



Full wwPDB X-ray Structure Validation Report ⓘ

Nov 2, 2021 – 03:05 AM EDT

PDB ID : 1ZYE
Title : Crystal structure analysis of Bovine Mitochondrial Peroxiredoxin III
Authors : Cao, Z.; Roszak, A.W.; Gourlay, L.J.; Lindsay, J.G.; Isaacs, N.W.
Deposited on : 2005-06-10
Resolution : 3.30 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
Xtriage (Phenix) : 1.13
EDS : 2.23.2
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) : 2.23.2

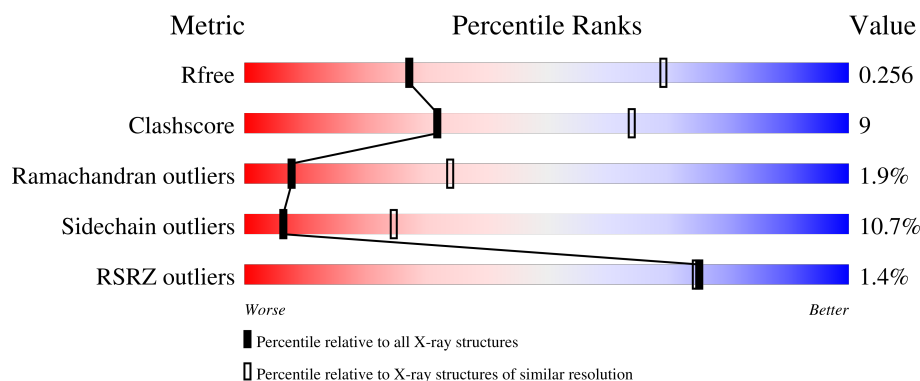
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.30 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	1149 (3.34-3.26)
Clashscore	141614	1205 (3.34-3.26)
Ramachandran outliers	138981	1183 (3.34-3.26)
Sidechain outliers	138945	1182 (3.34-3.26)
RSRZ outliers	127900	1115 (3.34-3.26)

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	220	
1	B	220	
1	C	220	
1	D	220	
1	E	220	

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Mol	Chain	Length	Quality of chain
1	F	220	
1	G	220	
1	H	220	
1	I	220	
1	J	220	
1	K	220	
1	L	220	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 15201 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 Thioredoxin-dependent peroxide reductase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	162	Total	C	N	O	S	0	0	0
			1266	824	208	231	3			
1	B	162	Total	C	N	O	S	0	0	0
			1266	824	208	231	3			
1	C	162	Total	C	N	O	S	0	0	0
			1266	824	208	231	3			
1	D	162	Total	C	N	O	S	0	0	0
			1266	824	208	231	3			
1	E	162	Total	C	N	O	S	0	0	0
			1266	824	208	231	3			
1	F	162	Total	C	N	O	S	0	0	0
			1266	824	208	231	3			
1	G	162	Total	C	N	O	S	0	0	0
			1266	824	208	231	3			
1	H	162	Total	C	N	O	S	0	0	0
			1266	824	208	231	3			
1	I	162	Total	C	N	O	S	0	0	0
			1266	824	208	231	3			
1	J	162	Total	C	N	O	S	0	0	0
			1266	824	208	231	3			
1	K	162	Total	C	N	O	S	0	0	0
			1266	824	208	231	3			
1	L	162	Total	C	N	O	S	0	0	0
			1266	824	208	231	3			

There are 312 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-24	MET	-	expression tag	UNP P35705
A	-23	GLY	-	expression tag	UNP P35705
A	-22	SER	-	expression tag	UNP P35705
A	-21	SER	-	expression tag	UNP P35705
A	-20	HIS	-	expression tag	UNP P35705

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Chain	Residue	Modelled	Actual	Comment	Reference
A	-19	HIS	-	expression tag	UNP P35705
A	-18	HIS	-	expression tag	UNP P35705
A	-17	HIS	-	expression tag	UNP P35705
A	-16	HIS	-	expression tag	UNP P35705
A	-15	HIS	-	expression tag	UNP P35705
A	-14	SER	-	expression tag	UNP P35705
A	-13	SER	-	expression tag	UNP P35705
A	-12	GLY	-	expression tag	UNP P35705
A	-11	LEU	-	expression tag	UNP P35705
A	-10	VAL	-	expression tag	UNP P35705
A	-9	PRO	-	expression tag	UNP P35705
A	-8	ARG	-	expression tag	UNP P35705
A	-7	GLY	-	expression tag	UNP P35705
A	-6	SER	-	expression tag	UNP P35705
A	-5	HIS	-	expression tag	UNP P35705
A	-4	MET	-	expression tag	UNP P35705
A	-3	LEU	-	expression tag	UNP P35705
A	-2	GLU	-	expression tag	UNP P35705
A	-1	ASP	-	expression tag	UNP P35705
A	0	PRO	-	expression tag	UNP P35705
A	168	SER	CYS	engineered mutation	UNP P35705
B	-24	MET	-	expression tag	UNP P35705
B	-23	GLY	-	expression tag	UNP P35705
B	-22	SER	-	expression tag	UNP P35705
B	-21	SER	-	expression tag	UNP P35705
B	-20	HIS	-	expression tag	UNP P35705
B	-19	HIS	-	expression tag	UNP P35705
B	-18	HIS	-	expression tag	UNP P35705
B	-17	HIS	-	expression tag	UNP P35705
B	-16	HIS	-	expression tag	UNP P35705
B	-15	HIS	-	expression tag	UNP P35705
B	-14	SER	-	expression tag	UNP P35705
B	-13	SER	-	expression tag	UNP P35705
B	-12	GLY	-	expression tag	UNP P35705
B	-11	LEU	-	expression tag	UNP P35705
B	-10	VAL	-	expression tag	UNP P35705
B	-9	PRO	-	expression tag	UNP P35705
B	-8	ARG	-	expression tag	UNP P35705
B	-7	GLY	-	expression tag	UNP P35705
B	-6	SER	-	expression tag	UNP P35705
B	-5	HIS	-	expression tag	UNP P35705
B	-4	MET	-	expression tag	UNP P35705

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-3	LEU	-	expression tag	UNP P35705
B	-2	GLU	-	expression tag	UNP P35705
B	-1	ASP	-	expression tag	UNP P35705
B	0	PRO	-	expression tag	UNP P35705
B	168	SER	CYS	engineered mutation	UNP P35705
C	-24	MET	-	expression tag	UNP P35705
C	-23	GLY	-	expression tag	UNP P35705
C	-22	SER	-	expression tag	UNP P35705
C	-21	SER	-	expression tag	UNP P35705
C	-20	HIS	-	expression tag	UNP P35705
C	-19	HIS	-	expression tag	UNP P35705
C	-18	HIS	-	expression tag	UNP P35705
C	-17	HIS	-	expression tag	UNP P35705
C	-16	HIS	-	expression tag	UNP P35705
C	-15	HIS	-	expression tag	UNP P35705
C	-14	SER	-	expression tag	UNP P35705
C	-13	SER	-	expression tag	UNP P35705
C	-12	GLY	-	expression tag	UNP P35705
C	-11	LEU	-	expression tag	UNP P35705
C	-10	VAL	-	expression tag	UNP P35705
C	-9	PRO	-	expression tag	UNP P35705
C	-8	ARG	-	expression tag	UNP P35705
C	-7	GLY	-	expression tag	UNP P35705
C	-6	SER	-	expression tag	UNP P35705
C	-5	HIS	-	expression tag	UNP P35705
C	-4	MET	-	expression tag	UNP P35705
C	-3	LEU	-	expression tag	UNP P35705
C	-2	GLU	-	expression tag	UNP P35705
C	-1	ASP	-	expression tag	UNP P35705
C	0	PRO	-	expression tag	UNP P35705
C	168	SER	CYS	engineered mutation	UNP P35705
D	-24	MET	-	expression tag	UNP P35705
D	-23	GLY	-	expression tag	UNP P35705
D	-22	SER	-	expression tag	UNP P35705
D	-21	SER	-	expression tag	UNP P35705
D	-20	HIS	-	expression tag	UNP P35705
D	-19	HIS	-	expression tag	UNP P35705
D	-18	HIS	-	expression tag	UNP P35705
D	-17	HIS	-	expression tag	UNP P35705
D	-16	HIS	-	expression tag	UNP P35705
D	-15	HIS	-	expression tag	UNP P35705
D	-14	SER	-	expression tag	UNP P35705

