



Full wwPDB X-ray Structure Validation Report ⓘ

Oct 23, 2024 – 08:02 AM EDT

PDB ID : 2ZF8
Title : Crystal structure of MotY
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Deposited on : 2007-12-25
Resolution : 2.85 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 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	:	4.02b-467
Xtriage (Phenix)	:	1.20.1
EDS	:	3.0
Percentile statistics	:	20231227.v01 (using entries in the PDB archive December 27th 2023)
CCP4	:	9.0.003 (Gargrove)
Density-Fitness	:	1.0.11
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.39

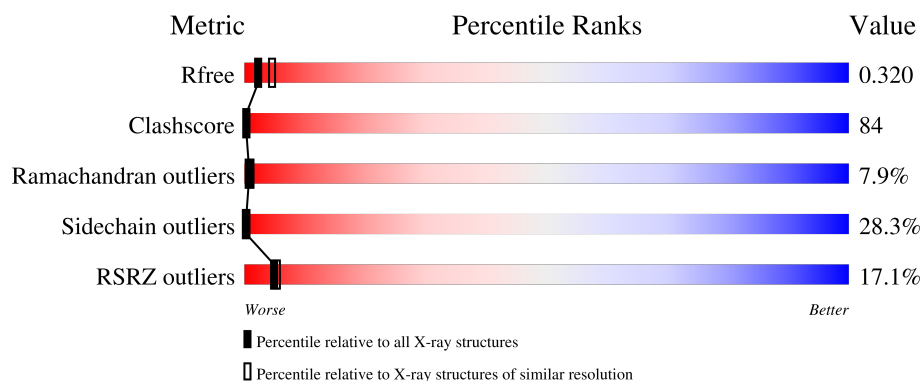
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.85 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}	164625	1268 (2.88-2.84)
Clashscore	180529	1351 (2.88-2.84)
Ramachandran outliers	177936	1318 (2.88-2.84)
Sidechain outliers	177891	1319 (2.88-2.84)
RSRZ outliers	164620	1269 (2.88-2.84)

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 of 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	278	<div> <div>15%</div> <div>23%</div> <div>42%</div> <div>20%</div> <div>•</div> <div>12%</div> </div>

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 2012 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 Component of sodium-driven polar flagellar motor.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	246	Total	C	N	O	S	0	0	0
			2000	1266	355	372	7			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	273	HIS	-	expression tag	UNP P74944
A	274	HIS	-	expression tag	UNP P74944
A	275	HIS	-	expression tag	UNP P74944
A	276	HIS	-	expression tag	UNP P74944
A	277	HIS	-	expression tag	UNP P74944
A	278	HIS	-	expression tag	UNP P74944

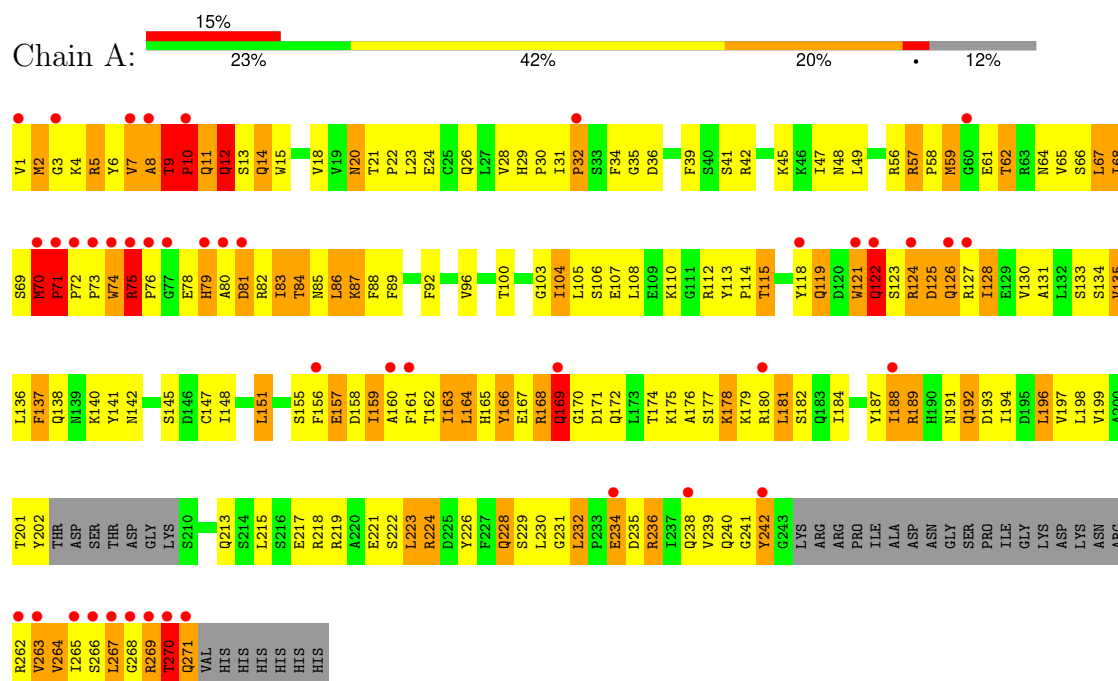
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	12	Total	O	0	0
			12	12		

