:GLY:O	1:A:32:LEU:HD23	2.13	0.47
1:A:68:ASN:HB3	1:A:71:VAL:CG2	2.44	0.47
1:A:219:TYR:HA	1:A:223:LEU:HD12	1.96	0.47
1:A:35:ALA:O	1:A:36:PRO:C	2.58	0.47
1:A:147:ALA:HB1	1:A:153:ARG:NH2	2.30	0.47
1:A:152:ARG:HH11	1:A:152:ARG:CB	2.26	0.47
1:A:199:THR:C	1:A:201:HIS:N	2.73	0.46
1:A:71:VAL:HG21	1:A:141:PHE:CE2	2.50	0.46
1:A:156:ARG:N	1:A:156:ARG:CD	2.78	0.46
1:A:226:ARG:NH1	1:A:226:ARG:N	2.64	0.46
1:A:41:VAL:HG11	1:A:177:TYR:OH	2.15	0.46
1:A:226:ARG:HD2	1:A:233:ARG:HH11	1.81	0.46
1:A:69:PRO:HA	1:A:93:GLN:HE22	1.80	0.45
1:A:142:VAL:HG21	1:A:212:ALA:HA	1.98	0.45
1:A:176:MET:HE3	1:A:180:GLY:HA2	1.97	0.45
1:A:89:TYR:HB3	1:A:93:GLN:OE1	2.16	0.45
1:A:226:ARG:NH1	1:A:226:ARG:H	2.15	0.45
1:A:84:LEU:HD12	1:A:84:LEU:C	2.41	0.45
1:A:52:LEU:HD21	1:A:141:PHE:CZ	2.52	0.45
1:A:139:LEU:HD23	1:A:139:LEU:O	2.17	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:119:ASN:ND2	1:A:181:ALA:O	2.49	0.44
1:A:121:LEU:HD13	1:A:122:HIS:N	2.32	0.44
1:A:39:LEU:HG	1:A:41:VAL:HG12	1.98	0.44
1:A:71:VAL:HG21	1:A:141:PHE:HE2	1.83	0.44
1:A:39:LEU:HD12	1:A:39:LEU:HA	1.86	0.44
1:A:19:ALA:HA	1:A:55:LEU:CD1	2.48	0.43
1:A:151:GLU:O	1:A:152:ARG:C	2.61	0.43
1:A:138:THR:HG21	1:A:208:PRO:CB	2.48	0.43
1:A:134:GLU:OE1	1:A:137:LEU:HD12	2.17	0.43
1:A:9:PHE:CD1	1:A:9:PHE:C	2.97	0.43
1:A:222:LEU:O	1:A:223:LEU:C	2.61	0.43
1:A:39:LEU:C	1:A:41:VAL:H	2.26	0.42
1:A:44:VAL:HG13	1:A:45:ALA:N	2.34	0.42
1:A:58:ALA:O	1:A:62:ILE:HD13	2.19	0.42
1:A:16:GLU:OE1	1:A:66:HIS:N	2.53	0.42
1:A:19:ALA:HA	1:A:55:LEU:HD12	2.00	0.42
1:A:139:LEU:HD23	1:A:139:LEU:C	2.44	0.42
1:A:19:ALA:HB2	1:A:59:VAL:CG2	2.31	0.42
1:A:13:ILE:CD1	1:A:88:CYS:HB3	2.46	0.42
1:A:152:ARG:NH2	1:A:226:ARG:HG3	2.34	0.42
1:A:113:ARG:HD2	1:A:113:ARG:O	2.18	0.42
1:A:21:LEU:C	1:A:21:LEU:HD23	2.45	0.42
1:A:39:LEU:C	1:A:41:VAL:N	2.78	0.42
1:A:227:LEU:HD12	1:A:227:LEU:H	1.83	0.42
1:A:197:ASN:HD22	1:A:197:ASN:HA	1.61	0.42
1:A:20:THR:HG21	1:A:93:GLN:O	2.20	0.42
1:A:152:ARG:HH21	1:A:226:ARG:HG3	1.85	0.42
1:A:127:VAL:HG11	1:A:202:TRP:CZ2	2.55	0.42
1:A:138:THR:OG1	1:A:211:GLY:HA3	2.20	0.41
1:A:233:ARG:C	1:A:235:SER:N	2.78	0.41
1:A:62:ILE:O	1:A:63:SER:HB2	2.19	0.41
1:A:122:HIS:O	1:A:123:PRO:C	2.63	0.41
1:A:140:GLN:NE2	1:A:141:PHE:HB2	2.36	0.41
1:A:46:LEU:HD13	1:A:46:LEU:O	2.20	0.41
1:A:193:ILE:C	1:A:195:THR:H	2.27	0.41
1:A:76:LEU:HD22	1:A:86:ALA:CB	2.49	0.41
1:A:138:THR:HG21	1:A:208:PRO:HB3	2.01	0.41
1:A:76:LEU:HB2	1:A:81:MET:CE	2.51	0.41
1:A:99:ALA:O	1:A:103:VAL:HG13	2.21	0.40
1:A:218:LEU:HD12	1:A:222:LEU:HD12	2.03	0.40
1:A:39:LEU:O	1:A:41:VAL:N	2.54	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	233/235 (99%)	193 (83%)	26 (11%)	14 (6%)	1 7

All (14) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	36	PRO
1	A	80	GLN
1	A	123	PRO
1	A	196	ARG
1	A	223	LEU
1	A	119	ASN
1	A	155	GLY
1	A	181	ALA
1	A	32	LEU
1	A	40	HIS
1	A	150	ASP
1	A	159	SER
1	A	65	ALA
1	A	234	LEU

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	178/180 (99%)	164 (92%)	14 (8%)	11	39

All (14) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	10	TRP
1	A	16	GLU
1	A	36	PRO
1	A	95	LEU
1	A	121	LEU
1	A	123	PRO
1	A	132	ILE
1	A	135	ILE
1	A	146	PHE
1	A	152	ARG
1	A	156	ARG
1	A	203	VAL
1	A	208	PRO
1	A	226	ARG

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

Mol	Chain	Res	Type
1	A	43	GLN
1	A	57	GLN
1	A	115	ASN
1	A	129	GLN
1	A	140	GLN
1	A	197	ASN

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 [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 [i](#)

There are no chain breaks in this entry.