



Full wwPDB X-ray Structure Validation Report ⓘ

Sep 6, 2023 – 05:08 AM EDT

PDB ID : 4DEJ
Title : Crystal structure of glutathione transferase-like protein IL0419 (Target EFI-501089) from *Idiomarina loihiensis* L2TR
Authors : Patskovsky, Y.; Toro, R.; Bhosle, R.; Zencheck, W.D.; Hillerich, B.; Seidel, R.D.; Washington, E.; Scott Glenn, A.; Chowdhury, S.; Evans, B.; Hammonds, J.; Imker, H.J.; Armstrong, R.N.; Gerlt, J.A.; Almo, S.C.; Enzyme Function Initiative (EFI)
Deposited on : 2012-01-20
Resolution : 2.90 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 : 4.02b-467
Xtrriage (Phenix) : 1.13
EDS : 2.35
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Rfmac : 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) : 2.35

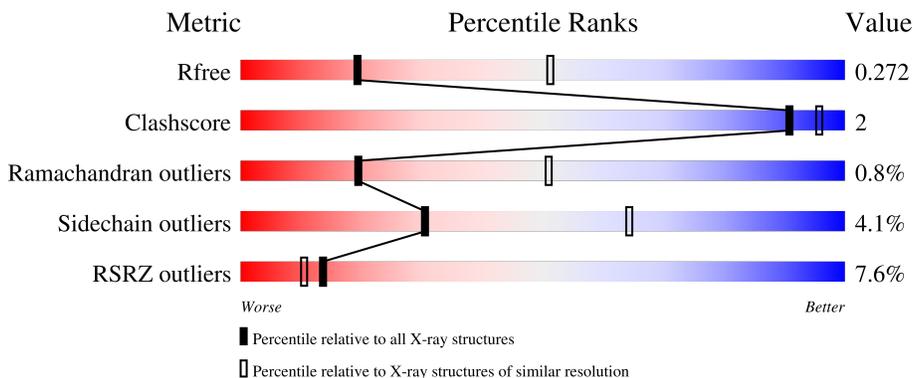
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.90 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.



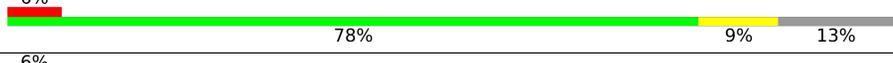
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	1957 (2.90-2.90)
Clashscore	141614	2172 (2.90-2.90)
Ramachandran outliers	138981	2115 (2.90-2.90)
Sidechain outliers	138945	2117 (2.90-2.90)
RSRZ outliers	127900	1906 (2.90-2.90)

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	231	 7% 78% 8% • 13%
1	B	231	 81% 5% • 13%
1	C	231	 3% 79% 8% 13%
1	D	231	 6% 83% • 12%

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Mol	Chain	Length	Quality of chain
1	E	231	 <p>4% 79% 6% 13%</p>
1	F	231	 <p>7% 79% 6% 15%</p>
1	G	231	 <p>6% 78% 9% 13%</p>
1	H	231	 <p>6% 75% 12% 13%</p>
1	I	231	 <p>5% 77% 9% 13%</p>
1	J	231	 <p>6% 77% 11% 13%</p>
1	K	231	 <p>15% 83% 1% 13%</p>
1	L	231	 <p>13% 73% 14% 13%</p>

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 19591 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 Glutathione S-transferase related protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	202	Total 1639	C 1051	N 268	O 312	S 8	0	0	0
1	B	201	Total 1634	C 1048	N 267	O 311	S 8	0	0	0
1	C	201	Total 1634	C 1048	N 267	O 311	S 8	0	0	0
1	D	203	Total 1653	C 1058	N 273	O 314	S 8	0	0	0
1	E	200	Total 1625	C 1042	N 265	O 310	S 8	0	0	0
1	F	197	Total 1601	C 1030	N 258	O 305	S 8	0	0	0
1	G	201	Total 1625	C 1044	N 263	O 310	S 8	0	0	0
1	H	201	Total 1636	C 1048	N 269	O 311	S 8	0	0	0
1	I	200	Total 1629	C 1045	N 266	O 310	S 8	0	0	0
1	J	202	Total 1643	C 1053	N 268	O 314	S 8	0	0	0
1	K	201	Total 1640	C 1051	N 270	O 311	S 8	0	0	0
1	L	200	Total 1618	C 1038	N 263	O 309	S 8	0	0	0

There are 288 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	MET	-	expression tag	UNP Q5R0K1
A	0	VAL	-	expression tag	UNP Q5R0K1
A	208	ALA	-	expression tag	UNP Q5R0K1
A	209	GLU	-	expression tag	UNP Q5R0K1
A	210	ASN	-	expression tag	UNP Q5R0K1

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Chain	Residue	Modelled	Actual	Comment	Reference
A	211	LEU	-	expression tag	UNP Q5R0K1
A	212	TYR	-	expression tag	UNP Q5R0K1
A	213	PHE	-	expression tag	UNP Q5R0K1
A	214	GLN	-	expression tag	UNP Q5R0K1
A	215	SER	-	expression tag	UNP Q5R0K1
A	216	HIS	-	expression tag	UNP Q5R0K1
A	217	HIS	-	expression tag	UNP Q5R0K1
A	218	HIS	-	expression tag	UNP Q5R0K1
A	219	HIS	-	expression tag	UNP Q5R0K1
A	220	HIS	-	expression tag	UNP Q5R0K1
A	221	HIS	-	expression tag	UNP Q5R0K1
A	222	TRP	-	expression tag	UNP Q5R0K1
A	223	SER	-	expression tag	UNP Q5R0K1
A	224	HIS	-	expression tag	UNP Q5R0K1
A	225	PRO	-	expression tag	UNP Q5R0K1
A	226	GLN	-	expression tag	UNP Q5R0K1
A	227	PHE	-	expression tag	UNP Q5R0K1
A	228	GLU	-	expression tag	UNP Q5R0K1
A	229	LYS	-	expression tag	UNP Q5R0K1
B	-1	MET	-	expression tag	UNP Q5R0K1
B	0	VAL	-	expression tag	UNP Q5R0K1
B	208	ALA	-	expression tag	UNP Q5R0K1
B	209	GLU	-	expression tag	UNP Q5R0K1
B	210	ASN	-	expression tag	UNP Q5R0K1
B	211	LEU	-	expression tag	UNP Q5R0K1
B	212	TYR	-	expression tag	UNP Q5R0K1
B	213	PHE	-	expression tag	UNP Q5R0K1
B	214	GLN	-	expression tag	UNP Q5R0K1
B	215	SER	-	expression tag	UNP Q5R0K1
B	216	HIS	-	expression tag	UNP Q5R0K1
B	217	HIS	-	expression tag	UNP Q5R0K1
B	218	HIS	-	expression tag	UNP Q5R0K1
B	219	HIS	-	expression tag	UNP Q5R0K1
B	220	HIS	-	expression tag	UNP Q5R0K1
B	221	HIS	-	expression tag	UNP Q5R0K1
B	222	TRP	-	expression tag	UNP Q5R0K1
B	223	SER	-	expression tag	UNP Q5R0K1
B	224	HIS	-	expression tag	UNP Q5R0K1
B	225	PRO	-	expression tag	UNP Q5R0K1
B	226	GLN	-	expression tag	UNP Q5R0K1
B	227	PHE	-	expression tag	UNP Q5R0K1
B	228	GLU	-	expression tag	UNP Q5R0K1