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 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: Component of sodium-driven polar flagellar motor



4 Data and refinement statistics

Property	Value	Source
Space group	P 65 2 2	Depositor
Cell constants a, b, c, α , β , γ	104.08Å 104.08Å 133.10Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	41.00 – 2.85 41.00 – 2.85	Depositor EDS
% Data completeness (in resolution range)	92.5 (41.00-2.85) 99.4 (41.00-2.85)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.68 (at 2.86Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.290 , 0.312 0.300 , 0.320	Depositor DCC
R_{free} test set	1027 reflections (9.86%)	wwPDB-VP
Wilson B-factor (Å ²)	65.8	Xtriage
Anisotropy	0.631	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 51.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.88	EDS
Total number of atoms	2012	wwPDB-VP
Average B, all atoms (Å ²)	73.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.92% 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	0.59	0/2043	1.01	13/2757 (0.5%)

There are no bond length outliers.

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	59	MET	N-CA-C	-9.46	85.47	111.00
1	A	8	ALA	CA-C-N	-8.58	98.33	117.20
1	A	267	LEU	CA-CB-CG	-8.36	96.08	115.30
1	A	7	VAL	C-N-CA	-7.59	102.71	121.70
1	A	71	PRO	N-CA-C	6.62	129.32	112.10
1	A	12	GLN	N-CA-C	6.27	127.92	111.00
1	A	10	PRO	N-CA-C	6.26	128.38	112.10
1	A	263	VAL	N-CA-C	-6.00	94.81	111.00
1	A	57	ARG	N-CA-C	-5.60	95.88	111.00
1	A	8	ALA	O-C-N	5.51	131.51	122.70
1	A	8	ALA	N-CA-C	5.46	125.73	111.00
1	A	70	MET	N-CA-C	5.28	125.25	111.00
1	A	9	THR	N-CA-C	-5.05	97.38	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	2000	0	1971	334	2
2	A	12	0	0	4	0
All	All	2012	0	1971	334	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 84.

All (334) 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:194:ILE:HA	1:A:270:THR:OG1	1.29	1.28
1:A:11:GLN:CG	1:A:12:GLN:H	1.44	1.24
1:A:8:ALA:HB2	1:A:32:PRO:CG	1.71	1.20
1:A:13:SER:OG	1:A:135:VAL:HG21	1.46	1.15
1:A:270:THR:CG2	1:A:271:GLN:HG2	1.77	1.15
1:A:11:GLN:HG3	1:A:12:GLN:N	1.35	1.14
1:A:75:ARG:HB3	1:A:76:PRO:HD2	1.25	1.14
1:A:8:ALA:CB	1:A:32:PRO:HG3	1.77	1.12
1:A:179:LYS:HB3	1:A:180:ARG:HH11	1.06	1.11
1:A:69:SER:O	1:A:70:MET:HG3	1.52	1.09
1:A:170:GLY:HA2	1:A:219:ARG:NH1	1.66	1.09
1:A:270:THR:HG22	1:A:271:GLN:N	1.52	1.08
1:A:135:VAL:HG23	1:A:136:LEU:H	1.05	1.08
1:A:135:VAL:HG23	1:A:136:LEU:N	1.61	1.08
1:A:135:VAL:CG2	1:A:136:LEU:H	1.65	1.07
1:A:74:TRP:HA	1:A:75:ARG:HH11	0.93	1.07
1:A:75:ARG:HB3	1:A:76:PRO:CD	1.85	1.07
1:A:270:THR:HG21	1:A:271:GLN:HG2	1.34	1.06
1:A:13:SER:OG	1:A:135:VAL:CG2	2.05	1.05
1:A:31:ILE:HG21	1:A:34:PHE:HB3	1.38	1.05
1:A:11:GLN:HE21	1:A:12:GLN:HB3	1.16	1.04
1:A:8:ALA:HB2	1:A:32:PRO:HG3	1.09	1.02
1:A:74:TRP:HA	1:A:75:ARG:NH1	1.74	1.00
1:A:135:VAL:HG23	1:A:136:LEU:HG	1.40	0.99
1:A:133:SER:OG	1:A:135:VAL:HG13	1.62	0.99
1:A:6:TYR:OH	1:A:57:ARG:NH2	1.98	0.95
1:A:4:LYS:HD2	1:A:121:TRP:HE1	1.29	0.94
1:A:79:HIS:HD2	1:A:80:ALA:N	1.63	0.94
1:A:13:SER:HG	1:A:135:VAL:HG21	1.14	0.94
1:A:179:LYS:HB3	1:A:180:ARG:NH1	1.82	0.94
1:A:70:MET:HA	1:A:112:ARG:HH21	1.32	0.93
1:A:194:ILE:CA	1:A:270:THR:OG1	2.16	0.92

