



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 1, 2026 – 08:14 AM UTC

PDB ID : 2F7Y / pdb_00002f7y
Title : Crystal structure of Molybdenum cofactor biosynthesis protein Mog from *Shewanella oneidensis*
Authors : Chang, C.; Mulligan, R.; Collart, F.; Joachimiak, A.; Midwest Center for Structural Genomics (MCSG)
Deposited on : 2005-12-01
Resolution : 2.30 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

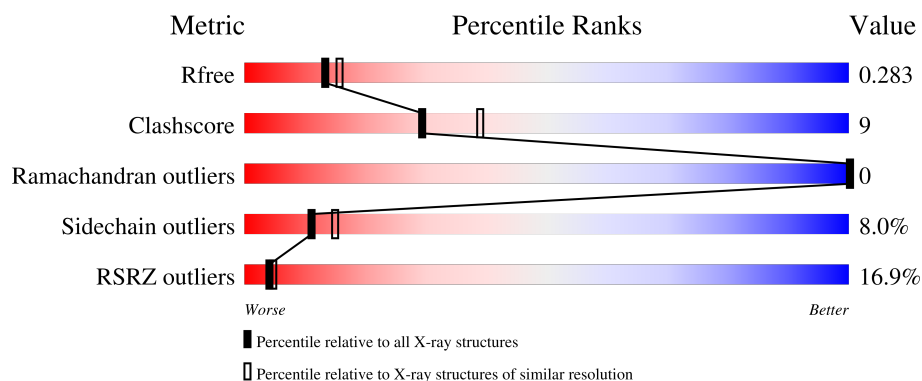
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	6319 (2.30-2.30)
Clashscore	190562	6919 (2.30-2.30)
Ramachandran outliers	187476	6854 (2.30-2.30)
Sidechain outliers	187428	6854 (2.30-2.30)
RSRZ outliers	180081	6325 (2.30-2.30)

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	177	16% 82% 12% ...
1	B	177	16% 67% 23% 5%
1	C	177	15% 81% 12% ...

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 4026 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 molybdenum cofactor biosynthesis protein Mog.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	173	Total	C	N	O	S	Se	0	0	0
			1316	832	214	259	6	5			
1	B	168	Total	C	N	O	S	Se	0	0	0
			1286	814	208	253	6	5			
1	C	171	Total	C	N	O	S	Se	0	0	0
			1302	823	211	257	6	5			

There are 18 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MSE	MET	modified residue	GB 24371665
A	61	MSE	MET	modified residue	GB 24371665
A	95	MSE	MET	modified residue	GB 24371665
A	96	MSE	MET	modified residue	GB 24371665
A	103	MSE	MET	modified residue	GB 24371665
A	157	MSE	MET	modified residue	GB 24371665
B	1	MSE	MET	modified residue	GB 24371665
B	61	MSE	MET	modified residue	GB 24371665
B	95	MSE	MET	modified residue	GB 24371665
B	96	MSE	MET	modified residue	GB 24371665
B	103	MSE	MET	modified residue	GB 24371665
B	157	MSE	MET	modified residue	GB 24371665
C	1	MSE	MET	modified residue	GB 24371665
C	61	MSE	MET	modified residue	GB 24371665
C	95	MSE	MET	modified residue	GB 24371665
C	96	MSE	MET	modified residue	GB 24371665
C	103	MSE	MET	modified residue	GB 24371665
C	157	MSE	MET	modified residue	GB 24371665

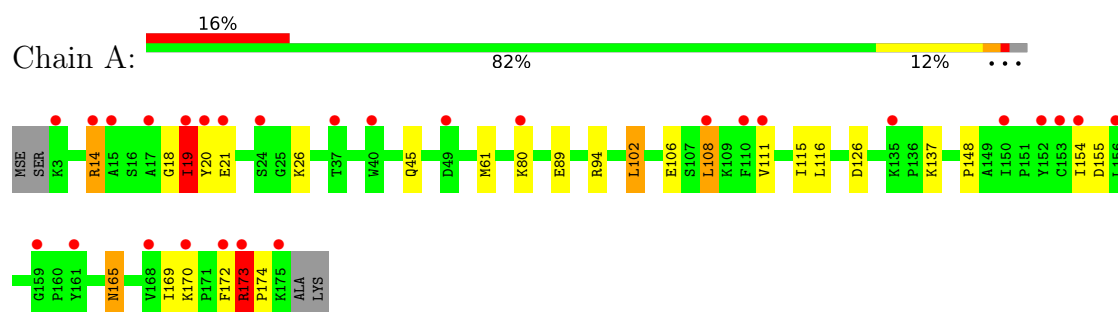
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	52	Total 52	O 52	0	0
2	B	21	Total 21	O 21	0	0
2	C	49	Total 49	O 49	0	0

