



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 06:15 AM UTC

PDB ID : 1O67 / pdb_00001o67
Title : Crystal structure of an hypothetical protein
Authors : Structural GenomiX
Deposited on : 2003-10-23
Resolution : 2.54 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
Mogul	:	2022.3.0, CSD as543be (2022)
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

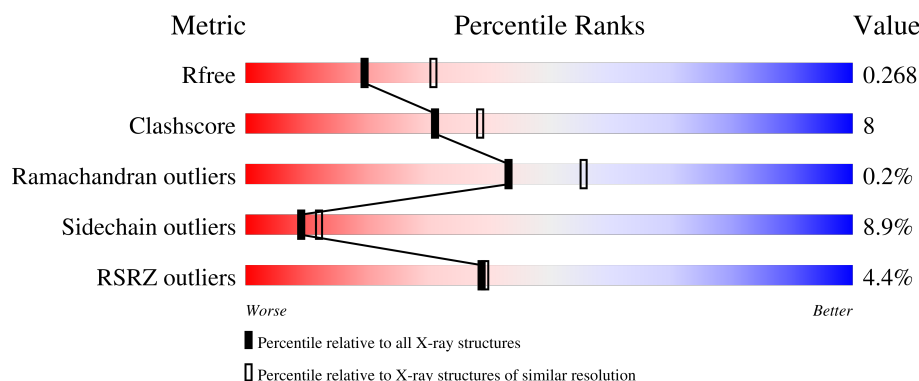
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.54 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}	180053	1091 (2.54-2.54)
Clashscore	190562	1120 (2.54-2.54)
Ramachandran outliers	187476	1106 (2.54-2.54)
Sidechain outliers	187428	1106 (2.54-2.54)
RSRZ outliers	180081	1091 (2.54-2.54)

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	246	2% 68% 16% • 13%
1	B	246	7% 68% 17% • 13%
1	C	246	2% 63% 18% • 15%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 5124 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 Hypothetical protein yiiM.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	214	Total	C	N	O	S	Se	0	0	0
			1709	1087	294	320	2	6			
1	B	214	Total	C	N	O	S	Se	0	1	0
			1703	1080	297	318	2	6			
1	C	210	Total	C	N	O	S	Se	0	0	0
			1682	1067	290	317	2	6			

There are 57 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	MSE	-	cloning artifact	UNP P32157
A	0	SER	-	cloning artifact	UNP P32157
A	1	LEU	-	cloning artifact	UNP P32157
A	11	MSE	MET	modified residue	UNP P32157
A	43	MSE	MET	modified residue	UNP P32157
A	110	MSE	ILE	engineered mutation	UNP P32157
A	146	MSE	MET	modified residue	UNP P32157
A	191	MSE	MET	modified residue	UNP P32157
A	215	MSE	MET	modified residue	UNP P32157
A	235	GLU	-	cloning artifact	UNP P32157
A	236	GLY	-	cloning artifact	UNP P32157
A	237	GLY	-	cloning artifact	UNP P32157
A	238	SER	-	cloning artifact	UNP P32157
A	239	HIS	-	cloning artifact	UNP P32157
A	240	HIS	-	cloning artifact	UNP P32157
A	241	HIS	-	cloning artifact	UNP P32157
A	242	HIS	-	cloning artifact	UNP P32157
A	243	HIS	-	cloning artifact	UNP P32157
A	244	HIS	-	cloning artifact	UNP P32157
B	-1	MSE	-	cloning artifact	UNP P32157
B	0	SER	-	cloning artifact	UNP P32157
B	1	LEU	-	cloning artifact	UNP P32157
B	11	MSE	MET	modified residue	UNP P32157

