



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 02:10 AM UTC

PDB ID : 3IAO / pdb_00003iao
Title : Conformational plasticity of the coiled coil domain of BmrR is required for bmr promoter binding-the unliganded structure of BmrR
Authors : Kumaraswami, M.; Newberry, K.J.; Brennan, R.G.
Deposited on : 2009-07-14
Resolution : 2.80 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	:	4-5-2 with Phenix2.0
Xtriage (Phenix)	:	2.0
EDS	:	3.0
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

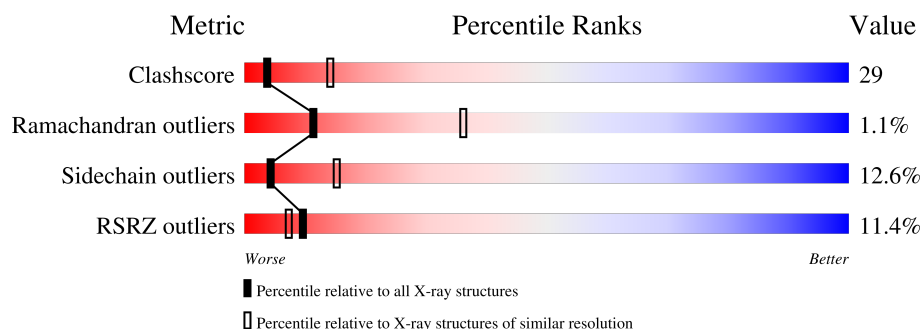
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 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))
Clashscore	190562	4276 (2.80-2.80)
Ramachandran outliers	187476	4196 (2.80-2.80)
Sidechain outliers	187428	4198 (2.80-2.80)
RSRZ outliers	180081	3869 (2.80-2.80)

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	11% 40% 46% 12% ..

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 2400 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 Multidrug-efflux transporter 1 regulator.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	273	Total	C	N	O	S	0	0	0
			2233	1442	353	430	8			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	253	GLN	GLU	engineered mutation	UNP P39075
A	275	GLU	ARG	engineered mutation	UNP P39075

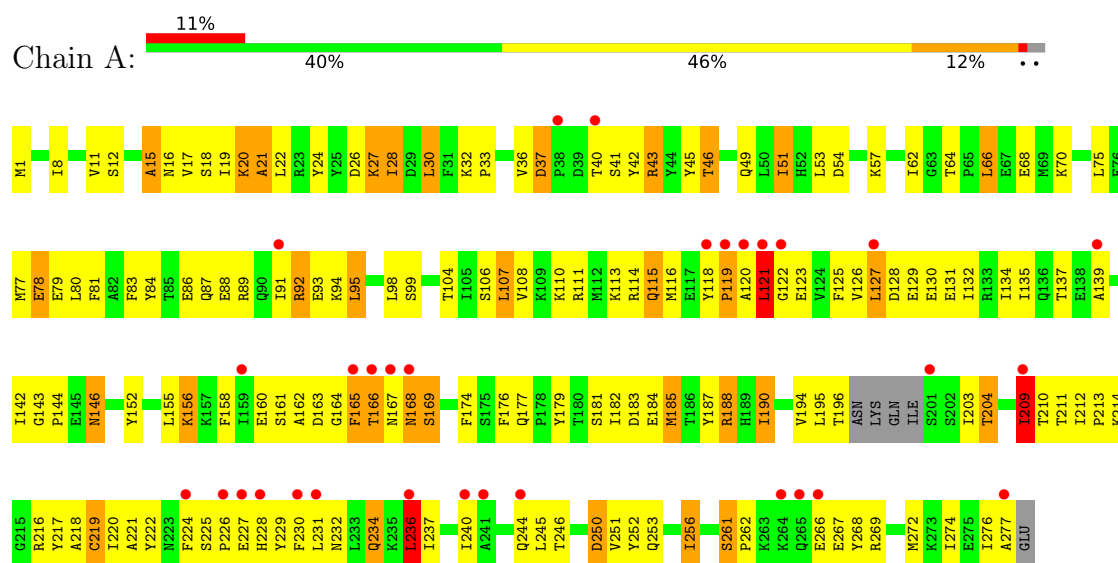
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	167	Total	O	0	0
			167	167		