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:4:LYS:CD	1:A:121:TRP:HE1	1.82	0.92
1:A:75:ARG:N	1:A:75:ARG:CD	2.33	0.91
1:A:75:ARG:N	1:A:75:ARG:HD2	1.85	0.89
1:A:31:ILE:CG2	1:A:34:PHE:HB3	2.02	0.89
1:A:11:GLN:NE2	1:A:12:GLN:HB3	1.88	0.88
1:A:174:THR:O	1:A:178:LYS:HE2	1.74	0.88
1:A:184:ILE:O	1:A:188:ILE:HG12	1.73	0.88
1:A:4:LYS:CE	1:A:121:TRP:HE1	1.87	0.88
1:A:15:TRP:O	1:A:140:LYS:HE3	1.74	0.87
1:A:270:THR:CG2	1:A:271:GLN:N	2.17	0.86
1:A:20:ASN:ND2	1:A:147:CYS:SG	2.48	0.86
1:A:69:SER:C	1:A:70:MET:HG3	1.95	0.85
1:A:170:GLY:HA2	1:A:219:ARG:HH12	1.40	0.85
1:A:270:THR:HG22	1:A:271:GLN:CA	2.07	0.85
1:A:270:THR:HG22	1:A:271:GLN:HG2	1.55	0.84
1:A:176:ALA:O	1:A:180:ARG:HD3	1.79	0.83
1:A:79:HIS:CD2	1:A:80:ALA:N	2.46	0.82
1:A:170:GLY:CA	1:A:219:ARG:NH1	2.42	0.82
1:A:179:LYS:HD3	1:A:179:LYS:O	1.80	0.82
1:A:189:ARG:HB2	1:A:189:ARG:HH11	1.44	0.82
1:A:75:ARG:CB	1:A:76:PRO:HD2	2.08	0.82
1:A:8:ALA:CA	1:A:32:PRO:HG3	2.10	0.81
1:A:9:THR:HG22	1:A:12:GLN:HG2	1.62	0.81
1:A:70:MET:O	1:A:72:PRO:HD3	1.80	0.81
1:A:270:THR:CG2	1:A:271:GLN:CG	2.57	0.81
1:A:9:THR:CG2	1:A:12:GLN:HG2	2.10	0.81
1:A:194:ILE:O	1:A:236:ARG:NH1	2.15	0.79
1:A:13:SER:OG	1:A:135:VAL:HG11	1.82	0.79
1:A:45:LYS:NZ	1:A:157:GLU:OE1	2.15	0.78
1:A:20:ASN:HD21	1:A:147:CYS:HB2	1.49	0.78
1:A:270:THR:HG22	1:A:271:GLN:CG	2.13	0.78
1:A:8:ALA:HB2	1:A:32:PRO:CD	2.13	0.77
1:A:31:ILE:HG21	1:A:34:PHE:CB	2.14	0.76
1:A:213:GLN:O	1:A:217:GLU:HG3	1.86	0.76
1:A:121:TRP:HZ3	1:A:125:ASP:HB2	1.50	0.76
1:A:13:SER:OG	1:A:135:VAL:CG1	2.33	0.76
1:A:160:ALA:HB1	2:A:288:HOH:O	1.86	0.76
1:A:170:GLY:HA3	1:A:215:LEU:HD11	1.67	0.75
1:A:4:LYS:HD2	1:A:121:TRP:NE1	2.02	0.75
1:A:61:GLU:HA	2:A:283:HOH:O	1.85	0.75
1:A:128:ILE:HG22	2:A:279:HOH:O	1.88	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:135:VAL:CG2	1:A:136:LEU:HG	2.15	0.74
1:A:189:ARG:HH11	1:A:189:ARG:CB	2.02	0.73
1:A:5:ARG:HG3	1:A:7:VAL:CG2	2.19	0.73
1:A:223:LEU:N	1:A:223:LEU:HD23	2.03	0.73
1:A:62:THR:HG21	1:A:118:TYR:HE1	1.54	0.73
1:A:1:VAL:O	1:A:126:GLN:HA	1.88	0.72
1:A:6:TYR:CZ	1:A:57:ARG:NH2	2.52	0.72
1:A:83:ILE:O	1:A:83:ILE:HG22	1.88	0.72
1:A:121:TRP:CZ3	1:A:125:ASP:OD2	2.43	0.71
1:A:78:GLU:HG3	1:A:81:ASP:H	1.54	0.71
1:A:135:VAL:CG2	1:A:136:LEU:N	2.30	0.71
1:A:170:GLY:C	1:A:219:ARG:HH11	1.94	0.71
1:A:21:THR:HG23	1:A:22:PRO:HD2	1.72	0.70
1:A:24:GLU:HG3	1:A:42:ARG:HG2	1.74	0.69
1:A:177:SER:O	1:A:181:LEU:HD22	1.91	0.69
1:A:169:GLN:OE1	1:A:169:GLN:HA	1.92	0.69
1:A:133:SER:OG	1:A:135:VAL:CG1	2.41	0.69
1:A:189:ARG:HH11	1:A:189:ARG:CG	2.05	0.69
1:A:20:ASN:HD21	1:A:147:CYS:CB	2.07	0.68