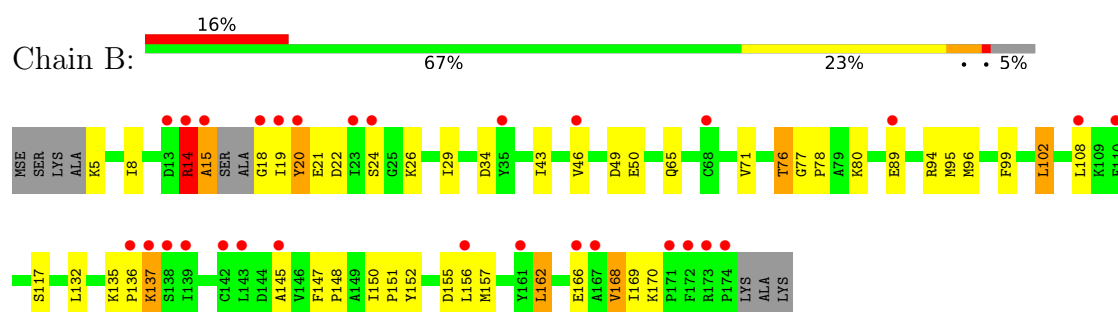
3 Residue-property plots [i](#)

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 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.

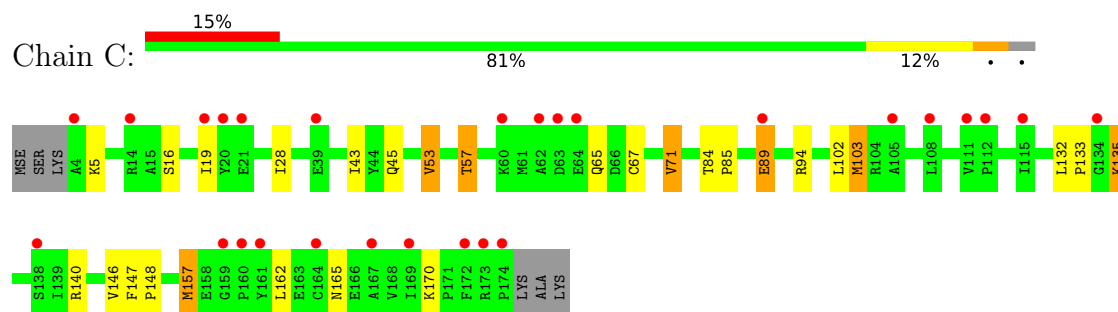
- Molecule 1: molybdenum cofactor biosynthesis protein Mog



- Molecule 1: molybdenum cofactor biosynthesis protein Mog



- Molecule 1: molybdenum cofactor biosynthesis protein Mog



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	85.57Å 215.36Å 40.84Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 2.30 50.00 – 2.30	Depositor EDS
% Data completeness (in resolution range)	98.6 (50.00-2.30) 98.5 (50.00-2.30)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.74 (at 2.29Å)	Xtriage
Refinement program	REFMAC 5.2.0005	Depositor
R, R_{free}	0.193 , 0.222 (Not available) , 0.283	Depositor DCC
R_{free} test set	1717 reflections (5.03%)	wwPDB-VP
Wilson B-factor (Å ²)	37.6	Xtriage
Anisotropy	0.331	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 57.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	4026	wwPDB-VP
Average B, all atoms (Å ²)	44.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.00% 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	1.25	2/1334 (0.1%)	1.27	9/1805 (0.5%)
1	B	1.40	11/1303 (0.8%)	1.16	6/1762 (0.3%)
1	C	1.17	3/1320 (0.2%)	1.18	5/1787 (0.3%)
All	All	1.28	16/3957 (0.4%)	1.21	20/5354 (0.4%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1
1	B	0	2
All	All	0	3