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Chain	Residue	Modelled	Actual	Comment	Reference
D	-13	SER	-	expression tag	UNP P35705
D	-12	GLY	-	expression tag	UNP P35705
D	-11	LEU	-	expression tag	UNP P35705
D	-10	VAL	-	expression tag	UNP P35705
D	-9	PRO	-	expression tag	UNP P35705
D	-8	ARG	-	expression tag	UNP P35705
D	-7	GLY	-	expression tag	UNP P35705
D	-6	SER	-	expression tag	UNP P35705
D	-5	HIS	-	expression tag	UNP P35705
D	-4	MET	-	expression tag	UNP P35705
D	-3	LEU	-	expression tag	UNP P35705
D	-2	GLU	-	expression tag	UNP P35705
D	-1	ASP	-	expression tag	UNP P35705
D	0	PRO	-	expression tag	UNP P35705
D	168	SER	CYS	engineered mutation	UNP P35705
E	-24	MET	-	expression tag	UNP P35705
E	-23	GLY	-	expression tag	UNP P35705
E	-22	SER	-	expression tag	UNP P35705
E	-21	SER	-	expression tag	UNP P35705
E	-20	HIS	-	expression tag	UNP P35705
E	-19	HIS	-	expression tag	UNP P35705
E	-18	HIS	-	expression tag	UNP P35705
E	-17	HIS	-	expression tag	UNP P35705
E	-16	HIS	-	expression tag	UNP P35705
E	-15	HIS	-	expression tag	UNP P35705
E	-14	SER	-	expression tag	UNP P35705
E	-13	SER	-	expression tag	UNP P35705
E	-12	GLY	-	expression tag	UNP P35705
E	-11	LEU	-	expression tag	UNP P35705
E	-10	VAL	-	expression tag	UNP P35705
E	-9	PRO	-	expression tag	UNP P35705
E	-8	ARG	-	expression tag	UNP P35705
E	-7	GLY	-	expression tag	UNP P35705
E	-6	SER	-	expression tag	UNP P35705
E	-5	HIS	-	expression tag	UNP P35705
E	-4	MET	-	expression tag	UNP P35705
E	-3	LEU	-	expression tag	UNP P35705
E	-2	GLU	-	expression tag	UNP P35705
E	-1	ASP	-	expression tag	UNP P35705
E	0	PRO	-	expression tag	UNP P35705
E	168	SER	CYS	engineered mutation	UNP P35705
F	-24	MET	-	expression tag	UNP P35705

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Chain	Residue	Modelled	Actual	Comment	Reference
F	-23	GLY	-	expression tag	UNP P35705
F	-22	SER	-	expression tag	UNP P35705
F	-21	SER	-	expression tag	UNP P35705
F	-20	HIS	-	expression tag	UNP P35705
F	-19	HIS	-	expression tag	UNP P35705
F	-18	HIS	-	expression tag	UNP P35705
F	-17	HIS	-	expression tag	UNP P35705
F	-16	HIS	-	expression tag	UNP P35705
F	-15	HIS	-	expression tag	UNP P35705
F	-14	SER	-	expression tag	UNP P35705
F	-13	SER	-	expression tag	UNP P35705
F	-12	GLY	-	expression tag	UNP P35705
F	-11	LEU	-	expression tag	UNP P35705
F	-10	VAL	-	expression tag	UNP P35705
F	-9	PRO	-	expression tag	UNP P35705
F	-8	ARG	-	expression tag	UNP P35705
F	-7	GLY	-	expression tag	UNP P35705
F	-6	SER	-	expression tag	UNP P35705
F	-5	HIS	-	expression tag	UNP P35705
F	-4	MET	-	expression tag	UNP P35705
F	-3	LEU	-	expression tag	UNP P35705
F	-2	GLU	-	expression tag	UNP P35705
F	-1	ASP	-	expression tag	UNP P35705
F	0	PRO	-	expression tag	UNP P35705
F	168	SER	CYS	engineered mutation	UNP P35705
G	-24	MET	-	expression tag	UNP P35705
G	-23	GLY	-	expression tag	UNP P35705
G	-22	SER	-	expression tag	UNP P35705
G	-21	SER	-	expression tag	UNP P35705
G	-20	HIS	-	expression tag	UNP P35705
G	-19	HIS	-	expression tag	UNP P35705
G	-18	HIS	-	expression tag	UNP P35705
G	-17	HIS	-	expression tag	UNP P35705
G	-16	HIS	-	expression tag	UNP P35705
G	-15	HIS	-	expression tag	UNP P35705
G	-14	SER	-	expression tag	UNP P35705
G	-13	SER	-	expression tag	UNP P35705
G	-12	GLY	-	expression tag	UNP P35705
G	-11	LEU	-	expression tag	UNP P35705
G	-10	VAL	-	expression tag	UNP P35705
G	-9	PRO	-	expression tag	UNP P35705
G	-8	ARG	-	expression tag	UNP P35705

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Chain	Residue	Modelled	Actual	Comment	Reference
G	-7	GLY	-	expression tag	UNP P35705
G	-6	SER	-	expression tag	UNP P35705
G	-5	HIS	-	expression tag	UNP P35705
G	-4	MET	-	expression tag	UNP P35705
G	-3	LEU	-	expression tag	UNP P35705
G	-2	GLU	-	expression tag	UNP P35705
G	-1	ASP	-	expression tag	UNP P35705
G	0	PRO	-	expression tag	UNP P35705
G	168	SER	CYS	engineered mutation	UNP P35705
H	-24	MET	-	expression tag	UNP P35705
H	-23	GLY	-	expression tag	UNP P35705
H	-22	SER	-	expression tag	UNP P35705
H	-21	SER	-	expression tag	UNP P35705
H	-20	HIS	-	expression tag	UNP P35705
H	-19	HIS	-	expression tag	UNP P35705
H	-18	HIS	-	expression tag	UNP P35705
H	-17	HIS	-	expression tag	UNP P35705
H	-16	HIS	-	expression tag	UNP P35705
H	-15	HIS	-	expression tag	UNP P35705
H	-14	SER	-	expression tag	UNP P35705
H	-13	SER	-	expression tag	UNP P35705
H	-12	GLY	-	expression tag	UNP P35705
H	-11	LEU	-	expression tag	UNP P35705
H	-10	VAL	-	expression tag	UNP P35705
H	-9	PRO	-	expression tag	UNP P35705
H	-8	ARG	-	expression tag	UNP P35705
H	-7	GLY	-	expression tag	UNP P35705
H	-6	SER	-	expression tag	UNP P35705
H	-5	HIS	-	expression tag	UNP P35705
H	-4	MET	-	expression tag	UNP P35705
H	-3	LEU	-	expression tag	UNP P35705
H	-2	GLU	-	expression tag	UNP P35705
H	-1	ASP	-	expression tag	UNP P35705
H	0	PRO	-	expression tag	UNP P35705
H	168	SER	CYS	engineered mutation	UNP P35705
I	-24	MET	-	expression tag	UNP P35705
I	-23	GLY	-	expression tag	UNP P35705
I	-22	SER	-	expression tag	UNP P35705
I	-21	SER	-	expression tag	UNP P35705
I	-20	HIS	-	expression tag	UNP P35705
I	-19	HIS	-	expression tag	UNP P35705
I	-18	HIS	-	expression tag	UNP P35705

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Chain	Residue	Modelled	Actual	Comment	Reference
I	-17	HIS	-	expression tag	UNP P35705
I	-16	HIS	-	expression tag	UNP P35705
I	-15	HIS	-	expression tag	UNP P35705
I	-14	SER	-	expression tag	UNP P35705
I	-13	SER	-	expression tag	UNP P35705
I	-12	GLY	-	expression tag	UNP P35705
I	-11	LEU	-	expression tag	UNP P35705
I	-10	VAL	-	expression tag	UNP P35705
I	-9	PRO	-	expression tag	UNP P35705
I	-8	ARG	-	expression tag	UNP P35705
I	-7	GLY	-	expression tag	UNP P35705
I	-6	SER	-	expression tag	UNP P35705
I	-5	HIS	-	expression tag	UNP P35705
I	-4	MET	-	expression tag	UNP P35705
I	-3	LEU	-	expression tag	UNP P35705
I	-2	GLU	-	expression tag	UNP P35705
I	-1	ASP	-	expression tag	UNP P35705
I	0	PRO	-	expression tag	UNP P35705
I	168	SER	CYS	engineered mutation	UNP P35705
J	-24	MET	-	expression tag	UNP P35705
J	-23	GLY	-	expression tag	UNP P35705
J	-22	SER	-	expression tag	UNP P35705
J	-21	SER	-	expression tag	UNP P35705
J	-20	HIS	-	expression tag	UNP P35705
J	-19	HIS	-	expression tag	UNP P35705
J	-18	HIS	-	expression tag	UNP P35705
J	-17	HIS	-	expression tag	UNP P35705
J	-16	HIS	-	expression tag	UNP P35705
J	-15	HIS	-	expression tag	UNP P35705
J	-14	SER	-	expression tag	UNP P35705
J	-13	SER	-	expression tag	UNP P35705
J	-12	GLY	-	expression tag	UNP P35705
J	-11	LEU	-	expression tag	UNP P35705
J	-10	VAL	-	expression tag	UNP P35705
J	-9	PRO	-	expression tag	UNP P35705
J	-8	ARG	-	expression tag	UNP P35705
J	-7	GLY	-	expression tag	UNP P35705
J	-6	SER	-	expression tag	UNP P35705
J	-5	HIS	-	expression tag	UNP P35705
J	-4	MET	-	expression tag	UNP P35705
J	-3	LEU	-	expression tag	UNP P35705
J	-2	GLU	-	expression tag	UNP P35705