1:A:215:LEU:HD21	1:A:219:ARG:CZ	2.24	0.68
1:A:121:TRP:HZ3	1:A:125:ASP:CB	2.06	0.68
1:A:79:HIS:O	1:A:83:ILE:HG13	1.94	0.68
1:A:179:LYS:CB	1:A:180:ARG:HH11	1.97	0.67
1:A:174:THR:HG22	1:A:175:LYS:N	2.09	0.67
1:A:70:MET:SD	1:A:115:THR:HG21	2.34	0.67
1:A:170:GLY:CA	1:A:219:ARG:HH11	2.08	0.67
1:A:47:ILE:HD12	1:A:47:ILE:H	1.60	0.67
1:A:5:ARG:CG	1:A:7:VAL:CG2	2.73	0.67
1:A:135:VAL:HG23	1:A:136:LEU:CG	2.20	0.67
1:A:128:ILE:HG12	1:A:128:ILE:O	1.93	0.67
1:A:5:ARG:CG	1:A:7:VAL:HG21	2.25	0.66
1:A:31:ILE:N	1:A:35:GLY:O	2.23	0.66
1:A:124:ARG:HD2	1:A:124:ARG:O	1.95	0.66
1:A:21:THR:HG22	1:A:23:LEU:N	2.11	0.66
1:A:70:MET:HA	1:A:112:ARG:NH2	2.08	0.66
1:A:196:LEU:HA	1:A:267:LEU:HD22	1.78	0.65
1:A:148:ILE:HA	1:A:151:LEU:HD22	1.78	0.65
1:A:119:GLN:NE2	1:A:123:SER:O	2.30	0.65
1:A:158:ASP:HA	1:A:161:PHE:CE1	2.31	0.65
1:A:75:ARG:H	1:A:75:ARG:HD3	1.62	0.64
1:A:5:ARG:HG3	1:A:7:VAL:HG23	1.79	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:4:LYS:CE	1:A:121:TRP:NE1	2.59	0.64
1:A:268:GLY:O	1:A:269:ARG:C	2.36	0.64
1:A:11:GLN:HG3	1:A:12:GLN:H	0.57	0.64
1:A:68:ILE:HG22	1:A:84:THR:O	1.98	0.64
1:A:162:THR:OG1	1:A:264:VAL:CG1	2.46	0.64
1:A:9:THR:HG22	1:A:12:GLN:CG	2.29	0.63
1:A:167:GLU:C	1:A:169:GLN:H	2.02	0.63
1:A:180:ARG:N	1:A:180:ARG:CD	2.61	0.63
1:A:31:ILE:HG22	1:A:34:PHE:H	1.63	0.63
1:A:121:TRP:HE3	1:A:121:TRP:O	1.82	0.63
1:A:148:ILE:O	1:A:151:LEU:HB2	1.98	0.63
1:A:270:THR:HG22	1:A:271:GLN:CB	2.28	0.63
1:A:11:GLN:CG	1:A:12:GLN:N	2.16	0.63
1:A:10:PRO:HB3	1:A:73:PRO:HB3	1.81	0.62
1:A:15:TRP:CE3	1:A:29:HIS:HB2	2.34	0.62
1:A:226:TYR:O	1:A:230:LEU:HG	1.99	0.62
1:A:6:TYR:C	1:A:7:VAL:HG23	2.20	0.61
1:A:201:THR:O	1:A:202:TYR:HB2	2.01	0.60
1:A:224:ARG:O	1:A:228:GLN:HB2	2.01	0.60
1:A:21:THR:C	1:A:23:LEU:H	2.04	0.60
1:A:62:THR:HG21	1:A:118:TYR:CE1	2.35	0.60
1:A:162:THR:O	1:A:164:LEU:CD1	2.49	0.60
1:A:194:ILE:HA	1:A:270:THR:HG1	1.60	0.60
1:A:80:ALA:O	1:A:81:ASP:C	2.40	0.60
1:A:189:ARG:HB2	1:A:189:ARG:NH1	2.17	0.59
1:A:69:SER:HB2	1:A:112:ARG:HB3	1.83	0.59
1:A:121:TRP:HZ3	1:A:125:ASP:CG	2.05	0.59
1:A:138:GLN:O	1:A:142:ASN:ND2	2.35	0.59
1:A:239:VAL:HG22	1:A:239:VAL:O	2.02	0.59
1:A:121:TRP:CZ3	1:A:125:ASP:HB2	2.34	0.59
1:A:192:GLN:HA	1:A:236:ARG:NH2	2.17	0.59
1:A:136:LEU:O	1:A:137:PHE:C	2.41	0.59
1:A:21:THR:HG22	1:A:23:LEU:H	1.67	0.58
1:A:74:TRP:C	1:A:74:TRP:HE3	2.07	0.58
1:A:155:SER:C	1:A:157:GLU:H	2.05	0.58
1:A:3:GLY:N	1:A:126:GLN:HB2	2.19	0.58
1:A:21:THR:HG22	1:A:22:PRO:N	2.19	0.58
1:A:69:SER:C	1:A:70:MET:CG	2.72	0.57
1:A:115:THR:HB	1:A:131:ALA:HB2	1.86	0.57
1:A:4:LYS:HE3	1:A:121:TRP:NE1	2.19	0.57