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: Multidrug-efflux transporter 1 regulator



4 Data and refinement statistics

Property	Value	Source
Space group	I 41 2 2	Depositor
Cell constants a, b, c, α , β , γ	168.00Å 168.00Å 63.10Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 2.80 50.00 – 2.80	Depositor EDS
% Data completeness (in resolution range)	99.9 (50.00-2.80) 99.2 (50.00-2.80)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.93 (at 2.82Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.241 , 0.275 0.255 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	62.6	Xtriage
Anisotropy	0.198	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 64.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	2400	wwPDB-VP
Average B, all atoms (Å ²)	58.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.03% 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

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	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	0.55	0/2280	1.14	27/3081 (0.9%)

There are no bond length outliers.

All (27) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	166	THR	N-CA-C	-9.06	96.51	109.29
1	A	15	ALA	N-CA-C	-8.25	99.42	110.55
1	A	261	SER	N-CA-C	7.52	119.54	110.07
1	A	236	LEU	N-CA-C	-7.46	102.52	111.69
1	A	250	ASP	N-CA-C	-7.22	100.38	110.35
1	A	37	ASP	CA-C-N	7.01	126.63	119.19
1	A	37	ASP	C-N-CA	7.01	126.63	119.19
1	A	204	THR	CA-C-N	6.94	127.52	119.47
1	A	204	THR	C-N-CA	6.94	127.52	119.47
1	A	209	ILE	N-CA-C	-6.70	98.93	108.58
1	A	20	LYS	N-CA-C	-6.21	105.61	113.43
1	A	21	ALA	N-CA-C	-5.94	104.93	111.82
1	A	68	GLU	N-CA-C	-5.76	106.59	113.97
1	A	188	ARG	N-CA-C	-5.68	103.86	112.04
1	A	169	SER	N-CA-C	5.48	117.31	109.14
1	A	77	MET	N-CA-C	5.47	120.40	111.37
1	A	244	GLN	N-CA-C	-5.42	103.97	111.28
1	A	30	LEU	N-CA-C	-5.41	106.61	113.43
1	A	168	ASN	N-CA-C	-5.32	105.38	111.07
1	A	121	LEU	N-CA-C	5.16	117.02	108.20
1	A	122	GLY	N-CA-C	-5.15	107.56	116.01
1	A	203	ILE	N-CA-C	-5.12	101.09	108.87
1	A	224	PHE	N-CA-C	5.08	121.63	110.80
1	A	185	MET	N-CA-C	5.03	118.00	109.76
1	A	190	ILE	N-CA-C	-5.02	101.02	108.45
1	A	27	LYS	N-CA-C	5.01	116.75	111.28
1	A	165	PHE	N-CA-C	-5.01	104.06	110.53

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	2233	0	2219	128	0
2	A	167	0	0	4	1
All	All	2400	0	2219	128	1

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 29.

