Full wwPDB X-ray Structure Validation Report ⓘ

May 27, 2020 – 07:37 pm BST

PDB ID	:	2F7Y
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 following versions of software and data (see references ⓘ) were used in the production of this report:

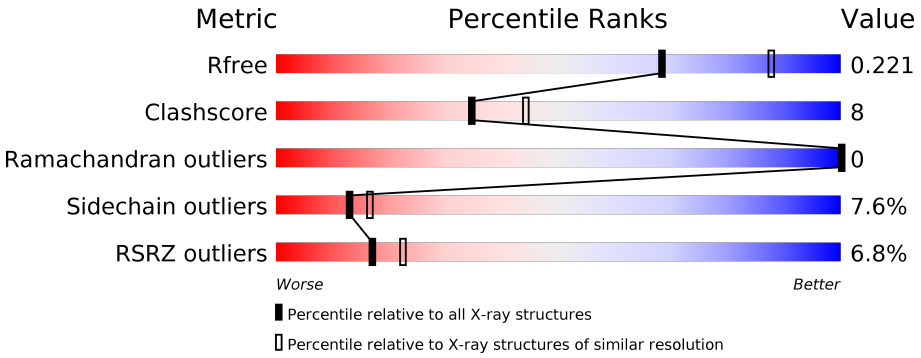
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.11
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.11

1 Overall quality at a glance (i)

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}	130704	5042 (2.30-2.30)
Clashscore	141614	5643 (2.30-2.30)
Ramachandran outliers	138981	5575 (2.30-2.30)
Sidechain outliers	138945	5575 (2.30-2.30)
RSRZ outliers	127900	4938 (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 on 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 <=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	5%, 82%, 11%
1	B	177	8%, 69%, 20%, 5% • 5%
1	C	177	6%, 83%, 10%

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.

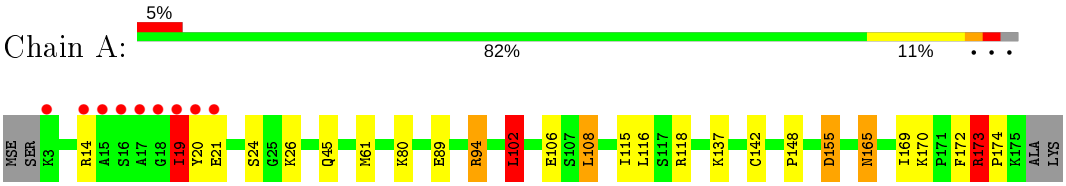
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	52	Total O 52 52	0	0
2	B	21	Total O 21 21	0	0
2	C	49	Total O 49 49	0	0

3 Residue-property plots (i)

These plots are drawn for all protein, RNA and DNA 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.

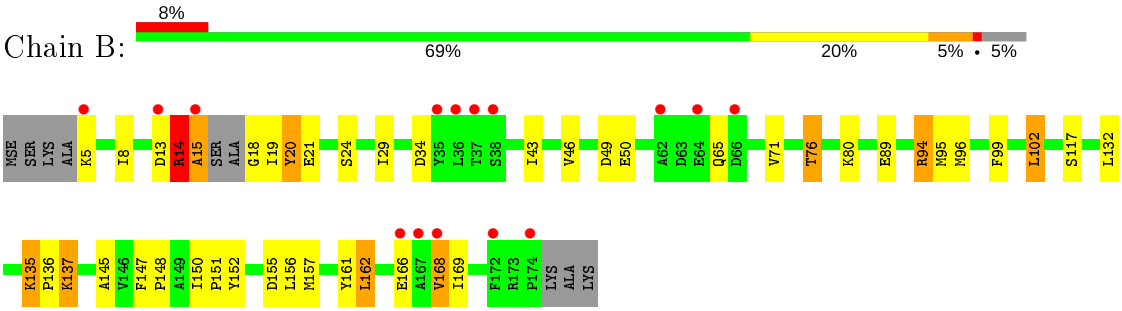
- Molecule 1: molybdenum cofactor biosynthesis protein Mog

Chain A:



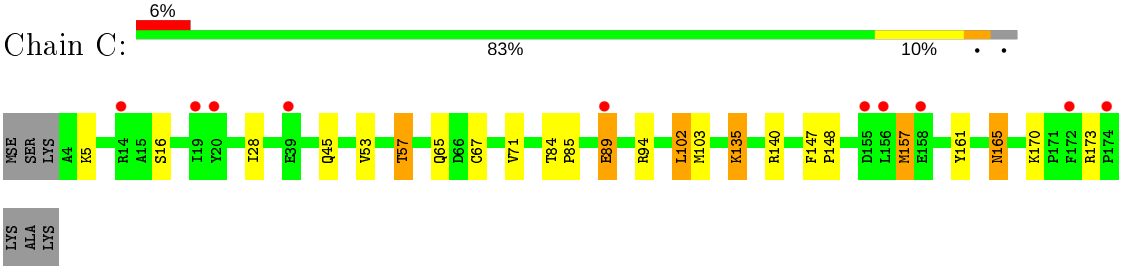
- Molecule 1: molybdenum cofactor biosynthesis protein Mog

Chain B:



- Molecule 1: molybdenum cofactor biosynthesis protein Mog

Chain C:



4 Data and refinement statistics (i)

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 36.86 – 2.30	Depositor EDS
% Data completeness (in resolution range)	98.6 (50.00-2.30) 98.6 (36.86-2.30)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	3.74 (at 2.29Å)	Xtriage
Refinement program	REFMAC 5.2.0005	Depositor
R, R_{free}	0.193 , 0.222 0.191 , 0.221	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)	0.35 , 58.0	EDS
L-test for twinning[2]	$< \|L\| > = 0.47$, $< L^2 > = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.94	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.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ 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	0.97	1/1334 (0.1%)	1.05	10/1805 (0.6%)
1	B	1.50	12/1303 (0.9%)	0.96	4/1762 (0.2%)
1	C	0.96	4/1320 (0.3%)	0.91	3/1787 (0.2%)
All	All	1.17	17/3957 (0.4%)	0.98	17/5354 (0.3%)

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 (17) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	15	ALA	C-O	29.03	1.78	1.23
1	B	21	GLU	CD-OE1	16.85	1.44	1.25
1	B	21	GLU	CD-OE2	16.23	1.43	1.25
1	B	20	TYR	CG-CD2	12.64	1.55	1.39
1	B	24	SER	CB-OG	12.21	1.58	1.42
1	C	89	GLU	CG-CD	9.39	1.66	1.51
1	B	20	TYR	CE1-CZ	9.27	1.50	1.38
1	B	20	TYR	CE2-CZ	7.93	1.48	1.38
1	B	161	TYR	CG-CD1	7.91	1.49	1.39
1	B	161	TYR	CE2-CZ	7.80	1.48	1.38
1	B	18	GLY	N-CA	7.78	1.57	1.46
1	C	161	TYR	CG-CD2	7.08	1.48	1.39
1	A	142	CYS	CB-SG	6.64	1.93	1.82
1	C	161	TYR	CE1-CZ	6.23	1.46	1.38
1	B	161	TYR	CE1-CZ	5.93	1.46	1.38

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	161	TYR	CG-CD2	5.80	1.46	1.39
1	C	89	GLU	CD-OE2	5.14	1.31	1.25

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	20	TYR	CB-CG-CD2	-12.77	113.34	121.00
1	A	173	ARG	NE-CZ-NH1	-11.93	114.33	120.30
1	A	118	ARG	NE-CZ-NH2	-9.35	115.63	120.30
1	B	20	TYR	CG-CD2-CE2	-9.09	114.03	121.30
1	A	155	ASP	CB-CG-OD2	-7.73	111.34	118.30
1	A	173	ARG	NE-CZ-NH2	7.39	124.00	120.30
1	C	173	ARG	NE-CZ-NH2	-6.87	116.86	120.30
1	A	155	ASP	CB-CG-OD1	6.79	124.42	118.30
1	A	102	LEU	CB-CG-CD2	6.72	122.42	111.00
1	A	118	ARG	NE-CZ-NH1	6.28	123.44	120.30
1	B	15	ALA	CA-C-O	-5.89	107.73	120.10
1	C	102	LEU	CB-CG-CD2	5.59	120.51	111.00
1	A	19	ILE	N-CA-C	5.44	125.70	111.00
1	A	94	ARG	NE-CZ-NH2	-5.21	117.69	120.30
1	B	21	GLU	OE1-CD-OE2	5.10	129.43	123.30
1	C	173	ARG	NE-CZ-NH1	5.08	122.84	120.30
1	A	24	SER	CB-CA-C	5.06	119.71	110.10

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 (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	1316	0	1324	14	0

Continued on next page...

Continued from previous page...