1:A:31:ILE:HD11	1:A:131:ALA:O	2.05	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:191:ASN:O	1:A:194:ILE:CG2	2.52	0.57
1:A:11:GLN:HG3	1:A:12:GLN:CA	2.29	0.57
1:A:15:TRP:CZ3	1:A:29:HIS:HB2	2.39	0.57
1:A:21:THR:CG2	1:A:23:LEU:HB2	2.35	0.57
1:A:270:THR:CG2	1:A:271:GLN:H	2.09	0.57
1:A:170:GLY:HA3	1:A:215:LEU:CD1	2.34	0.56
1:A:235:ASP:O	1:A:238:GLN:NE2	2.38	0.56
1:A:166:TYR:CE2	1:A:172:GLN:HB3	2.41	0.56
1:A:79:HIS:HD2	1:A:80:ALA:H	1.51	0.56
1:A:86:LEU:HD21	1:A:107:GLU:HG3	1.87	0.56
1:A:103:GLY:O	1:A:105:LEU:N	2.38	0.55
1:A:8:ALA:HB2	1:A:32:PRO:HD3	1.87	0.55
1:A:75:ARG:CD	1:A:75:ARG:H	2.12	0.55
1:A:68:ILE:HG22	1:A:85:ASN:HA	1.88	0.55
1:A:121:TRP:HZ3	1:A:125:ASP:OD2	1.88	0.55
1:A:168:ARG:C	1:A:170:GLY:N	2.59	0.55
1:A:166:TYR:HD2	1:A:166:TYR:O	1.90	0.55
1:A:6:TYR:C	1:A:7:VAL:CG2	2.74	0.54
1:A:176:ALA:O	1:A:179:LYS:N	2.40	0.54
1:A:8:ALA:O	1:A:9:THR:C	2.46	0.54
1:A:121:TRP:C	1:A:122:GLN:O	2.41	0.54
1:A:11:GLN:O	1:A:13:SER:N	2.41	0.54
1:A:176:ALA:HB1	1:A:180:ARG:CZ	2.39	0.54
1:A:174:THR:HG22	1:A:175:LYS:H	1.72	0.53
1:A:162:THR:O	1:A:164:LEU:HD13	2.08	0.53
1:A:189:ARG:CG	1:A:189:ARG:NH1	2.71	0.53
1:A:1:VAL:O	1:A:126:GLN:CA	2.57	0.53
1:A:168:ARG:O	1:A:170:GLY:N	2.42	0.53
1:A:75:ARG:CB	1:A:76:PRO:CD	2.65	0.52
1:A:242:TYR:HD2	1:A:242:TYR:O	1.93	0.52
1:A:21:THR:CG2	1:A:22:PRO:N	2.72	0.52
1:A:189:ARG:HH11	1:A:189:ARG:HG3	1.72	0.52
1:A:1:VAL:HB	1:A:126:GLN:NE2	2.24	0.52
1:A:162:THR:HA	1:A:264:VAL:CG1	2.40	0.52
1:A:234:GLU:C	1:A:234:GLU:OE2	2.48	0.52
1:A:74:TRP:C	1:A:75:ARG:HD2	2.30	0.51
1:A:70:MET:O	1:A:72:PRO:CD	2.54	0.51
1:A:68:ILE:O	1:A:114:PRO:HA	2.10	0.51
1:A:159:ILE:HG13	1:A:267:LEU:O	2.11	0.51
1:A:174:THR:CG2	1:A:175:LYS:N	2.73	0.51
1:A:179:LYS:HD3	1:A:179:LYS:C	2.30	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:121:TRP:CZ3	1:A:125:ASP:CG	2.84	0.51
1:A:74:TRP:HE3	1:A:74:TRP:O	1.94	0.50
1:A:9:THR:HG22	1:A:11:GLN:HG2	1.93	0.50
1:A:235:ASP:OD1	1:A:236:ARG:HG2	2.12	0.50
1:A:31:ILE:HG23	1:A:32:PRO:HD2	1.93	0.50
1:A:69:SER:O	1:A:70:MET:CG	2.42	0.50
1:A:121:TRP:CE3	1:A:125:ASP:OD2	2.65	0.50
1:A:8:ALA:HB2	1:A:32:PRO:CB	2.39	0.50
1:A:20:ASN:ND2	1:A:147:CYS:HB2	2.24	0.50
1:A:13:SER:OG	1:A:135:VAL:CB	2.59	0.50
1:A:20:ASN:ND2	1:A:147:CYS:CB	2.73	0.50
1:A:136:LEU:O	1:A:138:GLN:N	2.45	0.50
1:A:213:GLN:HG3	1:A:217:GLU:OE2	2.12	0.49
1:A:79:HIS:CD2	1:A:80:ALA:H	2.28	0.49
1:A:31:ILE:CG2	1:A:34:PHE:CB	2.81	0.49
1:A:164:LEU:O	1:A:174:THR:OG1	2.18	0.49
1:A:201:THR:O	1:A:201:THR:HG23	2.10	0.49
1:A:234:GLU:OE2	1:A:235:ASP:HB3	2.12	0.49
1:A:266:SER:O	1:A:266:SER:OG	2.26	0.49