All (128) 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:120:ALA:HB3	1:A:125:PHE:HZ	1.28	0.97
1:A:120:ALA:HB3	1:A:125:PHE:CZ	2.11	0.85
1:A:113:LYS:HA	1:A:116:MET:HE3	1.60	0.83
1:A:152:TYR:O	1:A:156:LYS:HG3	1.78	0.82
1:A:26:ASP:OD1	1:A:32:LYS:HE3	1.85	0.76
1:A:130:GLU:HA	1:A:214:LYS:HB3	1.68	0.74
1:A:229:TYR:OH	1:A:253:GLN:NE2	2.18	0.74
1:A:75:LEU:HD22	1:A:79:GLU:HB3	1.69	0.74
1:A:234:GLN:HA	1:A:237:ILE:HD12	1.69	0.74
1:A:139:ALA:HB2	1:A:190:ILE:HG23	1.71	0.72
1:A:181:SER:C	1:A:230:PHE:HE2	1.97	0.72
1:A:156:LYS:HD3	1:A:167:ASN:OD1	1.90	0.71
1:A:166:THR:HG22	1:A:167:ASN:H	1.57	0.69
1:A:30:LEU:HD13	1:A:70:LYS:HG3	1.74	0.68
1:A:222:TYR:OH	1:A:229:TYR:HA	1.93	0.68
1:A:36:VAL:HG22	1:A:43:ARG:CG	2.24	0.68
1:A:218:ALA:O	1:A:236:LEU:HD21	1.94	0.68
1:A:40:THR:HG23	1:A:42:TYR:H	1.59	0.67
1:A:115:GLN:O	1:A:269:ARG:NH2	2.27	0.67
1:A:163:ASP:C	1:A:165:PHE:H	2.04	0.66
1:A:156:LYS:C	1:A:158:PHE:H	2.03	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:163:ASP:OD1	1:A:196:THR:HG23	1.97	0.65
1:A:216:ARG:HG3	2:A:427:HOH:O	1.97	0.64
1:A:36:VAL:HG22	1:A:43:ARG:HG3	1.81	0.62
1:A:40:THR:O	1:A:41:SER:HB3	2.00	0.61
1:A:181:SER:C	1:A:230:PHE:CE2	2.78	0.61
1:A:266:GLU:HB2	1:A:268:TYR:CE1	2.35	0.61
1:A:46:THR:HG22	1:A:49:GLN:H	1.67	0.59
1:A:266:GLU:HB2	1:A:268:TYR:HE1	1.68	0.59
1:A:18:SER:C	1:A:20:LYS:H	2.10	0.59
1:A:246:THR:OG1	1:A:277:ALA:HB3	2.03	0.58
1:A:104:THR:O	1:A:108:VAL:HG23	2.04	0.57
1:A:46:THR:HB	1:A:49:GLN:HG2	1.85	0.57
1:A:236:LEU:HD13	1:A:274:ILE:HD11	1.87	0.57
1:A:168:ASN:N	1:A:168:ASN:OD1	2.36	0.57
1:A:194:VAL:HG12	1:A:196:THR:H	1.70	0.57
1:A:36:VAL:HG22	1:A:43:ARG:HG2	1.86	0.56
1:A:28:ILE:O	1:A:70:LYS:NZ	2.40	0.55
1:A:220:ILE:HD11	1:A:232:ASN:OD1	2.06	0.55
1:A:120:ALA:CB	1:A:125:PHE:HZ	2.12	0.55
1:A:62:ILE:HG13	1:A:64:THR:OG1	2.08	0.54
1:A:37:ASP:HB3	1:A:40:THR:CG2	2.38	0.54
1:A:114:ARG:O	1:A:115:GLN:C	2.51	0.53
1:A:78:GLU:O	1:A:81:PHE:HB3	2.09	0.53
1:A:176:PHE:CE1	1:A:240:ILE:HD12	2.44	0.53
1:A:127:LEU:HD21	1:A:219:CYS:SG	2.49	0.53
1:A:24:TYR:O	1:A:28:ILE:HG12	2.09	0.53
1:A:135:ILE:HG12	1:A:194:VAL:CG2	2.38	0.53
1:A:179:TYR:HD2	1:A:184:GLU:HB3	1.74	0.52
1:A:37:ASP:OD2	1:A:40:THR:HG22	2.10	0.52
1:A:94:LYS:O	1:A:98:LEU:HG	2.08	0.52
1:A:236:LEU:HD13	1:A:274:ILE:CD1	2.41	0.51
1:A:66:LEU:CD2	1:A:70:LYS:HD2	2.39	0.51
1:A:89:ARG:NH1	1:A:93:GLU:OE2	2.43	0.51
1:A:106:SER:O	1:A:110:LYS:HG3	2.11	0.51
1:A:176:PHE:CD1	1:A:274:ILE:HD13	2.46	0.51
1:A:115:GLN:HG3	1:A:269:ARG:HG2	1.92	0.51
1:A:160:GLU:O	1:A:164:GLY:N	2.40	0.50
1:A:195:LEU:HD12	1:A:195:LEU:N	2.27	0.50
1:A:176:PHE:HD1	1:A:274:ILE:HD13	1.77	0.50
1:A:11:VAL:HG21	1:A:45:TYR:CG	2.47	0.49
1:A:188:ARG:NH1	1:A:250:ASP:OD1	2.45	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:17:VAL:HG12	1:A:21:ALA:HB3	1.95	0.49