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Chain	Residue	Modelled	Actual	Comment	Reference
B	229	LYS	-	expression tag	UNP Q5R0K1
C	-1	MET	-	expression tag	UNP Q5R0K1
C	0	VAL	-	expression tag	UNP Q5R0K1
C	208	ALA	-	expression tag	UNP Q5R0K1
C	209	GLU	-	expression tag	UNP Q5R0K1
C	210	ASN	-	expression tag	UNP Q5R0K1
C	211	LEU	-	expression tag	UNP Q5R0K1
C	212	TYR	-	expression tag	UNP Q5R0K1
C	213	PHE	-	expression tag	UNP Q5R0K1
C	214	GLN	-	expression tag	UNP Q5R0K1
C	215	SER	-	expression tag	UNP Q5R0K1
C	216	HIS	-	expression tag	UNP Q5R0K1
C	217	HIS	-	expression tag	UNP Q5R0K1
C	218	HIS	-	expression tag	UNP Q5R0K1
C	219	HIS	-	expression tag	UNP Q5R0K1
C	220	HIS	-	expression tag	UNP Q5R0K1
C	221	HIS	-	expression tag	UNP Q5R0K1
C	222	TRP	-	expression tag	UNP Q5R0K1
C	223	SER	-	expression tag	UNP Q5R0K1
C	224	HIS	-	expression tag	UNP Q5R0K1
C	225	PRO	-	expression tag	UNP Q5R0K1
C	226	GLN	-	expression tag	UNP Q5R0K1
C	227	PHE	-	expression tag	UNP Q5R0K1
C	228	GLU	-	expression tag	UNP Q5R0K1
C	229	LYS	-	expression tag	UNP Q5R0K1
D	-1	MET	-	expression tag	UNP Q5R0K1
D	0	VAL	-	expression tag	UNP Q5R0K1
D	208	ALA	-	expression tag	UNP Q5R0K1
D	209	GLU	-	expression tag	UNP Q5R0K1
D	210	ASN	-	expression tag	UNP Q5R0K1
D	211	LEU	-	expression tag	UNP Q5R0K1
D	212	TYR	-	expression tag	UNP Q5R0K1
D	213	PHE	-	expression tag	UNP Q5R0K1
D	214	GLN	-	expression tag	UNP Q5R0K1
D	215	SER	-	expression tag	UNP Q5R0K1
D	216	HIS	-	expression tag	UNP Q5R0K1
D	217	HIS	-	expression tag	UNP Q5R0K1
D	218	HIS	-	expression tag	UNP Q5R0K1
D	219	HIS	-	expression tag	UNP Q5R0K1
D	220	HIS	-	expression tag	UNP Q5R0K1
D	221	HIS	-	expression tag	UNP Q5R0K1
D	222	TRP	-	expression tag	UNP Q5R0K1

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Chain	Residue	Modelled	Actual	Comment	Reference
D	223	SER	-	expression tag	UNP Q5R0K1
D	224	HIS	-	expression tag	UNP Q5R0K1
D	225	PRO	-	expression tag	UNP Q5R0K1
D	226	GLN	-	expression tag	UNP Q5R0K1
D	227	PHE	-	expression tag	UNP Q5R0K1
D	228	GLU	-	expression tag	UNP Q5R0K1
D	229	LYS	-	expression tag	UNP Q5R0K1
E	-1	MET	-	expression tag	UNP Q5R0K1
E	0	VAL	-	expression tag	UNP Q5R0K1
E	208	ALA	-	expression tag	UNP Q5R0K1
E	209	GLU	-	expression tag	UNP Q5R0K1
E	210	ASN	-	expression tag	UNP Q5R0K1
E	211	LEU	-	expression tag	UNP Q5R0K1
E	212	TYR	-	expression tag	UNP Q5R0K1
E	213	PHE	-	expression tag	UNP Q5R0K1
E	214	GLN	-	expression tag	UNP Q5R0K1
E	215	SER	-	expression tag	UNP Q5R0K1
E	216	HIS	-	expression tag	UNP Q5R0K1
E	217	HIS	-	expression tag	UNP Q5R0K1
E	218	HIS	-	expression tag	UNP Q5R0K1
E	219	HIS	-	expression tag	UNP Q5R0K1
E	220	HIS	-	expression tag	UNP Q5R0K1
E	221	HIS	-	expression tag	UNP Q5R0K1
E	222	TRP	-	expression tag	UNP Q5R0K1
E	223	SER	-	expression tag	UNP Q5R0K1
E	224	HIS	-	expression tag	UNP Q5R0K1
E	225	PRO	-	expression tag	UNP Q5R0K1
E	226	GLN	-	expression tag	UNP Q5R0K1
E	227	PHE	-	expression tag	UNP Q5R0K1
E	228	GLU	-	expression tag	UNP Q5R0K1
E	229	LYS	-	expression tag	UNP Q5R0K1
F	-1	MET	-	expression tag	UNP Q5R0K1
F	0	VAL	-	expression tag	UNP Q5R0K1
F	208	ALA	-	expression tag	UNP Q5R0K1
F	209	GLU	-	expression tag	UNP Q5R0K1
F	210	ASN	-	expression tag	UNP Q5R0K1
F	211	LEU	-	expression tag	UNP Q5R0K1
F	212	TYR	-	expression tag	UNP Q5R0K1
F	213	PHE	-	expression tag	UNP Q5R0K1
F	214	GLN	-	expression tag	UNP Q5R0K1
F	215	SER	-	expression tag	UNP Q5R0K1
F	216	HIS	-	expression tag	UNP Q5R0K1

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Chain	Residue	Modelled	Actual	Comment	Reference
F	217	HIS	-	expression tag	UNP Q5R0K1
F	218	HIS	-	expression tag	UNP Q5R0K1
F	219	HIS	-	expression tag	UNP Q5R0K1
F	220	HIS	-	expression tag	UNP Q5R0K1
F	221	HIS	-	expression tag	UNP Q5R0K1
F	222	TRP	-	expression tag	UNP Q5R0K1
F	223	SER	-	expression tag	UNP Q5R0K1
F	224	HIS	-	expression tag	UNP Q5R0K1
F	225	PRO	-	expression tag	UNP Q5R0K1
F	226	GLN	-	expression tag	UNP Q5R0K1
F	227	PHE	-	expression tag	UNP Q5R0K1
F	228	GLU	-	expression tag	UNP Q5R0K1
F	229	LYS	-	expression tag	UNP Q5R0K1
G	-1	MET	-	expression tag	UNP Q5R0K1
G	0	VAL	-	expression tag	UNP Q5R0K1
G	208	ALA	-	expression tag	UNP Q5R0K1
G	209	GLU	-	expression tag	UNP Q5R0K1
G	210	ASN	-	expression tag	UNP Q5R0K1
G	211	LEU	-	expression tag	UNP Q5R0K1
G	212	TYR	-	expression tag	UNP Q5R0K1
G	213	PHE	-	expression tag	UNP Q5R0K1
G	214	GLN	-	expression tag	UNP Q5R0K1
G	215	SER	-	expression tag	UNP Q5R0K1
G	216	HIS	-	expression tag	UNP Q5R0K1
G	217	HIS	-	expression tag	UNP Q5R0K1
G	218	HIS	-	expression tag	UNP Q5R0K1
G	219	HIS	-	expression tag	UNP Q5R0K1
G	220	HIS	-	expression tag	UNP Q5R0K1
G	221	HIS	-	expression tag	UNP Q5R0K1
G	222	TRP	-	expression tag	UNP Q5R0K1
G	223	SER	-	expression tag	UNP Q5R0K1
G	224	HIS	-	expression tag	UNP Q5R0K1
G	225	PRO	-	expression tag	UNP Q5R0K1
G	226	GLN	-	expression tag	UNP Q5R0K1
G	227	PHE	-	expression tag	UNP Q5R0K1
G	228	GLU	-	expression tag	UNP Q5R0K1
G	229	LYS	-	expression tag	UNP Q5R0K1
H	-1	MET	-	expression tag	UNP Q5R0K1
H	0	VAL	-	expression tag	UNP Q5R0K1
H	208	ALA	-	expression tag	UNP Q5R0K1
H	209	GLU	-	expression tag	UNP Q5R0K1
H	210	ASN	-	expression tag	UNP Q5R0K1