1:A:65:VAL:HG21	1:A:92:PHE:HB3	1.94	0.49
1:A:8:ALA:HA	1:A:32:PRO:HG3	1.90	0.49
1:A:166:TYR:CE2	1:A:172:GLN:CB	2.96	0.49
1:A:242:TYR:C	1:A:242:TYR:CD2	2.86	0.49
1:A:9:THR:HG23	1:A:10:PRO:HD2	1.94	0.48
1:A:9:THR:CG2	1:A:11:GLN:HG2	2.44	0.48
1:A:122:GLN:HB2	1:A:125:ASP:OD1	2.13	0.48
1:A:179:LYS:CB	1:A:180:ARG:NH1	2.64	0.48
1:A:170:GLY:CA	1:A:215:LEU:HD11	2.42	0.48
1:A:181:LEU:O	1:A:182:SER:C	2.51	0.48
1:A:162:THR:HA	1:A:264:VAL:HG12	1.95	0.47
1:A:189:ARG:NH1	1:A:189:ARG:HG3	2.29	0.47
1:A:2:MET:CE	1:A:5:ARG:NH1	2.77	0.47
1:A:174:THR:CG2	1:A:175:LYS:H	2.27	0.47
1:A:229:SER:OG	1:A:230:LEU:N	2.47	0.47
1:A:21:THR:HG22	1:A:23:LEU:HB2	1.97	0.47
1:A:180:ARG:N	1:A:180:ARG:HD2	2.28	0.47
1:A:103:GLY:O	1:A:106:SER:N	2.48	0.47
1:A:26:GLN:HA	1:A:39:PHE:O	2.15	0.47
1:A:5:ARG:CD	1:A:7:VAL:HG21	2.44	0.47
1:A:113:TYR:HE2	2:A:285:HOH:O	1.96	0.47
1:A:187:TYR:HB3	1:A:188:ILE:H	1.52	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:8:ALA:CB	1:A:32:PRO:HD3	2.45	0.47
1:A:71:PRO:O	1:A:74:TRP:CD1	2.68	0.47
1:A:75:ARG:N	1:A:75:ARG:HD3	2.15	0.47
1:A:167:GLU:O	1:A:169:GLN:N	2.43	0.47
1:A:180:ARG:HD3	1:A:180:ARG:N	2.28	0.47
1:A:215:LEU:O	1:A:219:ARG:HB2	2.15	0.47
1:A:5:ARG:CG	1:A:7:VAL:HG23	2.42	0.46
1:A:10:PRO:CB	1:A:73:PRO:HB3	2.45	0.46
1:A:242:TYR:O	1:A:242:TYR:CD2	2.68	0.46
1:A:11:GLN:HG3	1:A:12:GLN:CB	2.46	0.46
1:A:83:ILE:O	1:A:83:ILE:CG2	2.57	0.46
1:A:128:ILE:O	1:A:128:ILE:CG1	2.62	0.46
1:A:9:THR:CG2	1:A:12:GLN:CG	2.87	0.46
1:A:13:SER:O	1:A:135:VAL:HG21	2.15	0.46
1:A:79:HIS:CD2	1:A:79:HIS:C	2.89	0.46
1:A:106:SER:O	1:A:110:LYS:HG3	2.16	0.46
1:A:171:ASP:OD2	1:A:172:GLN:NE2	2.46	0.46
1:A:107:GLU:O	1:A:110:LYS:N	2.49	0.46
1:A:167:GLU:C	1:A:169:GLN:N	2.69	0.45
1:A:219:ARG:HG2	1:A:263:VAL:HG21	1.99	0.45
1:A:74:TRP:C	1:A:74:TRP:CE3	2.87	0.45
1:A:9:THR:HG22	1:A:11:GLN:CG	2.46	0.45
1:A:45:LYS:HB2	1:A:45:LYS:HE2	1.85	0.45
1:A:3:GLY:N	1:A:126:GLN:CB	2.80	0.45
1:A:191:ASN:O	1:A:194:ILE:HG22	2.17	0.45
1:A:21:THR:HG21	1:A:23:LEU:HB2	1.98	0.45
1:A:9:THR:HG21	1:A:12:GLN:HG2	1.97	0.44
1:A:162:THR:HG22	1:A:164:LEU:HD12	1.98	0.44
1:A:197:VAL:HA	1:A:238:GLN:O	2.18	0.44
1:A:164:LEU:CD1	1:A:164:LEU:N	2.80	0.44
1:A:266:SER:C	1:A:267:LEU:HD23	2.37	0.44
1:A:42:ARG:HH11	1:A:42:ARG:HG3	1.82	0.44
1:A:270:THR:HG21	1:A:271:GLN:CG	2.23	0.44
1:A:64:ASN:HA	1:A:88:PHE:CZ	2.53	0.44
1:A:103:GLY:O	1:A:104:ILE:C	2.55	0.44
1:A:71:PRO:C	1:A:73:PRO:HD2	2.37	0.43
1:A:82:ARG:O	1:A:112:ARG:NH1	2.41	0.43
1:A:196:LEU:N	1:A:236:ARG:O	2.50	0.43
1:A:14:GLN:O	1:A:30:PRO:HD2	2.18	0.43
1:A:21:THR:C	1:A:23:LEU:N	2.71	0.43