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Chain	Residue	Modelled	Actual	Comment	Reference
B	43	MSE	MET	modified residue	UNP P32157
B	110	MSE	ILE	engineered mutation	UNP P32157
B	146	MSE	MET	modified residue	UNP P32157
B	191	MSE	MET	modified residue	UNP P32157
B	215	MSE	MET	modified residue	UNP P32157
B	235	GLU	-	cloning artifact	UNP P32157
B	236	GLY	-	cloning artifact	UNP P32157
B	237	GLY	-	cloning artifact	UNP P32157
B	238	SER	-	cloning artifact	UNP P32157
B	239	HIS	-	cloning artifact	UNP P32157
B	240	HIS	-	cloning artifact	UNP P32157
B	241	HIS	-	cloning artifact	UNP P32157
B	242	HIS	-	cloning artifact	UNP P32157
B	243	HIS	-	cloning artifact	UNP P32157
B	244	HIS	-	cloning artifact	UNP P32157
C	-1	MSE	-	cloning artifact	UNP P32157
C	0	SER	-	cloning artifact	UNP P32157
C	1	LEU	-	cloning artifact	UNP P32157
C	11	MSE	MET	modified residue	UNP P32157
C	43	MSE	MET	modified residue	UNP P32157
C	110	MSE	ILE	engineered mutation	UNP P32157
C	146	MSE	MET	modified residue	UNP P32157
C	191	MSE	MET	modified residue	UNP P32157
C	215	MSE	MET	modified residue	UNP P32157
C	235	GLU	-	cloning artifact	UNP P32157
C	236	GLY	-	cloning artifact	UNP P32157
C	237	GLY	-	cloning artifact	UNP P32157
C	238	SER	-	cloning artifact	UNP P32157
C	239	HIS	-	cloning artifact	UNP P32157
C	240	HIS	-	cloning artifact	UNP P32157
C	241	HIS	-	cloning artifact	UNP P32157
C	242	HIS	-	cloning artifact	UNP P32157
C	243	HIS	-	cloning artifact	UNP P32157
C	244	HIS	-	cloning artifact	UNP P32157

- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	12	Total O 12 12	0	0
2	B	5	Total O 5 5	0	0

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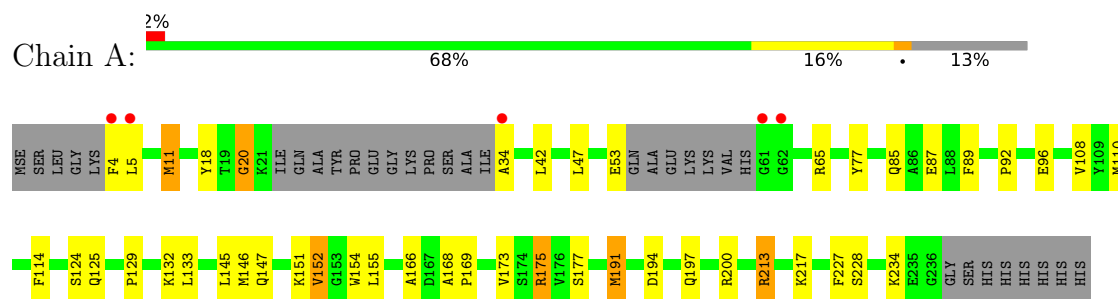
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	C	13	Total	O	0	0
			13	13		

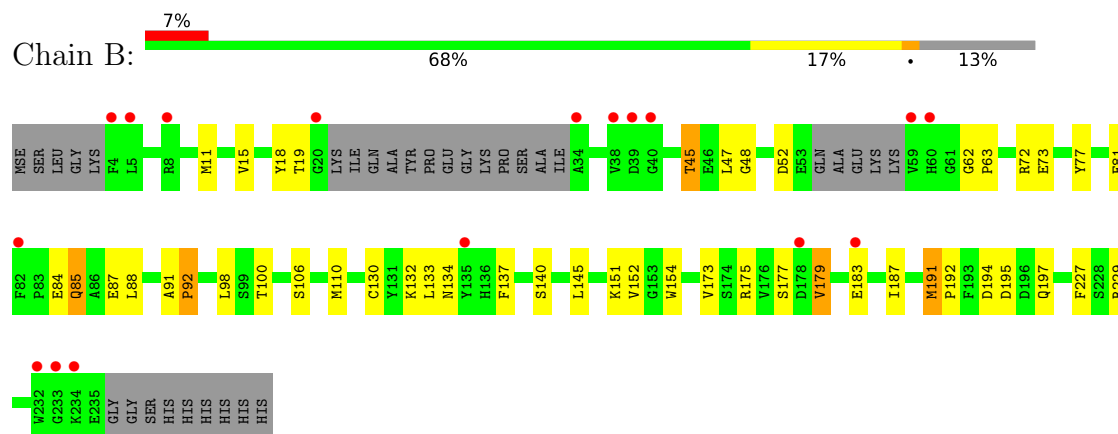
3 Residue-property plots [i](#)

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.