All (16) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	15	ALA	C-O	27.47	1.78	1.23
1	B	21	GLU	CD-OE1	9.91	1.44	1.25
1	B	21	GLU	CD-OE2	9.55	1.43	1.25
1	B	24	SER	CB-OG	7.98	1.58	1.42
1	B	20	TYR	CG-CD2	7.73	1.55	1.39
1	B	18	GLY	N-CA	7.61	1.57	1.45
1	C	103	MSE	SE-CE	7.43	2.17	1.95
1	B	22	ASP	C-N	7.42	1.43	1.34
1	B	21	GLU	C-N	6.55	1.42	1.33
1	A	154	ILE	CA-CB	6.28	1.62	1.54
1	B	22	ASP	C-O	5.91	1.30	1.23
1	C	89	GLU	CG-CD	5.59	1.66	1.52
1	B	21	GLU	C-O	5.51	1.30	1.23
1	B	20	TYR	CE1-CZ	5.14	1.50	1.38
1	C	19	ILE	CA-CB	5.05	1.61	1.54

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	126	ASP	C-O	-5.01	1.17	1.24

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	19	ILE	N-CA-C	15.34	125.70	110.82
1	A	18	GLY	N-CA-C	9.76	124.02	113.58
1	B	15	ALA	CA-C-O	-7.69	107.73	120.80
1	A	14	ARG	N-CA-C	7.42	120.24	111.71
1	A	173	ARG	NE-CZ-NH1	-7.17	114.33	121.50
1	B	14	ARG	CA-C-N	6.58	133.55	121.70
1	B	14	ARG	C-N-CA	6.58	133.55	121.70
1	A	19	ILE	CA-C-N	6.09	131.32	122.36
1	A	19	ILE	C-N-CA	6.09	131.32	122.36
1	C	133	PRO	CA-C-N	5.91	126.50	120.00
1	C	133	PRO	C-N-CA	5.91	126.50	120.00
1	C	53	VAL	N-CA-C	5.68	115.87	110.42
1	B	77	GLY	CA-C-N	5.43	125.13	119.64
1	B	77	GLY	C-N-CA	5.43	125.13	119.64
1	A	173	ARG	NE-CZ-NH2	5.33	124.00	119.20
1	C	132	LEU	CA-C-N	-5.33	115.28	120.98
1	C	132	LEU	C-N-CA	-5.33	115.28	120.98
1	A	111	VAL	CA-C-N	5.19	126.33	119.84
1	A	111	VAL	C-N-CA	5.19	126.33	119.84
1	B	78	PRO	N-CA-C	5.16	121.35	114.18

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	19	ILE	Peptide
1	B	14	ARG	Peptide
1	B	20	TYR	Sidechain

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	1316	0	1324	14	0
1	B	1286	0	1293	32	0
1	C	1302	0	1309	23	0
2	A	52	0	0	1	0
2	B	21	0	0	2	0
2	C	49	0	0	6	0
All	All	4026	0	3926	67	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 9.