1:A:181:SER:O	1:A:184:GLU:HB2	2.13	0.49
1:A:126:VAL:HA	1:A:217:TYR:O	2.13	0.49
1:A:230:PHE:CE1	1:A:234:GLN:NE2	2.81	0.49
1:A:174:PHE:CZ	1:A:251:VAL:HG11	2.47	0.48
1:A:163:ASP:CG	1:A:195:LEU:HB2	2.38	0.47
1:A:182:ILE:N	1:A:230:PHE:HE2	2.12	0.47
1:A:80:LEU:O	1:A:83:PHE:HB3	2.14	0.47
1:A:161:SER:HB3	2:A:399:HOH:O	2.15	0.47
1:A:8:ILE:HG13	1:A:22:LEU:HD13	1.96	0.47
1:A:24:TYR:O	1:A:28:ILE:CG1	2.64	0.46
1:A:113:LYS:HA	1:A:116:MET:CE	2.40	0.46
1:A:179:TYR:HA	1:A:184:GLU:OE1	2.16	0.46
1:A:135:ILE:HG12	1:A:194:VAL:HG23	1.97	0.46
1:A:32:LYS:O	1:A:33:PRO:C	2.59	0.45
1:A:163:ASP:OD1	1:A:195:LEU:HB2	2.16	0.45
1:A:88:GLU:HA	1:A:91:ILE:HG22	1.99	0.45
1:A:75:LEU:HD22	1:A:79:GLU:CB	2.42	0.45
1:A:115:GLN:OE1	1:A:115:GLN:HA	2.16	0.45
1:A:266:GLU:O	1:A:267:GLU:HG2	2.17	0.45
1:A:137:THR:HG22	1:A:155:LEU:HD11	1.99	0.45
1:A:228:HIS:HB3	1:A:232:ASN:ND2	2.32	0.45
1:A:118:TYR:O	1:A:119:PRO:C	2.59	0.44
1:A:234:GLN:O	1:A:237:ILE:HB	2.18	0.44
1:A:127:LEU:N	1:A:127:LEU:HD13	2.31	0.44
1:A:30:LEU:CD1	1:A:70:LYS:HG3	2.45	0.44
1:A:160:GLU:C	1:A:162:ALA:H	2.26	0.44
1:A:261:SER:HA	1:A:262:PRO:HD3	1.79	0.44
1:A:83:PHE:O	1:A:86:GLU:HB3	2.18	0.44
1:A:104:THR:HG21	2:A:379:HOH:O	2.18	0.44
1:A:128:ASP:OD2	1:A:128:ASP:N	2.51	0.43
1:A:118:TYR:CE2	1:A:221:ALA:HB3	2.53	0.43
1:A:225:SER:C	1:A:227:GLU:H	2.26	0.43
1:A:256:ILE:HG23	1:A:269:ARG:HB3	2.00	0.43
1:A:211:THR:OG1	1:A:212:ILE:N	2.52	0.43
1:A:84:TYR:O	1:A:87:GLN:N	2.47	0.43
1:A:245:LEU:HD22	1:A:276:ILE:HD12	2.01	0.43
1:A:12:SER:O	1:A:15:ALA:O	2.37	0.43
1:A:146:ASN:HD22	1:A:146:ASN:HA	1.54	0.43
1:A:182:ILE:O	1:A:185:MET:HG2	2.19	0.43
1:A:1:MET:N	2:A:390:HOH:O	2.33	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:107:LEU:HD13	1:A:195:LEU:HB3	2.00	0.43
1:A:53:LEU:O	1:A:57:LYS:HG3	2.20	0.42
1:A:135:ILE:CD1	1:A:209:ILE:HD13	2.50	0.42
1:A:163:ASP:C	1:A:165:PHE:N	2.73	0.42
1:A:115:GLN:HG3	1:A:269:ARG:NE	2.35	0.42
1:A:129:GLU:C	1:A:130:GLU:HG2	2.45	0.42
1:A:176:PHE:HE1	1:A:240:ILE:HD12	1.85	0.42
1:A:51:ILE:HD13	1:A:54:ASP:OD2	2.20	0.42
1:A:210:THR:HG23	1:A:211:THR:N	2.33	0.42
1:A:252:TYR:O	1:A:272:MET:HA	2.19	0.42
1:A:256:ILE:O	1:A:269:ARG:HB3	2.20	0.42
1:A:156:LYS:HD2	1:A:166:THR:CG2	2.50	0.42
1:A:12:SER:OG	1:A:17:VAL:O	2.34	0.41
1:A:37:ASP:HB3	1:A:40:THR:HG22	2.02	0.41
1:A:84:TYR:C	1:A:86:GLU:N	2.72	0.41
1:A:113:LYS:O	1:A:114:ARG:C	2.63	0.41
1:A:144:PRO:HA	1:A:187:TYR:OH	2.20	0.41
1:A:16:ASN:H	1:A:57:LYS:NZ	2.18	0.41
1:A:132:ILE:O	1:A:134:ILE:HG23	2.21	0.41
1:A:118:TYR:HD2	1:A:120:ALA:O	2.04	0.41
1:A:121:LEU:HG	1:A:222:TYR:HA	2.02	0.41
1:A:11:VAL:HG21	1:A:45:TYR:CB	2.51	0.40
1:A:156:LYS:C	1:A:158:PHE:N	2.71	0.40
1:A:92:ARG:NH2	1:A:95:LEU:HD22	2.37	0.40
1:A:143:GLY:H	1:A:146:ASN:HB2	1.87	0.40