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Chain	Residue	Modelled	Actual	Comment	Reference
H	211	LEU	-	expression tag	UNP Q5R0K1
H	212	TYR	-	expression tag	UNP Q5R0K1
H	213	PHE	-	expression tag	UNP Q5R0K1
H	214	GLN	-	expression tag	UNP Q5R0K1
H	215	SER	-	expression tag	UNP Q5R0K1
H	216	HIS	-	expression tag	UNP Q5R0K1
H	217	HIS	-	expression tag	UNP Q5R0K1
H	218	HIS	-	expression tag	UNP Q5R0K1
H	219	HIS	-	expression tag	UNP Q5R0K1
H	220	HIS	-	expression tag	UNP Q5R0K1
H	221	HIS	-	expression tag	UNP Q5R0K1
H	222	TRP	-	expression tag	UNP Q5R0K1
H	223	SER	-	expression tag	UNP Q5R0K1
H	224	HIS	-	expression tag	UNP Q5R0K1
H	225	PRO	-	expression tag	UNP Q5R0K1
H	226	GLN	-	expression tag	UNP Q5R0K1
H	227	PHE	-	expression tag	UNP Q5R0K1
H	228	GLU	-	expression tag	UNP Q5R0K1
H	229	LYS	-	expression tag	UNP Q5R0K1
I	-1	MET	-	expression tag	UNP Q5R0K1
I	0	VAL	-	expression tag	UNP Q5R0K1
I	208	ALA	-	expression tag	UNP Q5R0K1
I	209	GLU	-	expression tag	UNP Q5R0K1
I	210	ASN	-	expression tag	UNP Q5R0K1
I	211	LEU	-	expression tag	UNP Q5R0K1
I	212	TYR	-	expression tag	UNP Q5R0K1
I	213	PHE	-	expression tag	UNP Q5R0K1
I	214	GLN	-	expression tag	UNP Q5R0K1
I	215	SER	-	expression tag	UNP Q5R0K1
I	216	HIS	-	expression tag	UNP Q5R0K1
I	217	HIS	-	expression tag	UNP Q5R0K1
I	218	HIS	-	expression tag	UNP Q5R0K1
I	219	HIS	-	expression tag	UNP Q5R0K1
I	220	HIS	-	expression tag	UNP Q5R0K1
I	221	HIS	-	expression tag	UNP Q5R0K1
I	222	TRP	-	expression tag	UNP Q5R0K1
I	223	SER	-	expression tag	UNP Q5R0K1
I	224	HIS	-	expression tag	UNP Q5R0K1
I	225	PRO	-	expression tag	UNP Q5R0K1
I	226	GLN	-	expression tag	UNP Q5R0K1
I	227	PHE	-	expression tag	UNP Q5R0K1
I	228	GLU	-	expression tag	UNP Q5R0K1

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Chain	Residue	Modelled	Actual	Comment	Reference
I	229	LYS	-	expression tag	UNP Q5R0K1
J	-1	MET	-	expression tag	UNP Q5R0K1
J	0	VAL	-	expression tag	UNP Q5R0K1
J	208	ALA	-	expression tag	UNP Q5R0K1
J	209	GLU	-	expression tag	UNP Q5R0K1
J	210	ASN	-	expression tag	UNP Q5R0K1
J	211	LEU	-	expression tag	UNP Q5R0K1
J	212	TYR	-	expression tag	UNP Q5R0K1
J	213	PHE	-	expression tag	UNP Q5R0K1
J	214	GLN	-	expression tag	UNP Q5R0K1
J	215	SER	-	expression tag	UNP Q5R0K1
J	216	HIS	-	expression tag	UNP Q5R0K1
J	217	HIS	-	expression tag	UNP Q5R0K1
J	218	HIS	-	expression tag	UNP Q5R0K1
J	219	HIS	-	expression tag	UNP Q5R0K1
J	220	HIS	-	expression tag	UNP Q5R0K1
J	221	HIS	-	expression tag	UNP Q5R0K1
J	222	TRP	-	expression tag	UNP Q5R0K1
J	223	SER	-	expression tag	UNP Q5R0K1
J	224	HIS	-	expression tag	UNP Q5R0K1
J	225	PRO	-	expression tag	UNP Q5R0K1
J	226	GLN	-	expression tag	UNP Q5R0K1
J	227	PHE	-	expression tag	UNP Q5R0K1
J	228	GLU	-	expression tag	UNP Q5R0K1
J	229	LYS	-	expression tag	UNP Q5R0K1
K	-1	MET	-	expression tag	UNP Q5R0K1
K	0	VAL	-	expression tag	UNP Q5R0K1
K	208	ALA	-	expression tag	UNP Q5R0K1
K	209	GLU	-	expression tag	UNP Q5R0K1
K	210	ASN	-	expression tag	UNP Q5R0K1
K	211	LEU	-	expression tag	UNP Q5R0K1
K	212	TYR	-	expression tag	UNP Q5R0K1
K	213	PHE	-	expression tag	UNP Q5R0K1
K	214	GLN	-	expression tag	UNP Q5R0K1
K	215	SER	-	expression tag	UNP Q5R0K1
K	216	HIS	-	expression tag	UNP Q5R0K1
K	217	HIS	-	expression tag	UNP Q5R0K1
K	218	HIS	-	expression tag	UNP Q5R0K1
K	219	HIS	-	expression tag	UNP Q5R0K1
K	220	HIS	-	expression tag	UNP Q5R0K1
K	221	HIS	-	expression tag	UNP Q5R0K1
K	222	TRP	-	expression tag	UNP Q5R0K1