1:A:28:VAL:HG12	1:A:29:HIS:N	2.33	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:96:VAL:CG1	1:A:100:THR:HB	2.48	0.43
1:A:118:TYR:HB2	1:A:128:ILE:O	2.18	0.43
1:A:163:ILE:CG1	1:A:263:VAL:O	2.66	0.43
1:A:24:GLU:HA	1:A:41:SER:O	2.19	0.43
1:A:39:PHE:HZ	1:A:104:ILE:CG2	2.32	0.43
1:A:68:ILE:CG2	1:A:85:ASN:HA	2.48	0.43
1:A:103:GLY:C	1:A:105:LEU:N	2.71	0.43
1:A:266:SER:O	1:A:267:LEU:HD23	2.18	0.43
1:A:180:ARG:O	1:A:184:ILE:HG13	2.19	0.43
1:A:2:MET:HE2	1:A:2:MET:HB2	1.86	0.42
1:A:87:LYS:HD3	1:A:89:PHE:CE1	2.54	0.42
1:A:163:ILE:C	1:A:164:LEU:HD12	2.40	0.42
1:A:166:TYR:C	1:A:166:TYR:CD2	2.91	0.42
1:A:86:LEU:HD12	1:A:86:LEU:HA	1.82	0.42
1:A:199:VAL:HB	1:A:264:VAL:O	2.19	0.42
1:A:79:HIS:HD2	1:A:80:ALA:CA	2.29	0.42
1:A:124:ARG:NH1	1:A:124:ARG:HG2	2.33	0.42
1:A:155:SER:C	1:A:157:GLU:N	2.71	0.42
1:A:162:THR:HG22	1:A:164:LEU:CD1	2.50	0.42
1:A:232:LEU:HD23	1:A:232:LEU:O	2.19	0.42
1:A:168:ARG:O	1:A:169:GLN:C	2.57	0.42
1:A:262:ARG:HD2	1:A:264:VAL:HG21	2.02	0.42
1:A:73:PRO:HG3	1:A:113:TYR:OH	2.20	0.42
1:A:105:LEU:HG	1:A:141:TYR:CE2	2.54	0.42
1:A:166:TYR:HD2	1:A:166:TYR:C	2.22	0.42
1:A:105:LEU:HD12	1:A:105:LEU:HA	1.81	0.42
1:A:197:VAL:O	1:A:197:VAL:HG12	2.19	0.42
1:A:67:LEU:HD23	1:A:67:LEU:HA	1.89	0.42
1:A:73:PRO:O	1:A:75:ARG:NH1	2.53	0.42
1:A:223:LEU:HD23	1:A:223:LEU:H	1.82	0.41
1:A:1:VAL:CG1	1:A:126:GLN:NE2	2.84	0.41
1:A:6:TYR:O	1:A:7:VAL:CG2	2.68	0.41
1:A:71:PRO:O	1:A:74:TRP:NE1	2.54	0.41
1:A:87:LYS:HG3	1:A:88:PHE:N	2.35	0.41
1:A:15:TRP:O	1:A:140:LYS:CE	2.57	0.41
1:A:31:ILE:CD1	1:A:131:ALA:O	2.67	0.41
1:A:201:THR:O	1:A:201:THR:CG2	2.69	0.41
1:A:241:GLY:O	1:A:242:TYR:HB3	2.19	0.41
1:A:165:HIS:HB3	1:A:170:GLY:O	2.21	0.41
1:A:67:LEU:HB3	1:A:86:LEU:HB2	2.03	0.41
1:A:2:MET:HE1	1:A:5:ARG:NH1	2.36	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:34:PHE:O	1:A:56:ARG:HB3	2.20	0.41
1:A:121:TRP:O	1:A:122:GLN:O	2.39	0.41
1:A:158:ASP:OD1	1:A:158:ASP:C	2.59	0.41
1:A:18:VAL:HG21	1:A:28:VAL:HG21	2.02	0.41
1:A:176:ALA:O	1:A:177:SER:C	2.60	0.41
1:A:180:ARG:HD3	1:A:180:ARG:H	1.85	0.41
1:A:151:LEU:HD12	1:A:151:LEU:HA	1.98	0.40
1:A:21:THR:HG23	1:A:22:PRO:CD	2.46	0.40
1:A:267:LEU:HD23	1:A:267:LEU:HA	1.83	0.40
1:A:31:ILE:HD13	1:A:31:ILE:HA	1.73	0.40
1:A:108:LEU:O	1:A:134:SER:HB2	2.21	0.40
1:A:181:LEU:O	1:A:184:ILE:N	2.55	0.40
1:A:197:VAL:H	1:A:267:LEU:CD2	2.35	0.40
1:A:121:TRP:O	1:A:122:GLN:C	2.59	0.40
1:A:124:ARG:C	1:A:125:ASP:OD1	2.60	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:A:172:GLN:OE1	1:A:172:GLN:OE1[8_556]	2.08	0.12
1:A:171:ASP:OD2	1:A:222:SER:OG[8_556]	2.10	0.10