All (1) 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)
2:A:391:HOH:O	2:A:391:HOH:O[5_555]	0.51	1.69

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	269/278 (97%)	223 (83%)	43 (16%)	3 (1%)	11	36

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	213	PRO
1	A	226	PRO
1	A	19	ILE

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	246/255 (96%)	215 (87%)	31 (13%)	4	15

All (31) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	27	LYS
1	A	28	ILE
1	A	43	ARG
1	A	46	THR
1	A	51	ILE
1	A	66	LEU
1	A	78	GLU
1	A	92	ARG
1	A	95	LEU
1	A	99	SER
1	A	107	LEU
1	A	111	ARG
1	A	115	GLN
1	A	119	PRO
1	A	121	LEU
1	A	123	GLU

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Mol	Chain	Res	Type
1	A	127	LEU
1	A	131	GLU
1	A	142	ILE
1	A	146	ASN
1	A	156	LYS
1	A	169	SER
1	A	177	GLN
1	A	183	ASP
1	A	204	THR
1	A	209	ILE
1	A	219	CYS
1	A	231	LEU
1	A	234	GLN
1	A	236	LEU
1	A	256	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	136	GLN
1	A	146	ASN
1	A	177	GLN
1	A	234	GLN
1	A	253	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	273/278 (98%)	0.93	31 (11%) 10 7	33, 58, 79, 95	0

All (31) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	240	ILE	3.5
1	A	166	THR	3.2
1	A	40	THR	3.0
1	A	201	SER	2.8
1	A	231	LEU	2.7
1	A	122	GLY	2.6
1	A	127	LEU	2.6
1	A	38	PRO	2.5
1	A	120	ALA	2.4
1	A	228	HIS	2.4
1	A	167	ASN	2.4
1	A	165	PHE	2.4
1	A	224	PHE	2.4
1	A	168	ASN	2.3
1	A	277	ALA	2.3
1	A	226	PRO	2.3
1	A	266	GLU	2.3
1	A	118	TYR	2.3
1	A	91	ILE	2.3
1	A	230	PHE	2.3
1	A	227	GLU	2.2
1	A	241	ALA	2.2
1	A	121	LEU	2.2
1	A	139	ALA	2.2
1	A	265	GLN	2.1
1	A	236	LEU	2.1
1	A	209	ILE	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	159	ILE	2.1
1	A	244	GLN	2.0
1	A	119	PRO	2.0
1	A	264	LYS	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 oligosaccharides 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.