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	B	1286	0	1293	31	0
1	C	1302	0	1309	22	0
2	A	52	0	0	1	0
2	B	21	0	0	2	0
2	C	49	0	0	5	0
All	All	4026	0	3926	65	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 (65) 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:CE	1:C:103:MSE:SE	2.17	1.41
1:B:15:ALA:C	1:B:15:ALA:O	1.78	1.20
1:C:16:SER:HB3	2:C:184:HOH:O	1.42	1.20
1:A:173:ARG:HH11	1:A:173:ARG:HG2	1.13	1.07
1:B:14:ARG:HG3	1:B:14:ARG:HH11	1.21	1.02
1:A:173:ARG:CG	1:A:173:ARG:HH11	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:NH1	1:A:173:ARG:HG2	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:CG	1:B:14:ARG:HH11	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:HD12	1:B:162:LEU:H	1.63	0.63
1:B:135:LYS:HD3	1:B:136:PRO:HD2	1.81	0.62
1:A:173:ARG:NH1	1:A:173:ARG:CG	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:53:VAL:O	1:C:57:THR:HG22	2.05	0.57
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:B:5:LYS:HE2	1:B:65:GLN:O	2.05	0.56
1:B:135:LYS:HD2	1:B:137:LYS:H	1.71	0.56
1:B:102:LEU:HD13	1:B:145:ALA:HB1	1.88	0.56

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:96:MSE:HE3	1:B:99:PHE:HE1	1.73	0.54
1:C:28:ILE:HG21	1:C:71:VAL:CG2	2.40	0.52
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:A:155:ASP:CG	1:A:173:ARG:HH12	2.09	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:A:165:ASN:HD22	1:A:165:ASN:C	2.16	0.48
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:B:135:LYS:HD3	1:B:135:LYS:C	2.34	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:A:45:GLN:HG3	1:A:61:MSE:HE3	1.99	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:N	1:B:162:LEU:HD12	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:C:165:ASN:C	1:C:165:ASN:HD22	2.23	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 (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 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 (i)

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%)	11	14
1	B	144/144 (100%)	131 (91%)	13 (9%)	9	11
1	C	145/144 (101%)	137 (94%)	8 (6%)	21	30
All	All	435/432 (101%)	402 (92%)	33 (8%)	13	16

All (33) 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

Continued on next page...

Continued from previous page...

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	13	ASP
1	B	14	ARG
1	B	19	ILE
1	B	46	VAL
1	B	50	GLU
1	B	76	THR
1	B	94	ARG
1	B	102	LEU
1	B	135	LYS
1	B	137	LYS
1	B	162	LEU
1	B	166	GLU
1	B	168	VAL
1	C	45	GLN
1	C	57	THR
1	C	102	LEU
1	C	135	LYS
1	C	140	ARG
1	C	157	MSE
1	C	165	ASN
1	C	170	LYS

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

Mol	Chain	Res	Type
1	A	165	ASN
1	B	119	GLN
1	B	165	ASN
1	C	165	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

There are no carbohydrates 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, 95^{th} 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%)	0.42	9 (5%) 25 32	35, 45, 59, 64	0
1	B	163/177 (92%)	0.59	15 (9%) 9 12	31, 41, 58, 68	0
1	C	166/177 (93%)	0.49	10 (6%) 21 28	29, 43, 57, 61	0
All	All	497/531 (93%)	0.50	34 (6%) 17 22	29, 43, 58, 68	0

All (34) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	20	TYR	10.4
1	A	19	ILE	6.8
1	A	17	ALA	6.4
1	C	19	ILE	6.2
1	C	20	TYR	5.1
1	B	167	ALA	4.5
1	B	166	GLU	4.2
1	B	66	ASP	3.8
1	C	39	GLU	3.8
1	C	174	PRO	3.7
1	B	172	PHE	3.4
1	B	5	LYS	3.4
1	B	15	ALA	3.3
1	C	172	PHE	3.3
1	B	36	LEU	3.3
1	B	35	TYR	3.2
1	B	174	PRO	3.1
1	A	18	GLY	3.1
1	A	14	ARG	3.0
1	B	168	VAL	3.0
1	B	64	GLU	2.9
1	A	21	GLU	2.8
1	B	38	SER	2.7

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	C	14	ARG	2.6
1	C	155	ASP	2.5
1	A	15	ALA	2.5
1	C	158	GLU	2.5
1	A	16	SER	2.5
1	C	89	GLU	2.5
1	B	62	ALA	2.4
1	B	13	ASP	2.3
1	A	3	LYS	2.3
1	B	37	THR	2.2
1	C	156	LEU	2.1

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 carbohydrates 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.