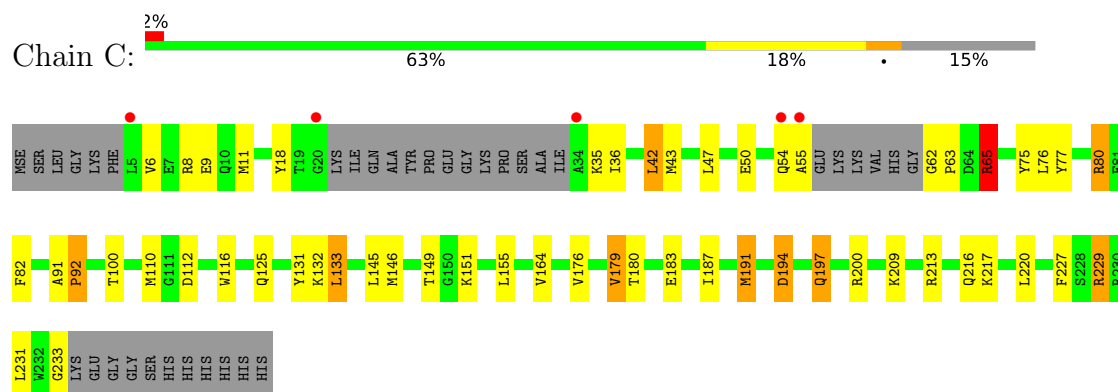
• Molecule 1: Hypothetical protein yiiM



• Molecule 1: Hypothetical protein yiiM



• Molecule 1: Hypothetical protein yiiM



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	74.53Å 99.33Å 99.41Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	59.63 – 2.54 59.63 – 2.54	Depositor EDS
% Data completeness (in resolution range)	(Not available) (59.63-2.54) 99.1 (59.63-2.54)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.46 (at 2.55Å)	Xtriage
Refinement program	REFMAC 4.0	Depositor
R, R_{free}	0.239 , 0.305 0.213 , 0.268	Depositor DCC
R_{free} test set	1257 reflections (5.09%)	wwPDB-VP
Wilson B-factor (Å ²)	42.5	Xtriage
Anisotropy	0.352	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 45.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.016 for -h,l,k	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	5124	wwPDB-VP
Average B, all atoms (Å ²)	39.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.24% 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	0.69	0/1745	1.45	11/2353 (0.5%)
1	B	0.63	0/1745	1.35	14/2356 (0.6%)
1	C	0.67	0/1717	1.43	12/2317 (0.5%)
All	All	0.67	0/5207	1.41	37/7026 (0.5%)

There are no bond length outliers.

All (37) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	20	GLY	N-CA-C	8.41	125.90	111.19
1	A	213	ARG	CD-NE-CZ	7.63	135.08	124.40
1	B	227	PHE	CA-CB-CG	7.25	121.05	113.80
1	C	65	ARG	CD-NE-CZ	7.08	134.31	124.40
1	A	89	PHE	CA-CB-CG	-6.47	107.33	113.80
1	A	227	PHE	CA-CB-CG	6.45	120.25	113.80
1	C	227	PHE	CA-CB-CG	6.13	119.93	113.80
1	A	96	GLU	N-CA-C	6.08	118.62	110.35
1	C	231	LEU	N-CA-C	6.03	117.93	111.36
1	C	82	PHE	CA-C-N	5.60	125.27	119.05
1	C	82	PHE	C-N-CA	5.60	125.27	119.05
1	A	228	SER	N-CA-C	5.55	117.77	111.11
1	B	137	PHE	N-CA-C	-5.53	106.58	113.38
1	B	152	VAL	N-CA-C	5.51	118.93	113.53
1	A	166	ALA	CA-C-N	5.50	131.77	122.26
1	A	166	ALA	C-N-CA	5.50	131.77	122.26
1	B	154	TRP	CA-C-O	-5.48	115.58	121.45
1	A	200	ARG	NE-CZ-NH1	-5.34	116.16	121.50
1	A	114	PHE	CA-CB-CG	5.28	119.08	113.80
1	C	112	ASP	CA-CB-CG	5.25	117.84	112.60
1	B	100	THR	N-CA-C	5.24	115.06	108.45
1	B	132	LYS	CA-C-N	5.21	127.26	120.28
1	B	132	LYS	C-N-CA	5.21	127.26	120.28
1	A	152	VAL	N-CA-CB	-5.21	106.48	112.21