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Chain	Residue	Modelled	Actual	Comment	Reference
J	-1	ASP	-	expression tag	UNP P35705
J	0	PRO	-	expression tag	UNP P35705
J	168	SER	CYS	engineered mutation	UNP P35705
K	-24	MET	-	expression tag	UNP P35705
K	-23	GLY	-	expression tag	UNP P35705
K	-22	SER	-	expression tag	UNP P35705
K	-21	SER	-	expression tag	UNP P35705
K	-20	HIS	-	expression tag	UNP P35705
K	-19	HIS	-	expression tag	UNP P35705
K	-18	HIS	-	expression tag	UNP P35705
K	-17	HIS	-	expression tag	UNP P35705
K	-16	HIS	-	expression tag	UNP P35705
K	-15	HIS	-	expression tag	UNP P35705
K	-14	SER	-	expression tag	UNP P35705
K	-13	SER	-	expression tag	UNP P35705
K	-12	GLY	-	expression tag	UNP P35705
K	-11	LEU	-	expression tag	UNP P35705
K	-10	VAL	-	expression tag	UNP P35705
K	-9	PRO	-	expression tag	UNP P35705
K	-8	ARG	-	expression tag	UNP P35705
K	-7	GLY	-	expression tag	UNP P35705
K	-6	SER	-	expression tag	UNP P35705
K	-5	HIS	-	expression tag	UNP P35705
K	-4	MET	-	expression tag	UNP P35705
K	-3	LEU	-	expression tag	UNP P35705
K	-2	GLU	-	expression tag	UNP P35705
K	-1	ASP	-	expression tag	UNP P35705
K	0	PRO	-	expression tag	UNP P35705
K	168	SER	CYS	engineered mutation	UNP P35705
L	-24	MET	-	expression tag	UNP P35705
L	-23	GLY	-	expression tag	UNP P35705
L	-22	SER	-	expression tag	UNP P35705
L	-21	SER	-	expression tag	UNP P35705
L	-20	HIS	-	expression tag	UNP P35705
L	-19	HIS	-	expression tag	UNP P35705
L	-18	HIS	-	expression tag	UNP P35705
L	-17	HIS	-	expression tag	UNP P35705
L	-16	HIS	-	expression tag	UNP P35705
L	-15	HIS	-	expression tag	UNP P35705
L	-14	SER	-	expression tag	UNP P35705
L	-13	SER	-	expression tag	UNP P35705
L	-12	GLY	-	expression tag	UNP P35705

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Chain	Residue	Modelled	Actual	Comment	Reference
L	-11	LEU	-	expression tag	UNP P35705
L	-10	VAL	-	expression tag	UNP P35705
L	-9	PRO	-	expression tag	UNP P35705
L	-8	ARG	-	expression tag	UNP P35705
L	-7	GLY	-	expression tag	UNP P35705
L	-6	SER	-	expression tag	UNP P35705
L	-5	HIS	-	expression tag	UNP P35705
L	-4	MET	-	expression tag	UNP P35705
L	-3	LEU	-	expression tag	UNP P35705
L	-2	GLU	-	expression tag	UNP P35705
L	-1	ASP	-	expression tag	UNP P35705
L	0	PRO	-	expression tag	UNP P35705
L	168	SER	CYS	engineered mutation	UNP P35705

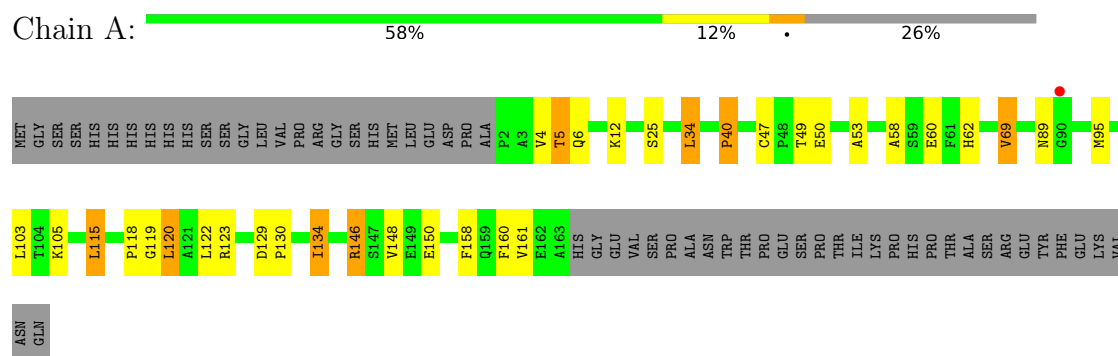
- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	2	Total O 2 2	0	0
2	B	1	Total O 1 1	0	0
2	E	1	Total O 1 1	0	0
2	H	1	Total O 1 1	0	0
2	I	1	Total O 1 1	0	0
2	J	1	Total O 1 1	0	0
2	K	1	Total O 1 1	0	0
2	L	1	Total O 1 1	0	0

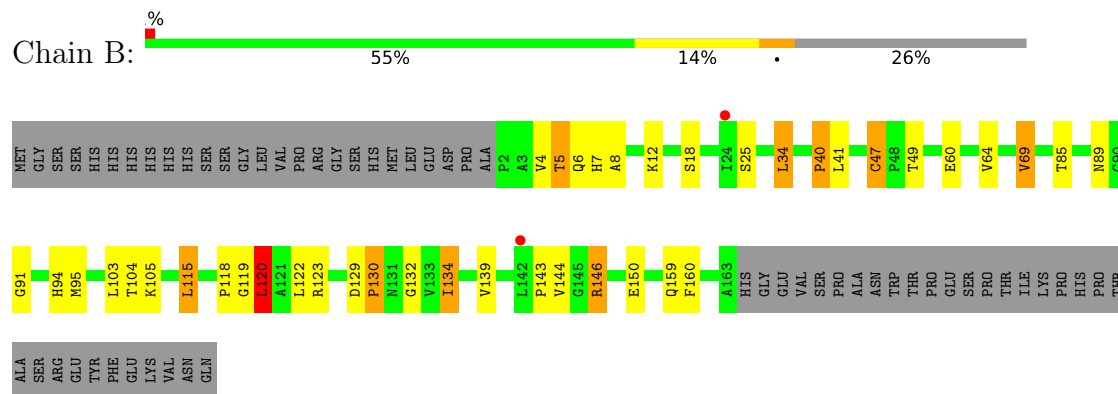
3 Residue-property plots [i](#)

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.

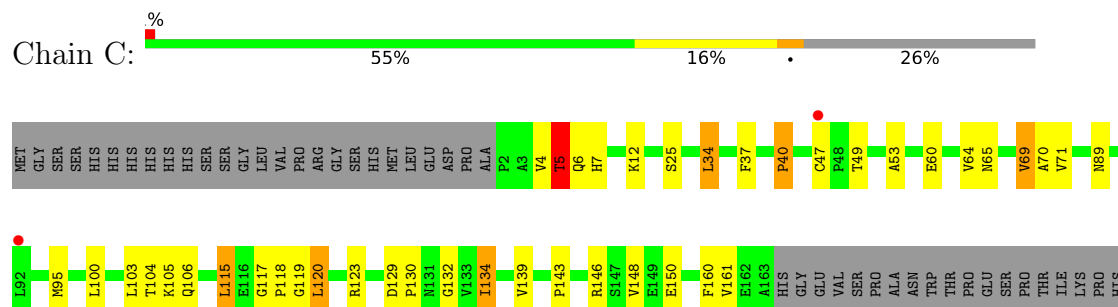
- Molecule 1: Thioredoxin-dependent peroxide reductase