All (67) 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:103:MSE:SE	1:C:103:MSE:CE	2.17	1.41
1:B:15:ALA:O	1:B:15:ALA:C	1.78	1.25
1:C:16:SER:HB3	2:C:184:HOH:O	1.42	1.20
1:A:173:ARG:HG2	1:A:173:ARG:HH11	1.13	1.07
1:B:14:ARG:HH11	1:B:14:ARG:HG3	1.21	1.02
1:A:173:ARG:HH11	1:A:173:ARG:CG	1.82	0.88
1:C:67:CYS:SG	2:C:206:HOH:O	2.33	0.87
1:A:155:ASP:OD2	1:A:173:ARG:NH1	2.07	0.86
1:A:173:ARG:HG2	1:A:173:ARG:NH1	1.84	0.86
2:A:182:HOH:O	1:B:94:ARG:HD2	1.81	0.80
1:C:53:VAL:O	1:C:57:THR:CG2	2.30	0.80
1:C:53:VAL:O	1:C:57:THR:HG23	1.83	0.79
1:B:14:ARG:HH11	1:B:14:ARG:CG	2.00	0.71
1:C:5:LYS:HE2	1:C:65:GLN:O	1.93	0.68
1:B:14:ARG:HG3	1:B:14:ARG:NH1	1.99	0.64
1:B:162:LEU:H	1:B:162:LEU:HD12	1.63	0.63
1:B:135:LYS:HD3	1:B:136:PRO:HD2	1.81	0.62
1:A:173:ARG:CG	1:A:173:ARG:NH1	2.50	0.61
1:B:135:LYS:HD3	1:B:136:PRO:N	2.16	0.61
1:C:94:ARG:NH2	1:C:157:MSE:HE1	2.16	0.60
1:B:135:LYS:HD3	1:B:136:PRO:CD	2.32	0.60
1:C:89:GLU:CG	2:C:216:HOH:O	2.52	0.58
1:C:28:ILE:HG21	1:C:71:VAL:HG21	1.85	0.57
1:C:28:ILE:HD13	1:C:71:VAL:HG23	1.86	0.57
1:C:53:VAL:O	1:C:57:THR:HG22	2.05	0.57
1:B:5:LYS:HE2	1:B:65:GLN:O	2.05	0.56
1:B:102:LEU:HD13	1:B:145:ALA:HB1	1.88	0.56
1:B:135:LYS:HD2	1:B:137:LYS:H	1.71	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:155:ASP:CG	1:A:173:ARG:HH12	2.09	0.55
1:B:96:MSE:HE3	1:B:99:PHE:HE1	1.73	0.54
1:A:165:ASN:HD22	1:A:165:ASN:C	2.16	0.53
1:C:28:ILE:HG21	1:C:71:VAL:CG2	2.40	0.52
1:B:135:LYS:HD3	1:B:135:LYS:C	2.34	0.51
1:B:71:VAL:HG11	1:B:132:LEU:CD1	2.41	0.51
1:B:34:ASP:O	1:B:168:VAL:HG21	2.10	0.51
1:A:148:PRO:HD3	1:A:169:ILE:HD11	1.93	0.51
1:B:49:ASP:CB	2:B:182:HOH:O	2.59	0.50
1:A:172:PHE:CE2	1:A:174:PRO:HG3	2.47	0.50
1:C:135:LYS:HB2	2:C:192:HOH:O	2.11	0.49
1:A:19:ILE:N	1:A:20:TYR:HB2	2.28	0.49
1:B:5:LYS:HG2	1:B:43:ILE:HD11	1.95	0.49
1:C:28:ILE:CG2	1:C:71:VAL:HG21	2.43	0.49
1:C:94:ARG:HH22	1:C:157:MSE:HE1	1.79	0.48
1:B:49:ASP:HB3	2:B:182:HOH:O	2.12	0.48
1:B:147:PHE:HA	1:B:150:ILE:HD12	1.97	0.47
1:A:116:LEU:HD22	1:B:96:MSE:HE1	1.97	0.46
1:A:102:LEU:HD22	1:A:106:GLU:HG2	1.97	0.46
1:B:148:PRO:HD3	1:B:169:ILE:HD11	1.97	0.46
1:B:151:PRO:HA	1:B:162:LEU:HD13	1.97	0.46
1:C:165:ASN:C	1:C:165:ASN:HD22	2.23	0.45
1:A:45:GLN:HG3	1:A:61:MSE:HE3	1.99	0.45
1:C:5:LYS:C	2:C:206:HOH:O	2.59	0.45
1:C:89:GLU:HG2	2:C:216:HOH:O	2.15	0.45
1:C:28:ILE:HD13	1:C:71:VAL:CG2	2.48	0.44
1:B:152:TYR:O	1:B:155:ASP:HB2	2.18	0.44
1:B:89:GLU:HG2	1:B:95:MSE:HE2	1.99	0.44
1:B:162:LEU:HD12	1:B:162:LEU:N	2.33	0.43
1:A:108:LEU:HD11	1:A:115:ILE:CG2	2.49	0.43
1:B:80:LYS:HD2	1:C:94:ARG:CZ	2.48	0.43
1:B:76:THR:CG2	1:B:117:SER:OG	2.67	0.42
1:B:156:LEU:HB2	1:B:157:MSE:HE2	2.01	0.42
1:C:147:PHE:N	1:C:148:PRO:CD	2.83	0.41
1:C:157:MSE:H	1:C:157:MSE:HG2	1.75	0.41
1:B:108:LEU:C	1:B:108:LEU:HD12	2.45	0.41
1:B:8:ILE:HD13	1:B:29:ILE:HG13	2.03	0.40
1:C:84:THR:N	1:C:85:PRO:CD	2.84	0.40
1:B:14:ARG:HB3	1:B:15:ALA:HB3	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	171/177 (97%)	167 (98%)	4 (2%)	0	100	100
1	B	164/177 (93%)	161 (98%)	3 (2%)	0	100	100
1	C	169/177 (96%)	163 (96%)	6 (4%)	0	100	100
All	All	504/531 (95%)	491 (97%)	13 (3%)	0	100	100

There are no Ramachandran outliers to report.