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	100	THR	N-CA-C	5.20	115.15	108.34
1	B	152	VAL	CB-CA-C	-5.15	105.60	110.65
1	B	72	ARG	CD-NE-CZ	5.14	131.60	124.40
1	B	152	VAL	N-CA-CB	-5.13	105.19	112.12
1	B	151	LYS	CA-C-N	5.10	127.83	121.71
1	B	151	LYS	C-N-CA	5.10	127.83	121.71
1	B	195	ASP	CA-CB-CG	-5.05	107.55	112.60
1	C	194	ASP	CA-C-N	5.04	127.35	120.54
1	C	194	ASP	C-N-CA	5.04	127.35	120.54
1	C	75	TYR	CA-C-N	5.04	127.29	120.44
1	C	75	TYR	C-N-CA	5.04	127.29	120.44
1	C	176	VAL	N-CA-C	5.00	117.18	111.88
1	B	175	ARG	NE-CZ-NH1	-5.00	116.50	121.50

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	1709	0	1641	22	0
1	B	1703	0	1619	19	0
1	C	1682	0	1613	41	0
2	A	12	0	0	0	0
2	B	5	0	0	0	0
2	C	13	0	0	0	0
All	All	5124	0	4873	82	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 8.

All (82) 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:C:65:ARG:HB3	1:C:65:ARG:HH11	1.28	0.99