- Molecule 1: Thioredoxin-dependent peroxide reductase

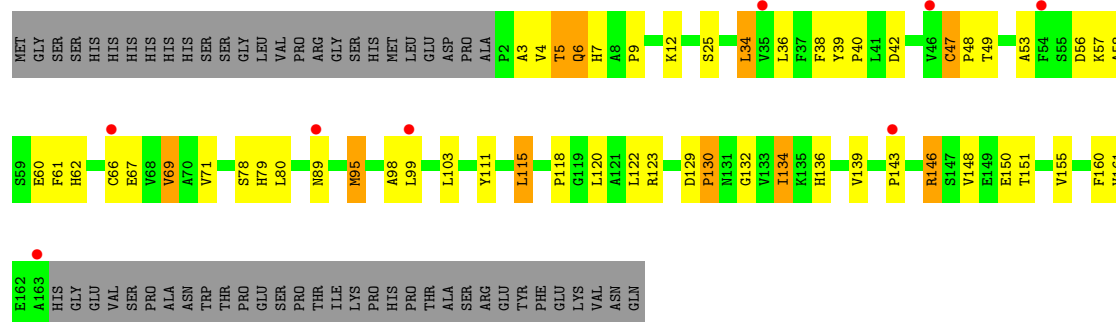


- Molecule 1: Thioredoxin-dependent peroxide reductase

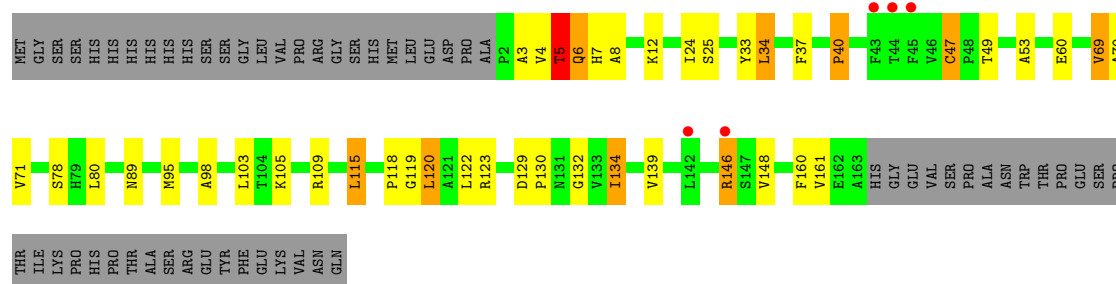


PRO
THR
ALA
SER
ARG
GLU
TYR
PHE
GLU
LYS
VAL
ASN
GLN

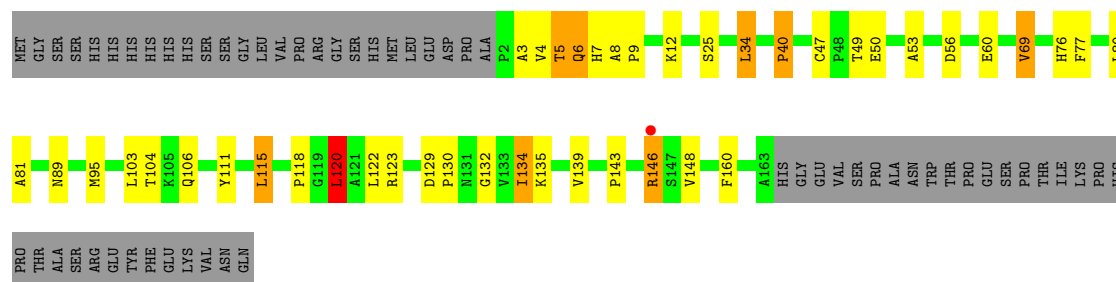
- Molecule 1: Thioredoxin-dependent peroxide reductase



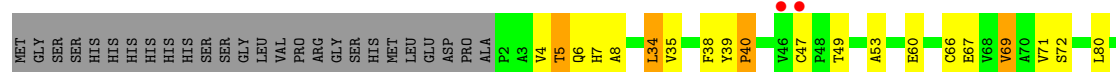
- Molecule 1: Thioredoxin-dependent peroxide reductase

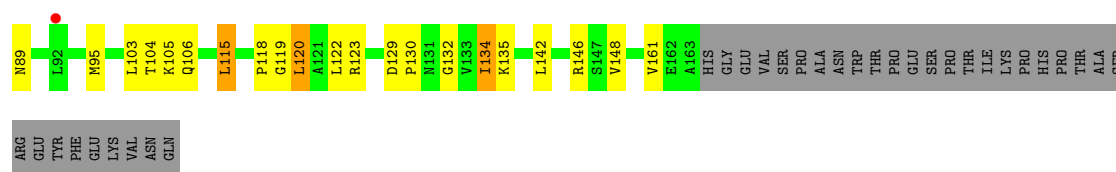


- Molecule 1: Thioredoxin-dependent peroxide reductase

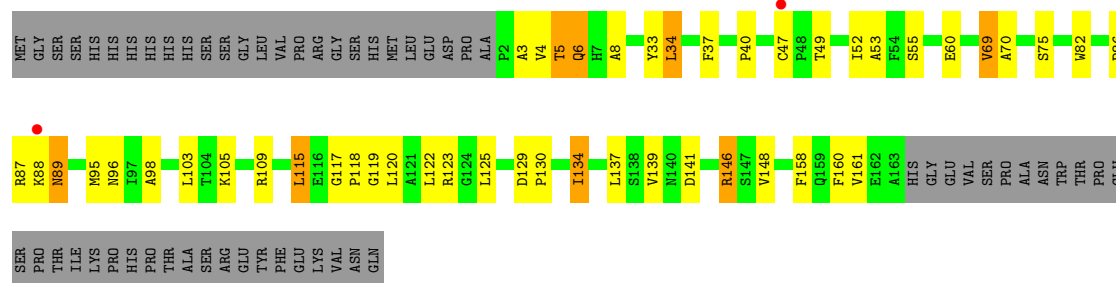


- Molecule 1: Thioredoxin-dependent peroxide reductase

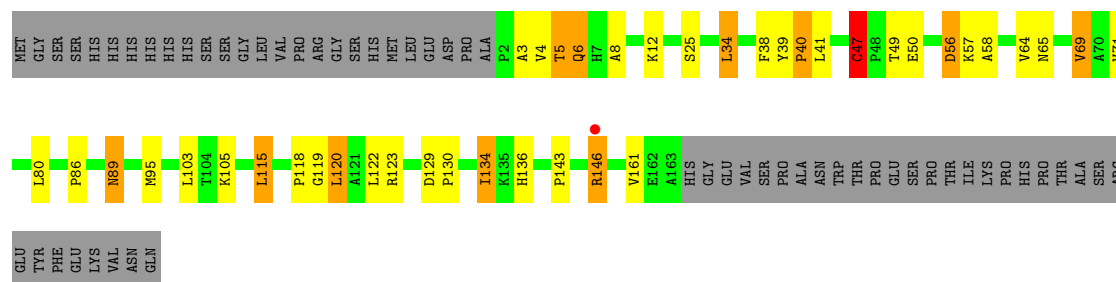




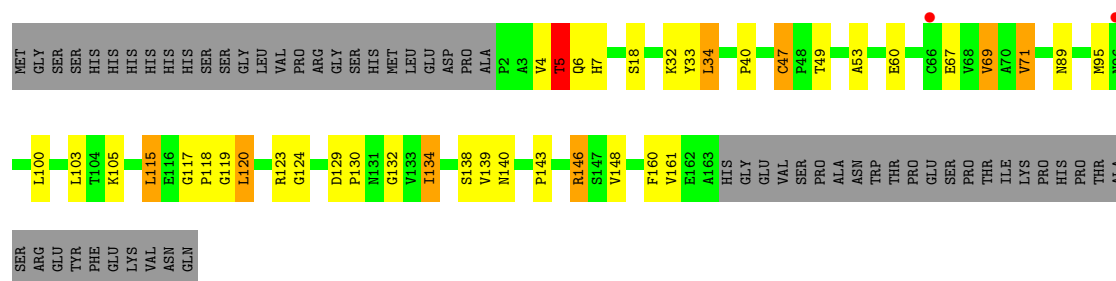
- Molecule 1: Thioredoxin-dependent peroxide reductase



- Molecule 1: Thioredoxin-dependent peroxide reductase

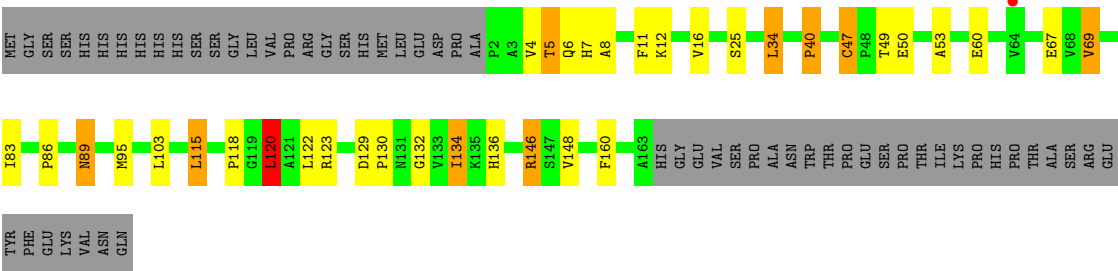


- Molecule 1: Thioredoxin-dependent peroxide reductase

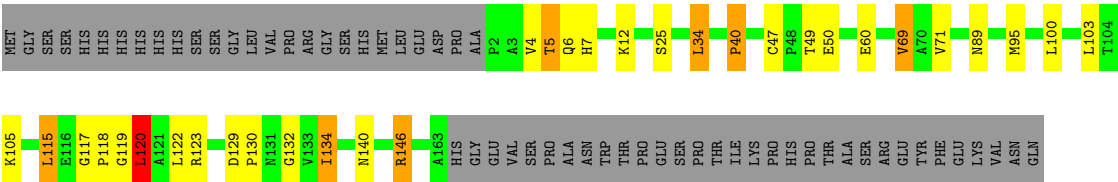


- Molecule 1: Thioredoxin-dependent peroxide reductase





• Molecule 1: Thioredoxin-dependent peroxide reductase



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	301.00Å 80.68Å 124.30Å 90.00° 112.76° 90.00°	Depositor
Resolution (Å)	35.00 – 3.30 138.78 – 3.30	Depositor EDS
% Data completeness (in resolution range)	100.0 (35.00-3.30) 99.9 (138.78-3.30)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.37 (at 3.33Å)	Xtriage
Refinement program	REFMAC 5.2.0005	Depositor
R, R_{free}	0.226 , 0.265 0.220 , 0.256	Depositor DCC
R_{free} test set	2109 reflections (5.04%)	wwPDB-VP
Wilson B-factor (Å ²)	80.0	Xtriage
Anisotropy	0.138	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 58.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.028 for -h-2*1,-k,l	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	15201	wwPDB-VP
Average B, all atoms (Å ²)	72.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.06% 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	0.62	0/1298	0.70	1/1762 (0.1%)
1	B	0.71	0/1298	0.76	1/1762 (0.1%)
1	C	0.72	0/1298	0.76	1/1762 (0.1%)
1	D	0.89	0/1298	0.80	1/1762 (0.1%)
1	E	0.81	0/1298	0.79	1/1762 (0.1%)
1	F	0.81	0/1298	0.77	1/1762 (0.1%)
1	G	0.79	0/1298	0.78	1/1762 (0.1%)
1	H	0.74	0/1298	0.76	0/1762
1	I	0.72	1/1298 (0.1%)	0.78	1/1762 (0.1%)
1	J	0.70	1/1298 (0.1%)	0.79	1/1762 (0.1%)
1	K	0.70	1/1298 (0.1%)	0.75	1/1762 (0.1%)
1	L	0.76	0/1298	0.78	1/1762 (0.1%)
All	All	0.75	3/15576 (0.0%)	0.77	11/21144 (0.1%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	K	47	CYS	CB-SG	-5.93	1.72	1.81
1	I	47	CYS	CB-SG	-5.38	1.73	1.81
1	J	47	CYS	CB-SG	-5.33	1.73	1.81