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Chain	Residue	Modelled	Actual	Comment	Reference
K	223	SER	-	expression tag	UNP Q5R0K1
K	224	HIS	-	expression tag	UNP Q5R0K1
K	225	PRO	-	expression tag	UNP Q5R0K1
K	226	GLN	-	expression tag	UNP Q5R0K1
K	227	PHE	-	expression tag	UNP Q5R0K1
K	228	GLU	-	expression tag	UNP Q5R0K1
K	229	LYS	-	expression tag	UNP Q5R0K1
L	-1	MET	-	expression tag	UNP Q5R0K1
L	0	VAL	-	expression tag	UNP Q5R0K1
L	208	ALA	-	expression tag	UNP Q5R0K1
L	209	GLU	-	expression tag	UNP Q5R0K1
L	210	ASN	-	expression tag	UNP Q5R0K1
L	211	LEU	-	expression tag	UNP Q5R0K1
L	212	TYR	-	expression tag	UNP Q5R0K1
L	213	PHE	-	expression tag	UNP Q5R0K1
L	214	GLN	-	expression tag	UNP Q5R0K1
L	215	SER	-	expression tag	UNP Q5R0K1
L	216	HIS	-	expression tag	UNP Q5R0K1
L	217	HIS	-	expression tag	UNP Q5R0K1
L	218	HIS	-	expression tag	UNP Q5R0K1
L	219	HIS	-	expression tag	UNP Q5R0K1
L	220	HIS	-	expression tag	UNP Q5R0K1
L	221	HIS	-	expression tag	UNP Q5R0K1
L	222	TRP	-	expression tag	UNP Q5R0K1
L	223	SER	-	expression tag	UNP Q5R0K1
L	224	HIS	-	expression tag	UNP Q5R0K1
L	225	PRO	-	expression tag	UNP Q5R0K1
L	226	GLN	-	expression tag	UNP Q5R0K1
L	227	PHE	-	expression tag	UNP Q5R0K1
L	228	GLU	-	expression tag	UNP Q5R0K1
L	229	LYS	-	expression tag	UNP Q5R0K1

- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	4	Total O 4 4	0	0
2	B	3	Total O 3 3	0	0
2	C	1	Total O 1 1	0	0
2	E	1	Total O 1 1	0	0

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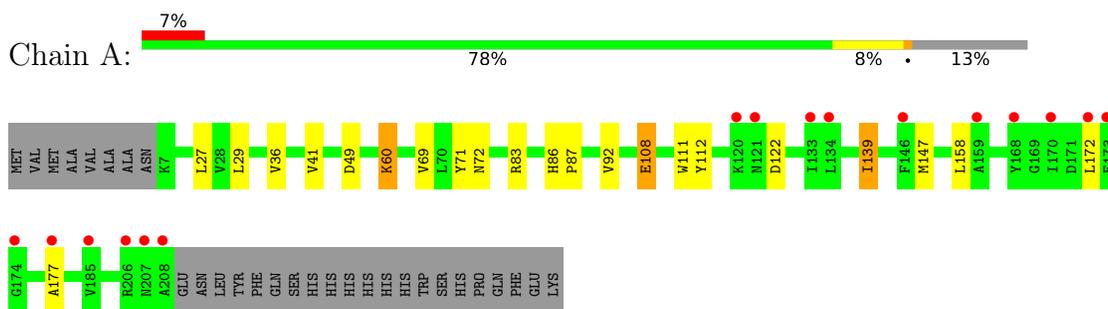
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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	F	1	Total O 1 1	0	0
2	I	2	Total O 2 2	0	0
2	J	1	Total O 1 1	0	0
2	L	1	Total O 1 1	0	0

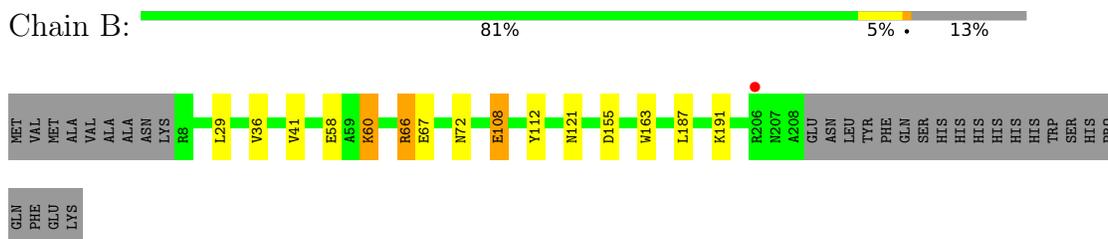
3 Residue-property plots [i](#)

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.

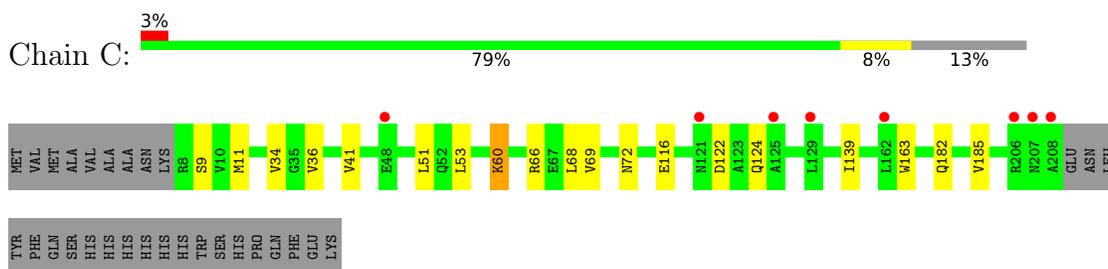
- Molecule 1: Glutathione S-transferase related protein



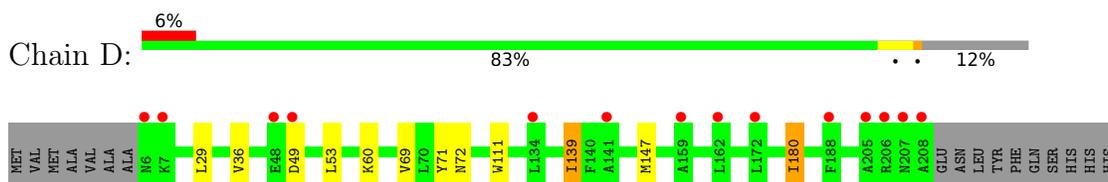
- Molecule 1: Glutathione S-transferase related protein



- Molecule 1: Glutathione S-transferase related protein

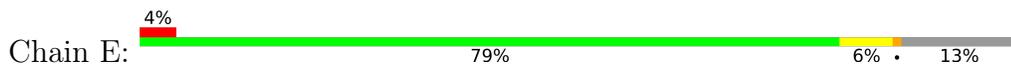


- Molecule 1: Glutathione S-transferase related protein



HIS
HIS
HIS
TRP
SER
HIS
PRO
GLN
GLU
LYS

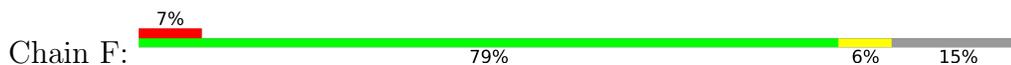
• Molecule 1: Glutathione S-transferase related protein



MET VAL MET MET VAL ALA ALA ALA ASN LYS HIS S9 LEU L20 Q24 L27 V28 L29 V34 G35 V36 V41 T42 S45 E48 D49 L50 L51 Q52 L53 K60 R66 E67 L68 V69 N72 I75 R83 I118 A123 Y157 I170 A208 ASN

LEU TYR PHE GLN SER HIS HIS HIS HIS HIS TRP SER HIS PRO GLN PHE GLU LYS

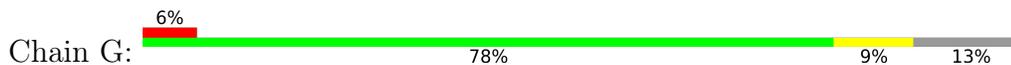
• Molecule 1: Glutathione S-transferase related protein