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	146/144 (101%)	134 (92%)	12 (8%)	10	14
1	B	144/144 (100%)	132 (92%)	12 (8%)	10	14
1	C	145/144 (101%)	134 (92%)	11 (8%)	12	16
All	All	435/432 (101%)	400 (92%)	35 (8%)	11	15

All (35) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	14	ARG
1	A	21	GLU
1	A	26	LYS
1	A	80	LYS
1	A	89	GLU

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Mol	Chain	Res	Type
1	A	94	ARG
1	A	102	LEU
1	A	108	LEU
1	A	137	LYS
1	A	165	ASN
1	A	170	LYS
1	A	173	ARG
1	B	14	ARG
1	B	19	ILE
1	B	26	LYS
1	B	46	VAL
1	B	50	GLU
1	B	76	THR
1	B	102	LEU
1	B	137	LYS
1	B	162	LEU
1	B	166	GLU
1	B	168	VAL
1	B	170	LYS
1	C	43	ILE
1	C	45	GLN
1	C	57	THR
1	C	71	VAL
1	C	102	LEU
1	C	135	LYS
1	C	140	ARG
1	C	146	VAL
1	C	157	MSE
1	C	162	LEU
1	C	170	LYS

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

Mol	Chain	Res	Type
1	A	65	GLN
1	A	165	ASN
1	B	165	ASN
1	C	165	ASN

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 [i](#)

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	168/177 (94%)	1.15	28 (16%) 4 5	35, 45, 59, 64	0
1	B	163/177 (92%)	1.15	29 (17%) 4 4	31, 41, 58, 68	0
1	C	166/177 (93%)	1.30	27 (16%) 4 5	29, 43, 57, 61	0
All	All	497/531 (93%)	1.20	84 (16%) 4 5	29, 43, 58, 68	0

All (84) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	20	TYR	6.4
1	B	15	ALA	5.9
1	B	20	TYR	5.9
1	B	19	ILE	5.6
1	A	19	ILE	5.1
1	B	174	PRO	4.6
1	C	20	TYR	4.4
1	A	152	TYR	4.2
1	B	137	LYS	4.1
1	B	138	SER	4.0
1	B	139	ILE	4.0
1	B	18	GLY	3.8
1	A	154	ILE	3.7
1	B	172	PHE	3.7
1	C	19	ILE	3.6
1	B	24	SER	3.4
1	A	156	LEU	3.4
1	C	39	GLU	3.4
1	C	174	PRO	3.4
1	B	142	CYS	3.4
1	C	115	ILE	3.3
1	A	17	ALA	3.2
1	A	14	ARG	3.2

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Mol	Chain	Res	Type	RSRZ
1	C	108	LEU	3.1
1	B	35	TYR	3.1
1	A	173	ARG	3.1
1	C	172	PHE	3.0
1	B	166	GLU	3.0
1	B	23	ILE	3.0
1	A	172	PHE	2.9
1	B	167	ALA	2.9
1	B	136	PRO	2.9
1	A	153	CYS	2.9
1	C	161	TYR	2.9
1	C	111	VAL	2.9
1	C	89	GLU	2.7
1	C	138	SER	2.7
1	A	175	LYS	2.7
1	A	161	TYR	2.7
1	B	161	TYR	2.7
1	B	68	CYS	2.7
1	B	143	LEU	2.6
1	A	110	PHE	2.6
1	C	112	PRO	2.6
1	C	164	CYS	2.6
1	C	4	ALA	2.6
1	C	160	PRO	2.5
1	A	111	VAL	2.5
1	B	156	LEU	2.5
1	C	167	ALA	2.5
1	A	3	LYS	2.5
1	A	150	ILE	2.4
1	C	105	ALA	2.4
1	A	40	TRP	2.4
1	A	108	LEU	2.4
1	B	14	ARG	2.4
1	C	21	GLU	2.4
1	A	15	ALA	2.3
1	A	37	THR	2.3
1	B	108	LEU	2.3
1	A	21	GLU	2.3
1	B	89	GLU	2.3
1	A	159	GLY	2.3
1	C	63	ASP	2.3
1	B	46	VAL	2.3

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Mol	Chain	Res	Type	RSRZ
1	A	135	LYS	2.3
1	B	13	ASP	2.3
1	B	171	PRO	2.3
1	A	24	SER	2.3
1	C	64	GLU	2.3
1	C	14	ARG	2.2
1	C	134	GLY	2.2
1	B	145	ALA	2.2
1	B	173	ARG	2.2
1	B	110	PHE	2.2
1	C	60	LYS	2.2
1	C	169	ILE	2.1
1	A	168	VAL	2.1
1	A	80	LYS	2.1
1	C	62	ALA	2.1
1	C	173	ARG	2.0
1	C	159	GLY	2.0
1	A	49	ASP	2.0
1	A	170	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.