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	120	LEU	CA-CB-CG	6.39	130.00	115.30
1	K	120	LEU	CA-CB-CG	6.32	129.84	115.30
1	B	120	LEU	CA-CB-CG	6.16	129.47	115.30
1	F	120	LEU	CA-CB-CG	6.13	129.41	115.30
1	G	120	LEU	CA-CB-CG	5.91	128.89	115.30
1	J	120	LEU	CA-CB-CG	5.89	128.84	115.30
1	A	120	LEU	CA-CB-CG	5.62	128.22	115.30
1	E	120	LEU	CA-CB-CG	5.56	128.09	115.30
1	D	95	MET	CG-SD-CE	5.54	109.06	100.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	L	120	LEU	CA-CB-CG	5.39	127.70	115.30
1	I	120	LEU	CA-CB-CG	5.24	127.34	115.30

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	1266	0	1261	21	0
1	B	1266	0	1261	23	0
1	C	1266	0	1261	25	0
1	D	1266	0	1261	33	0
1	E	1266	0	1261	28	0
1	F	1266	0	1261	28	0
1	G	1266	0	1261	22	0
1	H	1266	0	1261	24	0
1	I	1266	0	1261	32	0
1	J	1266	0	1261	21	0
1	K	1266	0	1261	21	0
1	L	1266	0	1261	19	0
2	A	2	0	0	0	0
2	B	1	0	0	0	0
2	E	1	0	0	0	0
2	H	1	0	0	0	0
2	I	1	0	0	0	0
2	J	1	0	0	0	0
2	K	1	0	0	0	0
2	L	1	0	0	0	0
All	All	15201	0	15132	271	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 (271) 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:L:129:ASP:HB2	1:L:130:PRO:HD2	1.57	0.86
1:C:150:GLU:HG2	1:D:150:GLU:HG2	1.57	0.86
1:C:129:ASP:HB2	1:C:130:PRO:HD2	1.65	0.77
1:I:129:ASP:HB2	1:I:130:PRO:HD2	1.72	0.71
1:B:4:VAL:HG22	1:B:134:ILE:HD11	1.70	0.71
1:K:12:LYS:HG3	1:K:25:SER:HB3	1.71	0.71
1:B:34:LEU:HD21	1:B:69:VAL:HG13	1.72	0.71
1:C:139:VAL:HB	1:D:4:VAL:HG21	1.73	0.71
1:G:34:LEU:HD21	1:G:69:VAL:HG13	1.73	0.70
1:E:129:ASP:HB2	1:E:130:PRO:HD2	1.73	0.70
1:L:105:LYS:HE3	1:L:119:GLY:O	1.94	0.68
1:I:34:LEU:HD21	1:I:69:VAL:HG13	1.75	0.67
1:E:34:LEU:HD21	1:E:69:VAL:HG13	1.77	0.67
1:A:129:ASP:HB2	1:A:130:PRO:HD2	1.77	0.66
1:C:161:VAL:HG21	1:D:143:PRO:HB2	1.76	0.66
1:K:129:ASP:HB2	1:K:130:PRO:HD2	1.75	0.66
1:H:34:LEU:HD21	1:H:69:VAL:HG13	1.77	0.66
1:G:129:ASP:HB2	1:G:130:PRO:HD2	1.78	0.66
1:H:4:VAL:HG22	1:H:134:ILE:HD11	1.76	0.65
1:I:4:VAL:HG21	1:J:139:VAL:HB	1.79	0.65
1:J:105:LYS:HE3	1:J:119:GLY:O	1.97	0.64
1:H:123:ARG:HB3	1:H:146:ARG:NH2	2.12	0.64
1:C:34:LEU:HD21	1:C:69:VAL:HG13	1.80	0.64
1:A:34:LEU:HD21	1:A:69:VAL:HG13	1.80	0.63
1:J:129:ASP:HB2	1:J:130:PRO:HD2	1.80	0.63
1:A:4:VAL:HG21	1:B:139:VAL:HB	1.81	0.63
1:A:150:GLU:HG2	1:B:150:GLU:HG2	1.80	0.63
1:G:105:LYS:HE3	1:G:119:GLY:O	1.99	0.63
1:B:12:LYS:HG3	1:B:25:SER:HB3	1.81	0.63
1:J:123:ARG:HB3	1:J:146:ARG:NH2	2.14	0.63
1:L:34:LEU:HD21	1:L:69:VAL:HG13	1.80	0.62
1:H:105:LYS:HE3	1:H:119:GLY:O	1.99	0.62
1:L:12:LYS:HG3	1:L:25:SER:HB3	1.80	0.62
1:K:34:LEU:HD21	1:K:69:VAL:HG13	1.80	0.62
1:C:105:LYS:HE3	1:C:119:GLY:O	1.99	0.62
1:E:37:PHE:CE1	1:E:70:ALA:HB2	2.35	0.61
1:B:129:ASP:HB2	1:B:130:PRO:HD2	1.82	0.61
1:J:4:VAL:HG22	1:J:134:ILE:HD11	1.82	0.61
1:E:123:ARG:HB3	1:E:146:ARG:NH2	2.16	0.61
1:J:34:LEU:HD21	1:J:69:VAL:HG13	1.83	0.61
1:I:123:ARG:HB3	1:I:146:ARG:NH2	2.15	0.60
1:F:129:ASP:HB2	1:F:130:PRO:HD2	1.83	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:12:LYS:HG3	1:D:25:SER:HB3	1.82	0.60
1:C:40:PRO:HD2	1:C:47:CYS:SG	2.42	0.60
1:C:115:LEU:HB3	1:C:118:PRO:HD2	1.85	0.59
1:I:4:VAL:HG22	1:I:134:ILE:HD11	1.84	0.59
1:G:4:VAL:HG22	1:G:134:ILE:HD11	1.83	0.59
1:C:143:PRO:HB2	1:D:161:VAL:HG21	1.84	0.59
1:L:50:GLU:OE1	1:L:123:ARG:NH1	2.36	0.58
1:F:3:ALA:O	1:F:6:GLN:HG3	2.02	0.58
1:A:58:ALA:O	1:A:62:HIS:ND1	2.35	0.58
1:E:53:ALA:HB1	1:E:148:VAL:HG21	1.84	0.58
1:K:50:GLU:OE1	1:K:123:ARG:NH1	2.36	0.58
1:C:4:VAL:HA	1:C:134:ILE:HG13	1.84	0.58
1:G:4:VAL:HG21	1:H:139:VAL:HB	1.86	0.58
1:D:129:ASP:HB2	1:D:130:PRO:HD2	1.85	0.57
1:F:40:PRO:HD2	1:F:47:CYS:SG	2.44	0.57
1:F:104:THR:HG23	1:F:106:GLN:H	1.69	0.57
1:F:34:LEU:HD21	1:F:69:VAL:HG13	1.87	0.57
1:A:105:LYS:HE3	1:A:119:GLY:O	2.05	0.57
1:E:4:VAL:HG21	1:F:139:VAL:CG1	2.34	0.56
1:A:4:VAL:HG22	1:A:134:ILE:HD11	1.88	0.56
1:D:151:THR:O	1:D:155:VAL:HG23	2.05	0.56
1:H:53:ALA:HB1	1:H:148:VAL:HG21	1.88	0.56
1:L:4:VAL:HG22	1:L:134:ILE:HD11	1.88	0.56
1:I:115:LEU:HB3	1:I:118:PRO:HD2	1.87	0.56
1:F:50:GLU:OE1	1:F:123:ARG:NH1	2.39	0.55
1:K:8:ALA:HB2	1:K:134:ILE:HG22	1.88	0.55
1:H:69:VAL:HB	1:H:98:ALA:HB3	1.87	0.55
1:C:4:VAL:HG21	1:D:139:VAL:HB	1.89	0.55
1:K:115:LEU:HB3	1:K:118:PRO:HD2	1.87	0.55
1:E:161:VAL:HG21	1:F:143:PRO:HB2	1.89	0.55
1:E:69:VAL:HB	1:E:98:ALA:HB3	1.89	0.54