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:191:MSE:HE1	1:B:194:ASP:HB3	1.63	0.78
1:C:179:VAL:HG23	1:C:183:GLU:HG3	1.67	0.76
1:A:191:MSE:HE1	1:A:194:ASP:HB3	1.70	0.73
1:A:65:ARG:HG2	1:A:155:LEU:HD13	1.69	0.73
1:C:191:MSE:SE	1:C:197:GLN:HG2	2.43	0.68
1:A:146:MSE:HE3	1:A:152:VAL:HA	1.75	0.68
1:A:110:MSE:HE1	1:A:125:GLN:HA	1.80	0.63
1:B:62:GLY:N	1:B:63:PRO:HD2	2.14	0.62
1:C:65:ARG:HB3	1:C:65:ARG:NH1	2.07	0.61
1:C:149:THR:HB	1:C:151:LYS:HE2	1.82	0.60
1:A:11:MSE:HA	1:A:11:MSE:HE2	1.84	0.60
1:C:80:ARG:NH1	1:C:80:ARG:HB2	2.17	0.59
1:A:213:ARG:O	1:A:217:LYS:HG2	2.02	0.59
1:A:77:TYR:CG	1:A:151:LYS:HE3	2.38	0.59
1:B:45:THR:HG23	1:B:47:LEU:H	1.67	0.59
1:A:4:PHE:CD1	1:A:175:ARG:HG2	2.39	0.56
1:C:62:GLY:N	1:C:63:PRO:CD	2.69	0.56
1:C:187:ILE:HG23	1:C:191:MSE:HE2	1.89	0.55
1:C:43:MSE:H	1:C:50:GLU:HG2	1.71	0.54
1:A:146:MSE:CE	1:A:152:VAL:HA	2.39	0.53
1:C:183:GLU:OE1	1:C:200:ARG:NH1	2.37	0.53
1:C:216:GLN:HE21	1:C:220:LEU:HG	1.73	0.53
1:C:133:LEU:HD23	1:C:146:MSE:SE	2.60	0.52
1:A:191:MSE:SE	1:A:197:GLN:HG3	2.61	0.51
1:B:191:MSE:SE	1:B:197:GLN:HG2	2.60	0.51
1:B:187:ILE:CD1	1:B:197:GLN:HG3	2.40	0.51
1:C:65:ARG:CB	1:C:155:LEU:HD13	2.41	0.50
1:A:108:VAL:HG21	1:A:154:TRP:CZ2	2.47	0.50
1:C:42:LEU:H	1:C:42:LEU:HD23	1.76	0.50
1:B:85:GLN:HB2	1:B:88:LEU:HD12	1.94	0.49
1:B:45:THR:HG22	1:B:48:GLY:O	2.12	0.49
1:C:191:MSE:HE1	1:C:197:GLN:HB3	1.94	0.48
1:C:213:ARG:NH2	1:C:217:LYS:HZ1	2.12	0.48
1:C:77:TYR:CE2	1:C:151:LYS:HE3	2.49	0.47
1:A:146:MSE:HE2	1:A:152:VAL:HG12	1.96	0.47
1:C:54:GLN:O	1:C:55:ALA:C	2.58	0.47
1:C:187:ILE:CD1	1:C:197:GLN:HG3	2.44	0.47
1:B:18:TYR:CE2	1:B:92:PRO:HD3	2.49	0.46
1:B:134:ASN:ND2	1:B:140:SER:O	2.48	0.46
1:C:42:LEU:HD23	1:C:42:LEU:N	2.31	0.46
1:C:77:TYR:CD2	1:C:151:LYS:HE3	2.51	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:180:THR:OG1	1:C:183:GLU:HG2	2.16	0.45
1:A:18:TYR:CD2	1:A:92:PRO:HG3	2.51	0.45
1:A:191:MSE:HE1	1:A:197:GLN:HB2	1.98	0.45
1:B:77:TYR:CE1	1:B:81:GLU:HG3	2.51	0.45
1:C:42:LEU:HB2	1:C:50:GLU:CG	2.46	0.45
1:A:20:GLY:HA2	1:A:34:ALA:O	2.17	0.44
1:A:4:PHE:HD1	1:A:175:ARG:HG2	1.81	0.44
1:B:15:VAL:HG21	1:B:98:LEU:HD22	2.00	0.44
1:B:62:GLY:N	1:B:63:PRO:CD	2.80	0.44
1:A:65:ARG:HG2	1:A:155:LEU:CD1	2.44	0.44
1:C:213:ARG:CZ	1:C:217:LYS:NZ	2.82	0.43
1:A:129:PRO:HG2	1:A:147:GLN:HB2	2.01	0.43
1:C:116:TRP:O	1:C:164:VAL:HG12	2.19	0.43
1:B:194:ASP:OD2	1:B:197:GLN:HB2	2.18	0.42
1:B:191:MSE:HE3	1:B:192:PRO:O	2.18	0.42
1:B:18:TYR:CD2	1:B:92:PRO:HG3	2.55	0.42
1:C:47:LEU:HD23	1:C:47:LEU:HA	1.79	0.42
1:A:4:PHE:CE1	1:A:175:ARG:HG2	2.55	0.42
1:C:229:ARG:CG	1:C:229:ARG:HH11	2.33	0.42
1:C:91:ALA:HA	1:C:92:PRO:HA	1.68	0.42
1:B:110:MSE:HG3	1:B:179:VAL:HG22	2.02	0.42
1:C:42:LEU:HB2	1:C:50:GLU:HG2	2.02	0.42
1:C:65:ARG:HB3	1:C:155:LEU:HD13	1.99	0.42
1:C:131:TYR:OH	1:C:132:LYS:HE3	2.20	0.42
1:B:183:GLU:O	1:B:187:ILE:HG13	2.20	0.42
1:C:80:ARG:HB2	1:C:80:ARG:HH11	1.84	0.42
1:C:110:MSE:HE1	1:C:125:GLN:HA	2.01	0.42
1:B:91:ALA:HA	1:B:92:PRO:HA	1.81	0.41
1:C:229:ARG:HD3	1:C:233:GLY:O	2.21	0.41
1:C:191:MSE:HE1	1:C:194:ASP:HB3	2.02	0.41
1:A:146:MSE:HE2	1:A:146:MSE:HB3	1.95	0.41
1:C:65:ARG:HB2	1:C:155:LEU:HD13	2.02	0.41
1:A:168:ALA:HA	1:A:169:PRO:HD2	1.96	0.41
1:C:213:ARG:NH2	1:C:217:LYS:NZ	2.68	0.41
1:C:229:ARG:HD3	1:C:229:ARG:HA	1.87	0.41
1:C:213:ARG:HG3	1:C:217:LYS:HE3	2.03	0.41
1:A:124:SER:O	1:A:125:GLN:HB3	2.21	0.40
1:C:18:TYR:HB3	1:C:35:LYS:HB3	2.03	0.40
1:B:19:THR:HB	1:B:52:ASP:HB2	2.04	0.40
1:C:187:ILE:HD13	1:C:197:GLN:HG3	2.03	0.40