MET VAL MET VAL VAL ALA ALA ASN LYS ARG S9 Q24 L27 L29 V36 V41 E44 S45 L53 K60 V69 N72 W111 N121 I134 S135 L136 I139 V154 Y157 L158 L161 L172 E173 G174 Q175 G176 A177 K178 E179 I180

V185 E199 A205 ARG ASN ALA GLU ASN ASN LEU TYR PHE GLN SER HIS HIS HIS HIS TRP SER PRO GLN PHE LYS

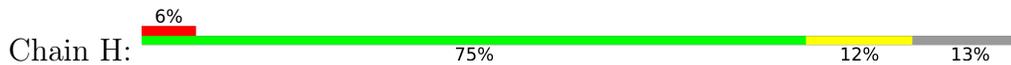
• Molecule 1: Glutathione S-transferase related protein



MET VAL MET ALA VAL ALA ALA ASN LYS RB V36 E44 L53 K60 R66 E67 L68 V69 L70 N72 A73 Q74 E78 R83 E108 Y112 A137 F138 I139 F140 A141 A159 L162 L165 L172 E173 G174 Q175 G176 A177 K178 E179 I180 D195

L204 A205 R206 N207 A208 GLU ASN LEU TYR PHE GLN SER HIS HIS HIS HIS HIS TRP SER PRO GLN PHE LYS

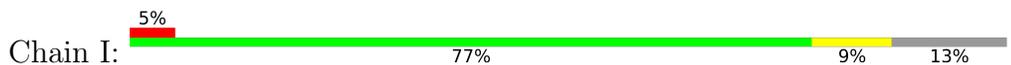
• Molecule 1: Glutathione S-transferase related protein

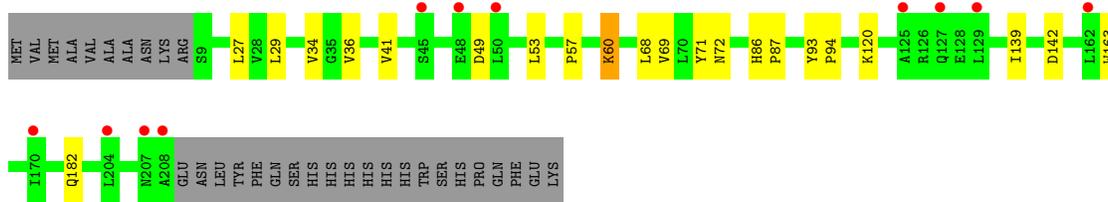


MET VAL MET VAL VAL ALA ALA ASN LYS R8 Q24 L27 V28 L29 V36 E44 D49 L53 K60 V64 D65 R66 Y71 N72 R83 V92 Y93 P94 E108 W111 Y112 K120 N121 D122 A123 Q124 K130 I133 L134 I139 Y157 L158

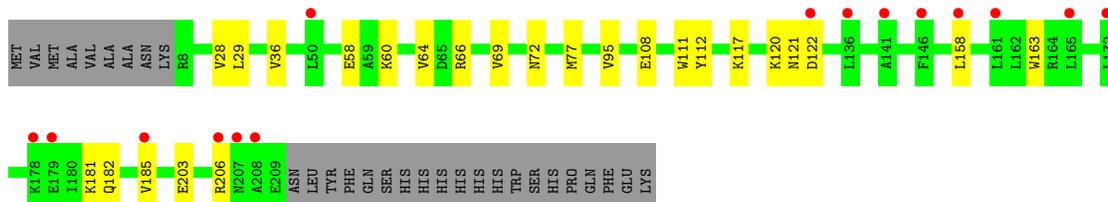
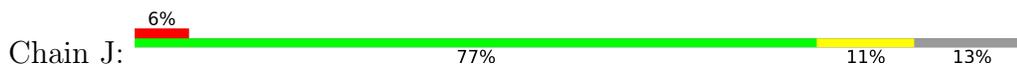
A159 L162 W163 R164 L165 L172 A177 K181 V185 K191 A208 GLU ASN LEU TYR PHE GLN SER HIS HIS HIS HIS HIS TRP SER PRO GLN PHE GLU LYS

• Molecule 1: Glutathione S-transferase related protein

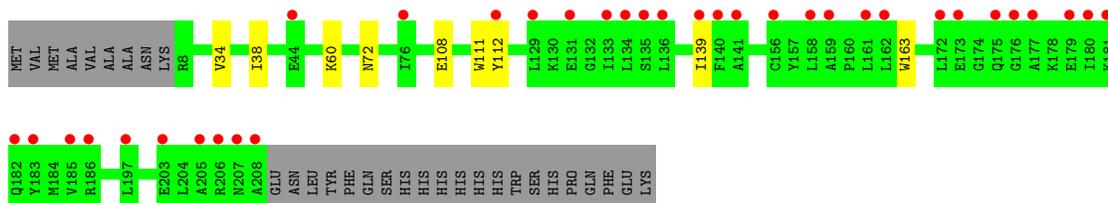
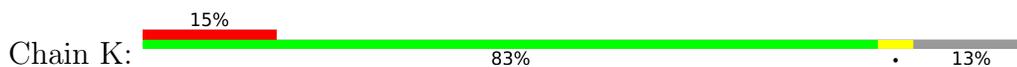




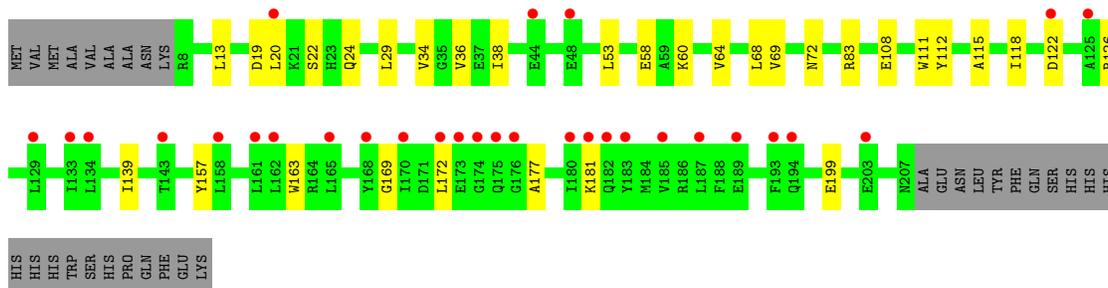
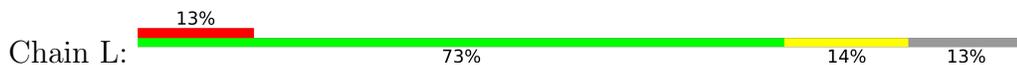
● Molecule 1: Glutathione S-transferase related protein



● Molecule 1: Glutathione S-transferase related protein



● Molecule 1: Glutathione S-transferase related protein



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	91.36Å 156.99Å 244.12Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	40.00 – 2.90 48.10 – 2.80	Depositor EDS
% Data completeness (in resolution range)	99.7 (40.00-2.90) 99.1 (48.10-2.80)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.08	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.13 (at 2.81Å)	Xtrriage
Refinement program	REFMAC 5.6.0117	Depositor
R, R_{free}	0.207 , 0.275 0.210 , 0.272	Depositor DCC
R_{free} test set	2617 reflections (3.02%)	wwPDB-VP
Wilson B-factor (Å ²)	87.9	Xtrriage
Anisotropy	0.166	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 75.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	19591	wwPDB-VP
Average B, all atoms (Å ²)	104.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.00% 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 [i](#)