1:F:12:LYS:HG3	1:F:25:SER:HB3	1.89	0.54
1:D:34:LEU:HD21	1:D:69:VAL:HG13	1.89	0.54
1:D:53:ALA:HB1	1:D:148:VAL:HG21	1.89	0.54
1:I:12:LYS:HG3	1:I:25:SER:HB3	1.89	0.54
1:L:40:PRO:HD2	1:L:47:CYS:SG	2.48	0.54
1:I:143:PRO:CG	1:J:161:VAL:HG21	2.38	0.54
1:L:4:VAL:HA	1:L:134:ILE:HG13	1.89	0.54
1:B:123:ARG:HB3	1:B:146:ARG:NH2	2.23	0.54
1:E:105:LYS:HE3	1:E:119:GLY:O	2.07	0.53
1:H:115:LEU:HD13	1:H:122:LEU:HG	1.91	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:L:115:LEU:HB3	1:L:118:PRO:HD2	1.90	0.53
1:A:115:LEU:HB3	1:A:118:PRO:HD2	1.90	0.53
1:F:47:CYS:HA	1:F:123:ARG:HH22	1.74	0.53
1:E:53:ALA:CB	1:E:148:VAL:HG21	2.39	0.53
1:F:104:THR:HG23	1:F:106:GLN:CB	2.39	0.52
1:J:4:VAL:O	1:J:5:THR:OG1	2.25	0.52
1:J:53:ALA:HB1	1:J:148:VAL:HG21	1.91	0.52
1:B:4:VAL:HA	1:B:134:ILE:HG13	1.90	0.52
1:D:123:ARG:HB3	1:D:146:ARG:NH2	2.24	0.52
1:E:47:CYS:HA	1:E:123:ARG:HH22	1.74	0.52
1:E:115:LEU:HB3	1:E:118:PRO:HD2	1.91	0.52
1:L:123:ARG:HB3	1:L:146:ARG:NH2	2.25	0.52
1:F:104:THR:HG23	1:F:106:GLN:HB3	1.91	0.52
1:G:115:LEU:HB3	1:G:118:PRO:HD2	1.91	0.52
1:A:161:VAL:HG21	1:B:143:PRO:HB2	1.90	0.52
1:H:8:ALA:HB2	1:H:134:ILE:HG22	1.91	0.52
1:I:50:GLU:OE1	1:I:123:ARG:NH1	2.43	0.52
1:G:67:GLU:OE1	1:G:67:GLU:HA	2.09	0.52
1:E:139:VAL:HB	1:F:4:VAL:HG21	1.92	0.51
1:A:12:LYS:HG3	1:A:25:SER:HB3	1.92	0.51
1:A:40:PRO:HD2	1:A:47:CYS:SG	2.50	0.51
1:D:7:HIS:HA	1:D:132:GLY:O	2.10	0.51
1:F:115:LEU:HB3	1:F:118:PRO:HD2	1.92	0.51
1:D:98:ALA:O	1:D:99:LEU:HD12	2.11	0.51
1:A:123:ARG:HB3	1:A:146:ARG:NH2	2.25	0.51
1:C:4:VAL:HG22	1:C:134:ILE:HD11	1.91	0.51
1:F:129:ASP:OD2	1:F:135:LYS:HD3	2.10	0.51
1:C:139:VAL:O	1:D:136:HIS:HA	2.10	0.51
1:K:122:LEU:HD13	1:L:4:VAL:HB	1.93	0.50
1:J:4:VAL:HA	1:J:134:ILE:HG13	1.92	0.50
1:H:115:LEU:HB3	1:H:118:PRO:HD2	1.93	0.50
1:H:129:ASP:HB2	1:H:130:PRO:HD2	1.94	0.50
1:D:67:GLU:HA	1:D:67:GLU:OE1	2.11	0.50
1:H:52:ILE:HG23	1:H:87:ARG:NH1	2.26	0.50
1:J:115:LEU:HB3	1:J:118:PRO:HD2	1.94	0.50
1:H:53:ALA:CB	1:H:148:VAL:HG21	2.41	0.50
1:D:4:VAL:HG22	1:D:134:ILE:HD11	1.94	0.49
1:D:4:VAL:HA	1:D:134:ILE:HG13	1.93	0.49
1:I:105:LYS:HE3	1:I:119:GLY:O	2.12	0.49
1:E:7:HIS:HA	1:E:132:GLY:O	2.13	0.49
1:K:123:ARG:HB3	1:K:146:ARG:NH2	2.28	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:71:VAL:HB	1:C:100:LEU:HB3	1.93	0.49
1:B:115:LEU:HB3	1:B:118:PRO:HD2	1.93	0.49
1:G:4:VAL:HA	1:G:134:ILE:HG13	1.95	0.49
1:B:40:PRO:HD2	1:B:47:CYS:SG	2.52	0.49
1:H:88:LYS:HD3	1:H:89:ASN:OD1	2.13	0.48
1:A:4:VAL:HA	1:A:134:ILE:HG13	1.95	0.48
1:C:7:HIS:HA	1:C:132:GLY:O	2.13	0.48
1:H:3:ALA:O	1:H:6:GLN:HG3	2.14	0.48
1:J:34:LEU:HD21	1:J:69:VAL:CG1	2.44	0.48
1:E:40:PRO:HD2	1:E:47:CYS:SG	2.54	0.48
1:I:134:ILE:HG13	1:I:134:ILE:O	2.14	0.48
1:C:161:VAL:HG21	1:D:143:PRO:CB	2.42	0.48
1:F:118:PRO:HG2	1:F:120:LEU:HD22	1.96	0.48
1:D:3:ALA:HB3	1:D:6:GLN:HE21	1.78	0.47
1:D:42:ASP:OD2	1:D:79:HIS:ND1	2.47	0.47
1:E:4:VAL:O	1:E:5:THR:OG1	2.29	0.47
1:E:8:ALA:HB2	1:E:134:ILE:HG22	1.96	0.47
1:D:39:TYR:HB2	1:D:47:CYS:SG	2.54	0.47
1:E:12:LYS:HG3	1:E:25:SER:HB3	1.97	0.47
1:D:56:ASP:C	1:D:58:ALA:H	2.18	0.47
1:I:64:VAL:O	1:I:65:ASN:HB2	2.14	0.47
1:I:161:VAL:HG21	1:J:143:PRO:HB2	1.96	0.47
1:K:11:PHE:O	1:K:25:SER:HA	2.15	0.47
1:A:62:HIS:HD1	1:A:62:HIS:H	1.63	0.47
1:B:7:HIS:HA	1:B:132:GLY:O	2.15	0.47
1:K:40:PRO:HD2	1:K:47:CYS:SG	2.55	0.47
1:D:115:LEU:HB3	1:D:118:PRO:HD2	1.96	0.46
1:I:40:PRO:HD2	1:I:47:CYS:SG	2.54	0.46
1:F:7:HIS:HA	1:F:132:GLY:O	2.16	0.46
1:G:34:LEU:HD21	1:G:69:VAL:CG1	2.43	0.46
1:K:86:PRO:HD2	1:K:89:ASN:ND2	2.30	0.46
1:K:118:PRO:HB2	1:K:120:LEU:HD22	1.97	0.46
1:D:9:PRO:HD2	1:D:111:TYR:CZ	2.51	0.46
1:C:4:VAL:O	1:C:5:THR:OG1	2.31	0.46
1:D:53:ALA:CB	1:D:148:VAL:HG21	2.46	0.46
1:L:115:LEU:HD13	1:L:122:LEU:HG	1.97	0.46
1:E:123:ARG:HD3	1:E:146:ARG:HH21	1.81	0.46
1:J:32:LYS:HE2	1:J:67:GLU:HG2	1.97	0.46
1:D:80:LEU:HD12	1:D:80:LEU:HA	1.65	0.45
1:H:3:ALA:HB3	1:H:6:GLN:HE21	1.80	0.45
1:K:4:VAL:HG22	1:K:134:ILE:HD11	1.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:115:LEU:HD13	1:G:122:LEU:HG	1.97	0.45
1:C:104:THR:OG1	1:C:106:GLN:HB2	2.16	0.45
1:H:47:CYS:HA	1:H:123:ARG:HH22	1.81	0.45
1:E:4:VAL:HG21	1:F:139:VAL:HB	1.99	0.45
1:E:37:PHE:CZ	1:E:70:ALA:HB2	2.52	0.45
1:G:40:PRO:HD2	1:G:47:CYS:SG	2.57	0.45
1:J:33:TYR:CD1	1:J:130:PRO:HD3	2.52	0.45
1:G:39:TYR:O	1:G:72:SER:HB2	2.16	0.45