There are no symmetry-related clashes.

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	208/246 (85%)	203 (98%)	5 (2%)	0	100	100
1	B	209/246 (85%)	203 (97%)	6 (3%)	0	100	100
1	C	204/246 (83%)	194 (95%)	9 (4%)	1 (0%)	24	34
All	All	621/738 (84%)	600 (97%)	20 (3%)	1 (0%)	43	56

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	92	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	178/199 (89%)	163 (92%)	15 (8%)	10	14
1	B	176/199 (88%)	160 (91%)	16 (9%)	9	11
1	C	176/199 (88%)	160 (91%)	16 (9%)	9	11
All	All	530/597 (89%)	483 (91%)	47 (9%)	9	12

All (47) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	5	LEU
1	A	11	MSE

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Mol	Chain	Res	Type
1	A	42	LEU
1	A	47	LEU
1	A	53	GLU
1	A	85	GLN
1	A	87	GLU
1	A	132	LYS
1	A	133	LEU
1	A	145	LEU
1	A	173	VAL
1	A	175	ARG
1	A	177	SER
1	A	191	MSE
1	A	234	LYS
1	B	11	MSE
1	B	45	THR
1	B	73	GLU
1	B	84	GLU
1	B	85	GLN
1	B	87	GLU
1	B	92	PRO
1	B	106	SER
1	B	130	CYS
1	B	133	LEU
1	B	145	LEU
1	B	173	VAL
1	B	177	SER
1	B	179	VAL
1	B	191	MSE
1	B	229	ARG
1	C	6	VAL
1	C	8	ARG
1	C	9	GLU
1	C	11	MSE
1	C	36	ILE
1	C	42	LEU
1	C	65	ARG
1	C	76	LEU
1	C	80	ARG
1	C	133	LEU
1	C	145	LEU
1	C	179	VAL
1	C	191	MSE

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Mol	Chain	Res	Type
1	C	197	GLN
1	C	209	LYS
1	C	229	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	85	GLN
1	A	136	HIS
1	A	182	GLN
1	A	197	GLN
1	B	10	GLN
1	B	122	GLN
1	C	197	GLN
1	C	216	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 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 ⓘ

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	208/246 (84%)	0.01	5 (2%) 59 61	18, 32, 60, 70	0
1	B	208/246 (84%)	0.42	17 (8%) 17 17	23, 42, 73, 87	1 (0%)
1	C	204/246 (82%)	0.08	5 (2%) 58 59	20, 35, 59, 96	0
All	All	620/738 (84%)	0.17	27 (4%) 39 39	18, 36, 68, 96	1 (0%)

All (27) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	55	ALA	4.3
1	B	5	LEU	4.1
1	B	4	PHE	3.9
1	B	234	LYS	3.4
1	A	4	PHE	3.3
1	B	59	VAL	3.3
1	B	34	ALA	3.2
1	A	61	GLY	2.7
1	A	34	ALA	2.6
1	B	40	GLY	2.6
1	C	34	ALA	2.4
1	B	8	ARG	2.4
1	B	135	TYR	2.4
1	C	5	LEU	2.4
1	C	54	GLN	2.4
1	B	38	VAL	2.4
1	A	62	GLY	2.3
1	A	5	LEU	2.3
1	B	39	ASP	2.3
1	B	20	GLY	2.3
1	B	178	ASP	2.3
1	B	60	HIS	2.2
1	C	20	GLY	2.2

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Mol	Chain	Res	Type	RSRZ
1	B	183	GLU	2.2
1	B	232	TRP	2.2
1	B	233	GLY	2.1
1	B	82	PHE	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.