5.1 Standard geometry [i](#)

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	0.43	1/1675 (0.1%)	0.62	0/2271
1	B	0.43	1/1670 (0.1%)	0.64	0/2264
1	C	0.42	1/1670 (0.1%)	0.62	0/2264
1	D	0.43	1/1689 (0.1%)	0.61	0/2289
1	E	0.43	0/1661	0.63	0/2253
1	F	0.43	1/1637 (0.1%)	0.60	0/2220
1	G	0.43	0/1661	0.62	0/2253
1	H	0.44	2/1672 (0.1%)	0.61	0/2267
1	I	0.42	1/1665 (0.1%)	0.61	0/2257
1	J	0.43	2/1679 (0.1%)	0.59	0/2276
1	K	0.43	2/1676 (0.1%)	0.60	0/2271
1	L	0.43	2/1654 (0.1%)	0.59	0/2245
All	All	0.43	14/20009 (0.1%)	0.61	0/27130

All (14) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	J	111	TRP	CD2-CE2	5.29	1.47	1.41
1	A	111	TRP	CD2-CE2	5.27	1.47	1.41
1	F	111	TRP	CD2-CE2	5.26	1.47	1.41
1	D	111	TRP	CD2-CE2	5.19	1.47	1.41
1	H	163	TRP	CD2-CE2	5.18	1.47	1.41
1	L	163	TRP	CD2-CE2	5.17	1.47	1.41
1	K	111	TRP	CD2-CE2	5.11	1.47	1.41
1	C	163	TRP	CD2-CE2	5.09	1.47	1.41
1	H	111	TRP	CD2-CE2	5.08	1.47	1.41
1	B	163	TRP	CD2-CE2	5.07	1.47	1.41
1	K	163	TRP	CD2-CE2	5.07	1.47	1.41
1	L	111	TRP	CD2-CE2	5.07	1.47	1.41
1	I	163	TRP	CD2-CE2	5.06	1.47	1.41
1	J	163	TRP	CD2-CE2	5.04	1.47	1.41

There are no bond angle outliers.

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	1639	0	1627	6	0
1	B	1634	0	1625	5	0
1	C	1634	0	1625	3	0
1	D	1653	0	1644	4	0
1	E	1625	0	1612	7	0
1	F	1601	0	1593	4	0
1	G	1625	0	1610	5	0
1	H	1636	0	1625	6	0
1	I	1629	0	1623	7	0
1	J	1643	0	1631	10	0
1	K	1640	0	1636	2	0
1	L	1618	0	1594	11	0
2	A	4	0	0	0	0
2	B	3	0	0	0	0
2	C	1	0	0	0	0
2	E	1	0	0	0	0
2	F	1	0	0	0	0
2	I	2	0	0	0	0
2	J	1	0	0	0	0
2	L	1	0	0	0	0
All	All	19591	0	19445	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 2.

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:J:121:ASN:O	1:J:121:ASN:ND2	2.30	0.64
1:I:29:LEU:HD13	1:I:36:VAL:HG11	1.85	0.57
1:E:29:LEU:HD13	1:E:36:VAL:HG11	1.87	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:J:120:LYS:O	1:J:121:ASN:HB3	2.06	0.56
1:L:29:LEU:HD13	1:L:36:VAL:HG11	1.88	0.56
1:A:172:LEU:HB2	1:A:177:ALA:HB1	1.88	0.55
1:B:29:LEU:HD13	1:B:36:VAL:HG11	1.88	0.55
1:A:108:GLU:HA	1:A:112:TYR:HB2	1.89	0.55
1:A:41:VAL:HG23	1:A:60:LYS:HE3	1.89	0.54
1:C:53:LEU:HB3	1:C:69:VAL:HG21	1.89	0.53
1:J:117:LYS:HB3	1:J:122:ASP:OD2	2.08	0.53
1:D:139:ILE:HD13	1:D:147:MET:HG3	1.91	0.53
1:A:29:LEU:HD13	1:A:36:VAL:HG11	1.89	0.52
1:J:29:LEU:HD13	1:J:36:VAL:HG11	1.91	0.52
1:L:53:LEU:HB3	1:L:69:VAL:HG21	1.92	0.52
1:E:42:THR:HG23	1:E:45:SER:H	1.76	0.51
1:D:29:LEU:HD13	1:D:36:VAL:HG11	1.94	0.50
1:I:41:VAL:HG23	1:I:60:LYS:HE3	1.93	0.50
1:F:29:LEU:HD13	1:F:36:VAL:HG11	1.93	0.50
1:D:53:LEU:HB3	1:D:69:VAL:HG21	1.94	0.50
1:G:108:GLU:HA	1:G:112:TYR:HB2	1.94	0.49
1:I:53:LEU:HB3	1:I:69:VAL:HG21	1.93	0.49
1:F:53:LEU:HB3	1:F:69:VAL:HG21	1.93	0.49
1:L:24:GLN:HE21	1:L:157:TYR:HA	1.78	0.48
1:G:53:LEU:HB3	1:G:69:VAL:HG21	1.95	0.48
1:H:24:GLN:HE21	1:H:157:TYR:HA	1.79	0.47
1:J:108:GLU:HA	1:J:112:TYR:HB2	1.96	0.47
1:E:34:VAL:HG12	1:L:34:VAL:HG13	1.97	0.47
1:A:139:ILE:HD13	1:A:147:MET:HG3	1.96	0.46
1:C:11:MET:HB2	1:C:36:VAL:HG12	1.97	0.46
1:L:108:GLU:HA	1:L:112:TYR:HB2	1.97	0.46
1:E:72:ASN:HB3	1:E:75:ILE:HD12	1.98	0.45
1:L:115:ALA:HA	1:L:118:ILE:HD12	1.98	0.45
1:C:41:VAL:HG23	1:C:60:LYS:HE3	1.98	0.45
1:H:29:LEU:HD13	1:H:36:VAL:HG11	1.98	0.45
1:L:126:ARG:HE	1:L:169:GLY:HA3	1.82	0.45
1:B:41:VAL:HG23	1:B:60:LYS:HE3	1.98	0.45
1:I:34:VAL:HG12	1:K:34:VAL:HG13	1.99	0.45
1:B:155:ASP:HB3	1:B:187:LEU:HD11	2.00	0.44
1:A:86:HIS:HA	1:A:87:PRO:HA	1.91	0.44
1:H:172:LEU:HB3	1:H:177:ALA:HB1	1.98	0.44
1:J:203:GLU:HG3	1:J:206:ARG:HH11	1.82	0.44
1:L:172:LEU:HD22	1:L:177:ALA:HB1	2.00	0.44
1:F:41:VAL:HG23	1:F:60:LYS:HE3	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:I:93:TYR:HA	1:I:94:PRO:HD3	1.87	0.43
1:E:41:VAL:HG23	1:E:60:LYS:HE3	1.99	0.43
1:J:28:VAL:HG21	1:J:77:MET:HG2	2.01	0.42
1:L:64:VAL:HG22	1:L:69:VAL:HG23	2.01	0.42
1:E:24:GLN:HE21	1:E:157:TYR:HA	1.84	0.42
1:I:68:LEU:HD22	1:J:95:VAL:HG13	2.01	0.42
1:K:108:GLU:HA	1:K:112:TYR:HB2	2.02	0.42
1:G:162:LEU:HA	1:G:165:LEU:HD13	2.01	0.42
1:L:13:LEU:HD23	1:L:38:ILE:HG12	2.01	0.42
1:H:53:LEU:HD13	1:H:64:VAL:HG21	2.01	0.42
1:B:66:ARG:HB2	1:B:67:GLU:H	1.73	0.42
1:G:137:ALA:HA	1:G:140:PHE:HD2	1.85	0.42
1:J:120:LYS:O	1:J:121:ASN:CB	2.65	0.42
1:G:74:GLN:HE21	1:G:78:GLU:HG3	1.84	0.41
1:H:93:TYR:HA	1:H:94:PRO:HD3	1.92	0.41
1:J:64:VAL:HG22	1:J:69:VAL:HG23	2.01	0.41
1:I:86:HIS:HA	1:I:87:PRO:HA	1.94	0.41
1:F:24:GLN:HE21	1:F:157:TYR:HA	1.85	0.41
1:H:108:GLU:HA	1:H:112:TYR:HB2	2.02	0.41
1:D:180:ILE:H	1:D:180:ILE:HG12	1.69	0.41
1:L:19:ASP:HB3	1:L:22:SER:HB2	2.03	0.41
1:E:53:LEU:HB3	1:E:69:VAL:HG21	2.03	0.40
1:B:108:GLU:HA	1:B:112:TYR:HB2	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	200/231 (87%)	191 (96%)	8 (4%)	1 (0%)	29 61
1	B	199/231 (86%)	195 (98%)	2 (1%)	2 (1%)	15 45