1:I:3:ALA:HB3	1:I:6:GLN:HE21	1.82	0.44
1:I:115:LEU:HD13	1:I:122:LEU:HG	1.97	0.44
1:A:53:ALA:HB1	1:A:148:VAL:HG21	1.98	0.44
1:B:34:LEU:HD21	1:B:69:VAL:CG1	2.45	0.44
1:E:115:LEU:HD13	1:E:122:LEU:HG	1.98	0.44
1:F:4:VAL:HA	1:F:134:ILE:HG13	1.99	0.44
1:F:9:PRO:HD2	1:F:111:TYR:CE2	2.52	0.44
1:H:37:PHE:CE1	1:H:70:ALA:HB2	2.52	0.44
1:H:86:PRO:HD2	1:H:89:ASN:ND2	2.32	0.44
1:I:8:ALA:HB2	1:I:134:ILE:HG22	1.99	0.44
1:B:41:LEU:HD11	1:B:120:LEU:HD12	1.99	0.44
1:K:136:HIS:ND1	1:L:140:ASN:OD1	2.44	0.44
1:F:53:ALA:HB1	1:F:148:VAL:HG21	2.00	0.44
1:H:75:SER:HB3	1:I:41:LEU:HD22	1.99	0.44
1:F:53:ALA:CB	1:F:148:VAL:HG21	2.48	0.44
1:I:129:ASP:CB	1:I:130:PRO:HD2	2.38	0.44
1:E:33:TYR:CD1	1:E:130:PRO:HD3	2.52	0.43
1:G:7:HIS:HA	1:G:132:GLY:O	2.17	0.43
1:I:56:ASP:C	1:I:58:ALA:H	2.21	0.43
1:I:38:PHE:HA	1:I:71:VAL:O	2.18	0.43
1:B:85:THR:HG22	1:B:91:GLY:HA3	1.99	0.43
1:E:129:ASP:CB	1:E:130:PRO:HD2	2.42	0.43
1:H:33:TYR:CD1	1:H:130:PRO:HD3	2.52	0.43
1:L:118:PRO:HG2	1:L:120:LEU:HD22	2.00	0.43
1:H:158:PHE:HA	1:H:161:VAL:HG12	1.98	0.43
1:D:58:ALA:O	1:D:62:HIS:ND1	2.48	0.43
1:C:37:PHE:CE1	1:C:70:ALA:HB2	2.54	0.43
1:E:24:ILE:O	1:E:24:ILE:HG13	2.18	0.43
1:G:35:VAL:HG23	1:G:66:CYS:SG	2.59	0.43
1:I:136:HIS:ND1	1:J:140:ASN:OD1	2.45	0.43
1:B:105:LYS:HE3	1:B:119:GLY:O	2.18	0.43
1:G:53:ALA:HB1	1:G:148:VAL:HG21	2.01	0.43
1:G:80:LEU:HD12	1:G:80:LEU:HA	1.81	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:I:56:ASP:O	1:I:58:ALA:N	2.52	0.43
1:K:53:ALA:CB	1:K:148:VAL:HG21	2.49	0.43
1:A:4:VAL:HB	1:B:122:LEU:HD13	2.01	0.42
1:D:38:PHE:HA	1:D:71:VAL:O	2.19	0.42
1:I:161:VAL:O	1:I:161:VAL:HG22	2.19	0.42
1:K:129:ASP:OD1	1:K:129:ASP:C	2.57	0.42
1:B:8:ALA:HB2	1:B:134:ILE:HG22	2.00	0.42
1:I:123:ARG:HD3	1:I:146:ARG:HH21	1.83	0.42
1:L:7:HIS:HA	1:L:132:GLY:O	2.20	0.42
1:B:123:ARG:HD3	1:B:146:ARG:HH21	1.85	0.42
1:D:61:PHE:HD2	1:D:66:CYS:HG	1.67	0.42
1:J:7:HIS:HA	1:J:132:GLY:O	2.19	0.42
1:K:4:VAL:HA	1:K:134:ILE:HG13	2.01	0.42
1:B:64:VAL:O	1:B:159:GLN:NE2	2.41	0.42
1:C:47:CYS:HA	1:C:123:ARG:HH22	1.84	0.42
1:D:58:ALA:C	1:D:62:HIS:HD1	2.22	0.42
1:F:4:VAL:HG22	1:F:134:ILE:HD11	2.00	0.42
1:G:161:VAL:O	1:G:161:VAL:HG22	2.19	0.42
1:I:34:LEU:HD21	1:I:69:VAL:CG1	2.47	0.42
1:J:47:CYS:HA	1:J:123:ARG:HH22	1.84	0.42
1:K:7:HIS:HA	1:K:132:GLY:O	2.19	0.42
1:K:67:GLU:OE1	1:K:67:GLU:HA	2.19	0.42
1:A:115:LEU:HD13	1:A:122:LEU:HG	2.02	0.42
1:L:129:ASP:CB	1:L:130:PRO:HD2	2.30	0.42
1:A:47:CYS:HA	1:A:123:ARG:HH22	1.85	0.42
1:A:158:PHE:CD1	1:B:144:VAL:HG21	2.55	0.42
1:E:3:ALA:HB3	1:E:6:GLN:HE21	1.85	0.42
1:I:47:CYS:HA	1:I:123:ARG:HH22	1.84	0.42
1:G:38:PHE:HA	1:G:71:VAL:O	2.20	0.42
1:L:120:LEU:HD23	1:L:120:LEU:O	2.20	0.42
1:I:161:VAL:HG21	1:J:143:PRO:CB	2.50	0.42
1:E:4:VAL:HB	1:F:122:LEU:HD13	2.01	0.41
1:I:86:PRO:HD2	1:I:89:ASN:ND2	2.35	0.41
1:F:104:THR:C	1:F:106:GLN:H	2.23	0.41
1:D:115:LEU:HD13	1:D:122:LEU:HG	2.01	0.41
1:G:104:THR:HG23	1:G:106:GLN:H	1.85	0.41
1:A:50:GLU:OE1	1:A:123:ARG:NH1	2.54	0.41
1:I:39:TYR:HB2	1:I:47:CYS:SG	2.61	0.41
1:K:47:CYS:HA	1:K:123:ARG:HH22	1.85	0.41
1:C:12:LYS:HG3	1:C:25:SER:HB3	2.03	0.41
1:I:80:LEU:HD12	1:I:80:LEU:HA	1.86	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:129:ASP:OD1	1:B:129:ASP:C	2.59	0.41
1:H:125:LEU:O	1:H:137:LEU:HA	2.21	0.41
1:B:104:THR:O	1:B:105:LYS:HB2	2.21	0.41
1:D:47:CYS:N	1:D:48:PRO:CD	2.83	0.41
1:F:50:GLU:OE2	1:F:146:ARG:CZ	2.68	0.41
1:J:71:VAL:HB	1:J:100:LEU:HB3	2.02	0.41
1:C:65:ASN:CG	1:C:65:ASN:O	2.59	0.41
1:K:16:VAL:HG11	1:K:83:ILE:CD1	2.51	0.41
1:L:47:CYS:HA	1:L:123:ARG:HH22	1.86	0.41
1:C:129:ASP:OD1	1:C:129:ASP:C	2.59	0.41
1:F:8:ALA:HB2	1:F:134:ILE:HG22	2.03	0.41
1:C:53:ALA:HB1	1:C:148:VAL:HG21	2.03	0.40
1:G:123:ARG:NE	1:G:142:LEU:O	2.38	0.40
1:C:129:ASP:CB	1:C:130:PRO:HD2	2.35	0.40
1:D:34:LEU:HD22	1:D:36:LEU:N	2.36	0.40
1:G:8:ALA:HB2	1:G:134:ILE:HG22	2.03	0.40
1:H:55:SER:O	1:H:96:ASN:HB2	2.21	0.40
1:L:71:VAL:HB	1:L:100:LEU:HB3	2.02	0.40
1:F:76:HIS:NE2	1:F:77:PHE:CE2	2.89	0.40
1:G:129:ASP:OD2	1:G:135:LYS:HD3	2.21	0.40
1:J:124:GLY:HA2	1:J:138:SER:O	2.22	0.40
1:A:53:ALA:CB	1:A:148:VAL:HG21	2.52	0.40
1:E:80:LEU:HD12	1:E:80:LEU:HA	1.74	0.40
1:I:86:PRO:HD2	1:I:89:ASN:HD22	1.86	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	160/220 (73%)	143 (89%)	15 (9%)	2 (1%)	12 40