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	240/278 (86%)	181 (75%)	40 (17%)	19 (8%)	1 1

All (19) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	12	GLN
1	A	81	ASP
1	A	137	PHE
1	A	188	ILE
1	A	234	GLU
1	A	270	THR
1	A	75	ARG
1	A	104	ILE
1	A	135	VAL
1	A	269	ARG
1	A	70	MET
1	A	169	GLN
1	A	71	PRO
1	A	122	GLN
1	A	11	GLN
1	A	9	THR
1	A	231	GLY
1	A	130	VAL
1	A	32	PRO

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	219/247 (89%)	157 (72%)	62 (28%)	0 0

All (62) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	2	MET
1	A	5	ARG
1	A	9	THR
1	A	10	PRO
1	A	14	GLN
1	A	20	ASN
1	A	36	ASP
1	A	48	ASN

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Mol	Chain	Res	Type
1	A	49	LEU
1	A	58	PRO
1	A	59	MET
1	A	62	THR
1	A	66	SER
1	A	67	LEU
1	A	68	ILE
1	A	70	MET
1	A	74	TRP
1	A	75	ARG
1	A	79	HIS
1	A	83	ILE
1	A	84	THR
1	A	86	LEU
1	A	87	LYS
1	A	115	THR
1	A	119	GLN
1	A	121	TRP
1	A	122	GLN
1	A	124	ARG
1	A	125	ASP
1	A	126	GLN
1	A	127	ARG
1	A	128	ILE
1	A	145	SER
1	A	151	LEU
1	A	156	PHE
1	A	157	GLU
1	A	159	ILE
1	A	163	ILE
1	A	164	LEU
1	A	166	TYR
1	A	168	ARG
1	A	169	GLN
1	A	178	LYS
1	A	181	LEU
1	A	189	ARG
1	A	192	GLN
1	A	193	ASP
1	A	196	LEU
1	A	198	LEU
1	A	218	ARG

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Mol	Chain	Res	Type
1	A	221	GLU
1	A	223	LEU
1	A	224	ARG
1	A	228	GLN
1	A	232	LEU
1	A	236	ARG
1	A	240	GLN
1	A	242	TYR
1	A	264	VAL
1	A	265	ILE
1	A	270	THR
1	A	271	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	11	GLN
1	A	14	GLN
1	A	20	ASN
1	A	64	ASN
1	A	79	HIS
1	A	126	GLN
1	A	142	ASN
1	A	238	GLN
1	A	240	GLN

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 oligosaccharides in this entry.

5.6 Ligand geometry

There are no ligands in this entry.

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	246/278 (88%)	0.86	42 (17%) 5 5	33, 73, 116, 129	0

All (42) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	124	ARG	5.8
1	A	169	GLN	5.4
1	A	71	PRO	4.8
1	A	74	TRP	4.6
1	A	80	ALA	4.3
1	A	73	PRO	4.1
1	A	72	PRO	4.1
1	A	267	LEU	3.6
1	A	126	GLN	3.6
1	A	75	ARG	3.5
1	A	263	VAL	3.4
1	A	180	ARG	3.1
1	A	268	GLY	2.9
1	A	70	MET	2.9
1	A	81	ASP	2.9
1	A	266	SER	2.8
1	A	121	TRP	2.8
1	A	127	ARG	2.8
1	A	122	GLN	2.7
1	A	7	VAL	2.7
1	A	238	GLN	2.6
1	A	271	GLN	2.6
1	A	3	GLY	2.6
1	A	270	THR	2.6
1	A	160	ALA	2.5
1	A	32	PRO	2.5
1	A	242	TYR	2.5

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Mol	Chain	Res	Type	RSRZ
1	A	79	HIS	2.5
1	A	262	ARG	2.4
1	A	269	ARG	2.4
1	A	77	GLY	2.3
1	A	8	ALA	2.3
1	A	60	GLY	2.3
1	A	10	PRO	2.3
1	A	156	PHE	2.2
1	A	76	PRO	2.2
1	A	265	ILE	2.2
1	A	188	ILE	2.2
1	A	234	GLU	2.1
1	A	1	VAL	2.1
1	A	161	PHE	2.1
1	A	118	TYR	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 monosaccharides in this entry.

6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.