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	C	199/231 (86%)	189 (95%)	8 (4%)	2 (1%)	15	45
1	D	201/231 (87%)	195 (97%)	5 (2%)	1 (0%)	29	61
1	E	198/231 (86%)	194 (98%)	2 (1%)	2 (1%)	15	45
1	F	195/231 (84%)	189 (97%)	5 (3%)	1 (0%)	29	61
1	G	199/231 (86%)	192 (96%)	5 (2%)	2 (1%)	15	45
1	H	199/231 (86%)	193 (97%)	3 (2%)	3 (2%)	10	34
1	I	198/231 (86%)	191 (96%)	5 (2%)	2 (1%)	15	45
1	J	200/231 (87%)	194 (97%)	4 (2%)	2 (1%)	15	45
1	K	199/231 (86%)	188 (94%)	10 (5%)	1 (0%)	29	61
1	L	198/231 (86%)	190 (96%)	7 (4%)	1 (0%)	29	61
All	All	2385/2772 (86%)	2301 (96%)	64 (3%)	20 (1%)	19	51

All (20) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	72	ASN
1	B	72	ASN
1	E	66	ARG
1	F	72	ASN
1	G	66	ARG
1	H	66	ARG
1	H	120	LYS
1	J	66	ARG
1	K	72	ASN
1	L	72	ASN
1	B	66	ARG
1	D	72	ASN
1	E	72	ASN
1	G	72	ASN
1	H	72	ASN
1	I	72	ASN
1	J	72	ASN
1	C	66	ARG
1	C	72	ASN
1	I	120	LYS

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	178/206 (86%)	167 (94%)	11 (6%)	18	47
1	B	178/206 (86%)	173 (97%)	5 (3%)	43	76
1	C	178/206 (86%)	167 (94%)	11 (6%)	18	47
1	D	180/206 (87%)	175 (97%)	5 (3%)	43	76
1	E	177/206 (86%)	172 (97%)	5 (3%)	43	76
1	F	175/206 (85%)	170 (97%)	5 (3%)	42	76
1	G	176/206 (85%)	166 (94%)	10 (6%)	20	51
1	H	178/206 (86%)	168 (94%)	10 (6%)	21	52
1	I	178/206 (86%)	170 (96%)	8 (4%)	27	61
1	J	179/206 (87%)	173 (97%)	6 (3%)	37	71
1	K	179/206 (87%)	176 (98%)	3 (2%)	60	86
1	L	175/206 (85%)	166 (95%)	9 (5%)	24	56
All	All	2131/2472 (86%)	2043 (96%)	88 (4%)	30	64

All (88) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	27	LEU
1	A	49	ASP
1	A	60	LYS
1	A	69	VAL
1	A	71	TYR
1	A	83	ARG
1	A	92	VAL
1	A	108	GLU
1	A	122	ASP
1	A	139	ILE
1	A	158	LEU
1	B	58	GLU
1	B	60	LYS
1	B	108	GLU

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Mol	Chain	Res	Type
1	B	121	ASN
1	B	191	LYS
1	C	9	SER
1	C	34	VAL
1	C	51	LEU
1	C	60	LYS
1	C	68	LEU
1	C	116	GLU
1	C	122	ASP
1	C	124	GLN
1	C	139	ILE
1	C	182	GLN
1	C	185	VAL
1	D	49	ASP
1	D	60	LYS
1	D	71	TYR
1	D	139	ILE
1	D	180	ILE
1	E	27	LEU
1	E	42	THR
1	E	49	ASP
1	E	60	LYS
1	E	83	ARG
1	F	27	LEU
1	F	60	LYS
1	F	121	ASN
1	F	139	ILE
1	F	158	LEU
1	G	36	VAL
1	G	60	LYS
1	G	68	LEU
1	G	71	TYR
1	G	83	ARG
1	G	108	GLU
1	G	139	ILE
1	G	172	LEU
1	G	175	GLN
1	G	195	ASP
1	H	8	ARG
1	H	27	LEU
1	H	49	ASP
1	H	60	LYS

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Mol	Chain	Res	Type
1	H	71	TYR
1	H	83	ARG
1	H	92	VAL
1	H	122	ASP
1	H	139	ILE
1	H	191	LYS
1	I	27	LEU
1	I	49	ASP
1	I	57	PRO
1	I	60	LYS
1	I	71	TYR
1	I	139	ILE
1	I	142	ASP
1	I	182	GLN
1	J	58	GLU
1	J	60	LYS
1	J	158	LEU
1	J	181	LYS
1	J	182	GLN
1	J	185	VAL
1	K	38	ILE
1	K	60	LYS
1	K	139	ILE
1	L	20	LEU
1	L	58	GLU
1	L	60	LYS
1	L	68	LEU
1	L	83	ARG
1	L	122	ASP
1	L	139	ILE
1	L	181	LYS
1	L	199	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	24	GLN
1	B	24	GLN
1	B	127	GLN
1	E	24	GLN
1	F	24	GLN
1	F	182	GLN

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Mol	Chain	Res	Type
1	G	182	GLN
1	H	24	GLN
1	H	119	GLN
1	I	24	GLN
1	I	119	GLN
1	J	52	GLN
1	J	121	ASN
1	J	182	GLN
1	K	182	GLN
1	L	24	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	202/231 (87%)	0.60	16 (7%) 12 10	57, 94, 148, 176	0
1	B	201/231 (87%)	0.31	1 (0%) 91 91	47, 77, 123, 184	0
1	C	201/231 (87%)	0.39	8 (3%) 38 33	56, 93, 142, 173	0
1	D	203/231 (87%)	0.56	14 (6%) 16 13	63, 95, 144, 173	0
1	E	200/231 (86%)	0.50	9 (4%) 33 29	56, 84, 140, 184	0
1	F	197/231 (85%)	0.63	16 (8%) 12 9	59, 103, 151, 169	0
1	G	201/231 (87%)	0.69	14 (6%) 16 12	55, 100, 149, 180	0
1	H	201/231 (87%)	0.52	15 (7%) 14 11	69, 105, 150, 169	0
1	I	200/231 (86%)	0.46	11 (5%) 25 21	67, 101, 147, 188	0
1	J	202/231 (87%)	0.53	15 (7%) 14 11	68, 104, 149, 187	0
1	K	201/231 (87%)	0.96	35 (17%) 1 1	76, 124, 163, 180	0
1	L	200/231 (86%)	0.85	30 (15%) 2 1	76, 126, 165, 182	0
All	All	2409/2772 (86%)	0.58	184 (7%) 13 10	47, 100, 155, 188	0