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	160/220 (73%)	144 (90%)	13 (8%)	3 (2%)	8	34
1	C	160/220 (73%)	141 (88%)	15 (9%)	4 (2%)	5	27
1	D	160/220 (73%)	141 (88%)	15 (9%)	4 (2%)	5	27
1	E	160/220 (73%)	142 (89%)	16 (10%)	2 (1%)	12	40
1	F	160/220 (73%)	140 (88%)	16 (10%)	4 (2%)	5	27
1	G	160/220 (73%)	143 (89%)	15 (9%)	2 (1%)	12	40
1	H	160/220 (73%)	139 (87%)	17 (11%)	4 (2%)	5	27
1	I	160/220 (73%)	139 (87%)	17 (11%)	4 (2%)	5	27
1	J	160/220 (73%)	140 (88%)	17 (11%)	3 (2%)	8	34
1	K	160/220 (73%)	143 (89%)	15 (9%)	2 (1%)	12	40
1	L	160/220 (73%)	146 (91%)	11 (7%)	3 (2%)	8	34
All	All	1920/2640 (73%)	1701 (89%)	182 (10%)	37 (2%)	8	34

All (37) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	5	THR
1	B	5	THR
1	C	5	THR
1	D	5	THR
1	E	5	THR
1	F	5	THR
1	G	5	THR
1	H	5	THR
1	I	5	THR
1	J	5	THR
1	K	5	THR
1	L	5	THR
1	B	40	PRO
1	I	57	LYS
1	J	40	PRO
1	K	40	PRO
1	A	40	PRO
1	C	40	PRO
1	D	40	PRO
1	D	57	LYS
1	E	40	PRO
1	F	40	PRO

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Mol	Chain	Res	Type
1	G	40	PRO
1	H	40	PRO
1	H	141	ASP
1	I	40	PRO
1	I	56	ASP
1	L	40	PRO
1	F	56	ASP
1	F	81	ALA
1	J	117	GLY
1	C	64	VAL
1	C	117	GLY
1	B	130	PRO
1	D	130	PRO
1	H	117	GLY
1	L	117	GLY

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	139/190 (73%)	125 (90%)	14 (10%)	7	27
1	B	139/190 (73%)	122 (88%)	17 (12%)	5	20
1	C	139/190 (73%)	125 (90%)	14 (10%)	7	27
1	D	139/190 (73%)	123 (88%)	16 (12%)	5	22
1	E	139/190 (73%)	121 (87%)	18 (13%)	4	18
1	F	139/190 (73%)	124 (89%)	15 (11%)	6	24
1	G	139/190 (73%)	126 (91%)	13 (9%)	8	30
1	H	139/190 (73%)	123 (88%)	16 (12%)	5	22
1	I	139/190 (73%)	126 (91%)	13 (9%)	8	30
1	J	139/190 (73%)	123 (88%)	16 (12%)	5	22
1	K	139/190 (73%)	125 (90%)	14 (10%)	7	27
1	L	139/190 (73%)	126 (91%)	13 (9%)	8	30

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
All	All	1668/2280 (73%)	1489 (89%)	179 (11%)	6 25

All (179) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	5	THR
1	A	6	GLN
1	A	34	LEU
1	A	49	THR
1	A	60	GLU
1	A	69	VAL
1	A	89	ASN
1	A	95	MET
1	A	103	LEU
1	A	115	LEU
1	A	120	LEU
1	A	134	ILE
1	A	146	ARG
1	A	160	PHE
1	B	5	THR
1	B	6	GLN
1	B	18	SER
1	B	34	LEU
1	B	47	CYS
1	B	49	THR
1	B	60	GLU
1	B	69	VAL
1	B	89	ASN
1	B	94	HIS
1	B	95	MET
1	B	103	LEU
1	B	115	LEU
1	B	120	LEU
1	B	134	ILE
1	B	146	ARG
1	B	160	PHE
1	C	5	THR
1	C	6	GLN
1	C	34	LEU
1	C	49	THR
1	C	60	GLU
1	C	69	VAL

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Mol	Chain	Res	Type
1	C	89	ASN
1	C	95	MET
1	C	103	LEU
1	C	115	LEU
1	C	120	LEU
1	C	134	ILE
1	C	146	ARG
1	C	160	PHE
1	D	5	THR
1	D	6	GLN
1	D	34	LEU
1	D	47	CYS
1	D	49	THR
1	D	60	GLU
1	D	69	VAL
1	D	78	SER
1	D	89	ASN
1	D	95	MET
1	D	103	LEU
1	D	115	LEU
1	D	120	LEU
1	D	134	ILE
1	D	146	ARG
1	D	160	PHE
1	E	5	THR
1	E	6	GLN
1	E	34	LEU
1	E	47	CYS
1	E	49	THR
1	E	60	GLU
1	E	69	VAL
1	E	71	VAL
1	E	78	SER
1	E	89	ASN
1	E	95	MET
1	E	103	LEU
1	E	109	ARG
1	E	115	LEU
1	E	120	LEU
1	E	134	ILE
1	E	146	ARG
1	E	160	PHE

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Mol	Chain	Res	Type
1	F	5	THR
1	F	6	GLN
1	F	34	LEU
1	F	49	THR
1	F	60	GLU
1	F	69	VAL
1	F	80	LEU
1	F	89	ASN
1	F	95	MET
1	F	103	LEU
1	F	115	LEU
1	F	120	LEU
1	F	134	ILE
1	F	146	ARG
1	F	160	PHE
1	G	5	THR
1	G	6	GLN
1	G	34	LEU
1	G	49	THR
1	G	60	GLU
1	G	69	VAL
1	G	89	ASN
1	G	95	MET
1	G	103	LEU
1	G	115	LEU
1	G	120	LEU
1	G	134	ILE
1	G	146	ARG
1	H	5	THR
1	H	6	GLN
1	H	34	LEU
1	H	49	THR
1	H	60	GLU
1	H	69	VAL
1	H	82	TRP
1	H	89	ASN
1	H	95	MET
1	H	103	LEU
1	H	109	ARG
1	H	115	LEU
1	H	120	LEU
1	H	134	ILE

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Mol	Chain	Res	Type
1	H	146	ARG
1	H	160	PHE
1	I	5	THR
1	I	6	GLN
1	I	34	LEU
1	I	47	CYS
1	I	49	THR
1	I	69	VAL
1	I	89	ASN
1	I	95	MET
1	I	103	LEU
1	I	115	LEU
1	I	120	LEU
1	I	134	ILE
1	I	146	ARG
1	J	5	THR
1	J	6	GLN
1	J	18	SER
1	J	34	LEU
1	J	49	THR
1	J	60	GLU
1	J	69	VAL
1	J	71	VAL
1	J	89	ASN
1	J	95	MET
1	J	103	LEU
1	J	115	LEU
1	J	120	LEU
1	J	134	ILE
1	J	146	ARG
1	J	160	PHE
1	K	5	THR
1	K	6	GLN
1	K	34	LEU
1	K	49	THR
1	K	60	GLU
1	K	69	VAL
1	K	89	ASN
1	K	95	MET
1	K	103	LEU
1	K	115	LEU
1	K	120	LEU

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Mol	Chain	Res	Type
1	K	134	ILE
1	K	146	ARG
1	K	160	PHE
1	L	5	THR
1	L	6	GLN
1	L	34	LEU
1	L	49	THR
1	L	60	GLU
1	L	69	VAL
1	L	89	ASN
1	L	95	MET
1	L	103	LEU
1	L	115	LEU
1	L	120	LEU
1	L	134	ILE
1	L	146	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	6	GLN
1	B	6	GLN
1	D	6	GLN
1	E	6	GLN
1	E	89	ASN
1	G	6	GLN
1	H	6	GLN
1	I	6	GLN
1	I	89	ASN
1	K	89	ASN
1	L	6	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 monosaccharides 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	162/220 (73%)	0.51	1 (0%) 89 90	58, 74, 87, 92	0
1	B	162/220 (73%)	0.40	2 (1%) 79 78	58, 74, 86, 92	0
1	C	162/220 (73%)	0.41	2 (1%) 79 78	57, 74, 86, 92	0
1	D	162/220 (73%)	0.58	8 (4%) 29 27	58, 74, 87, 91	0
1	E	162/220 (73%)	0.43	5 (3%) 49 48	56, 74, 87, 92	0
1	F	162/220 (73%)	0.41	1 (0%) 89 90	58, 74, 87, 92	0
1	G	162/220 (73%)	0.38	3 (1%) 66 65	58, 74, 86, 92	0
1	H	162/220 (73%)	0.36	2 (1%) 79 78	57, 74, 86, 92	0
1	I	162/220 (73%)	0.35	1 (0%) 89 90	58, 74, 87, 90	0
1	J	162/220 (73%)	0.41	2 (1%) 79 78	58, 73, 86, 92	0
1	K	162/220 (73%)	0.40	1 (0%) 89 90	58, 74, 86, 92	0
1	L	162/220 (73%)	0.47	0 100 100	58, 73, 86, 92	0
All	All	1944/2640 (73%)	0.43	28 (1%) 75 75	56, 74, 87, 92	0

All (28) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	44	THR	3.6
1	G	46	VAL	3.3
1	H	88	LYS	3.2
1	F	146	ARG	3.1
1	K	64	VAL	3.0
1	E	45	PHE	2.8
1	D	143	PRO	2.8
1	D	66	CYS	2.7
1	G	92	LEU	2.6
1	D	89	ASN	2.5
1	I	146	ARG	2.4

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Mol	Chain	Res	Type	RSRZ
1	E	43	PHE	2.3
1	D	46	VAL	2.3
1	D	54	PHE	2.3
1	D	35	VAL	2.3
1	C	47	CYS	2.2
1	C	92	LEU	2.2
1	A	90	GLY	2.2
1	E	142	LEU	2.2
1	B	24	ILE	2.1
1	J	66	CYS	2.1
1	E	146	ARG	2.1
1	D	163	ALA	2.1
1	G	47	CYS	2.1
1	B	142	LEU	2.0
1	D	99	LEU	2.0
1	J	96	ASN	2.0
1	H	47	CYS	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.