All (184) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	K	208	ALA	8.0
1	G	207	ASN	6.9
1	L	183	TYR	6.3
1	D	6	ASN	6.0
1	K	185	VAL	6.0
1	L	168	TYR	5.8
1	A	121	ASN	5.8
1	K	172	LEU	5.5
1	G	173	GLU	5.3
1	L	162	LEU	5.2
1	A	173	GLU	4.9

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Mol	Chain	Res	Type	RSRZ
1	C	207	ASN	4.8
1	D	206	ARG	4.7
1	C	121	ASN	4.5
1	F	176	GLY	4.5
1	E	48	GLU	4.4
1	L	175	GLN	4.4
1	L	125	ALA	4.4
1	F	180	ILE	4.3
1	L	165	LEU	4.2
1	L	176	GLY	4.2
1	L	158	LEU	4.1
1	L	172	LEU	4.1
1	J	208	ALA	4.1
1	J	207	ASN	4.1
1	L	129	LEU	4.0
1	H	134	LEU	4.0
1	G	208	ALA	3.9
1	J	136	LEU	3.9
1	G	174	GLY	3.8
1	I	45	SER	3.8
1	G	141	ALA	3.8
1	G	140	PHE	3.7
1	D	48	GLU	3.7
1	D	205	ALA	3.7
1	K	156	CYS	3.6
1	D	208	ALA	3.6
1	A	207	ASN	3.6
1	L	161	LEU	3.4
1	K	131	GLU	3.4
1	E	123	ALA	3.4
1	L	44	GLU	3.4
1	L	133	ILE	3.4
1	K	181	LYS	3.4
1	K	206	ARG	3.3
1	K	162	LEU	3.3
1	K	180	ILE	3.3
1	K	44	GLU	3.2
1	L	187	LEU	3.2
1	K	183	TYR	3.2
1	J	185	VAL	3.2
1	I	207	ASN	3.2
1	L	173	GLU	3.2

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Mol	Chain	Res	Type	RSRZ
1	A	174	GLY	3.2
1	G	204	LEU	3.2
1	I	208	ALA	3.2
1	E	50	LEU	3.2
1	D	172	LEU	3.2
1	A	134	LEU	3.2
1	I	127	GLN	3.1
1	G	206	ARG	3.1
1	D	162	LEU	3.1
1	C	206	ARG	3.1
1	L	122	ASP	3.1
1	A	172	LEU	3.0
1	K	129	LEU	3.0
1	A	133	ILE	3.0
1	D	49	ASP	2.9
1	E	45	SER	2.9
1	K	175	GLN	2.9
1	F	161	LEU	2.9
1	F	177	ALA	2.9
1	F	44	GLU	2.9
1	H	121	ASN	2.9
1	L	48	GLU	2.9
1	L	181	LYS	2.9
1	A	146	PHE	2.9
1	A	208	ALA	2.8
1	F	174	GLY	2.8
1	D	207	ASN	2.8
1	K	133	ILE	2.8
1	F	185	VAL	2.8
1	G	44	GLU	2.8
1	K	203	GLU	2.8
1	K	140	PHE	2.8
1	L	185	VAL	2.8
1	F	199	GLU	2.7
1	A	168	TYR	2.7
1	I	204	LEU	2.7
1	K	134	LEU	2.7
1	E	67	GLU	2.7
1	K	186	ARG	2.7
1	L	134	LEU	2.7
1	D	141	ALA	2.7
1	K	177	ALA	2.7

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Mol	Chain	Res	Type	RSRZ
1	F	45	SER	2.7
1	A	170	ILE	2.7
1	F	172	LEU	2.7
1	K	136	LEU	2.6
1	J	178	LYS	2.6
1	I	170	ILE	2.6
1	L	174	GLY	2.6
1	C	125	ALA	2.6
1	G	159	ALA	2.6
1	B	206	ARG	2.6
1	H	133	ILE	2.6
1	D	159	ALA	2.6
1	K	182	GLN	2.6
1	F	134	LEU	2.6
1	J	172	LEU	2.6
1	G	178	LYS	2.5
1	K	141	ALA	2.5
1	L	180	ILE	2.5
1	J	179	GLU	2.5
1	K	173	GLU	2.5
1	K	179	GLU	2.5
1	I	50	LEU	2.5
1	I	129	LEU	2.5
1	D	134	LEU	2.4
1	K	139	ILE	2.4
1	L	143	THR	2.4
1	K	76	ILE	2.4
1	K	205	ALA	2.4
1	J	206	ARG	2.4
1	F	121	ASN	2.4
1	C	48	GLU	2.4
1	G	176	GLY	2.4
1	K	197	LEU	2.4
1	E	20	LEU	2.3
1	I	48	GLU	2.3
1	K	176	GLY	2.3
1	L	193	PHE	2.3
1	H	53	LEU	2.3
1	J	122	ASP	2.3
1	D	188	PHE	2.3
1	C	162	LEU	2.3
1	F	136	LEU	2.3

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Mol	Chain	Res	Type	RSRZ
1	H	165	LEU	2.3
1	K	161	LEU	2.3
1	A	206	ARG	2.3
1	C	129	LEU	2.3
1	H	162	LEU	2.3
1	J	141	ALA	2.3
1	H	181	LYS	2.3
1	A	177	ALA	2.3
1	J	158	LEU	2.3
1	L	203	GLU	2.2
1	C	208	ALA	2.2
1	G	175	GLN	2.2
1	H	185	VAL	2.2
1	H	158	LEU	2.2
1	I	125	ALA	2.2
1	F	173	GLU	2.2
1	H	208	ALA	2.2
1	A	120	LYS	2.2
1	L	170	ILE	2.2
1	I	162	LEU	2.2
1	J	161	LEU	2.2
1	A	159	ALA	2.2
1	J	146	PHE	2.2
1	E	51	LEU	2.1
1	L	20	LEU	2.1
1	A	185	VAL	2.1
1	L	182	GLN	2.1
1	H	44	GLU	2.1
1	G	180	ILE	2.1
1	D	7	LYS	2.1
1	F	178	LYS	2.1
1	H	130	LYS	2.1
1	K	207	ASN	2.1
1	J	165	LEU	2.1
1	H	124	GLN	2.0
1	E	170	ILE	2.0
1	L	189	GLU	2.0
1	K	159	ALA	2.0
1	K	158	LEU	2.0
1	E	118	ILE	2.0
1	H	159	ALA	2.0
1	K	112	TYR	2.0

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Mol	Chain	Res	Type	RSRZ
1	F	154	VAL	2.0
1	J	50	LEU	2.0
1	L	194	GLN	2.0
1	H	177	ALA	2.0
1	K	135	SER	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.