



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 03:40 PM UTC

PDB ID : 6UTL / pdb_00006utl
Title : Yeast Thiol Specific antioxidant 2 with C171S mutation and catalytic cysteine alkylated with iodoacetamide
Authors : Tairum, C.A.; Bannitz-Fernandes, R.; Tonoli, C.C.C.; Murakami, M.T.; de Oliveira, M.A.; Netto, L.E.S.
Deposited on : 2019-10-29
Resolution : 2.60 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

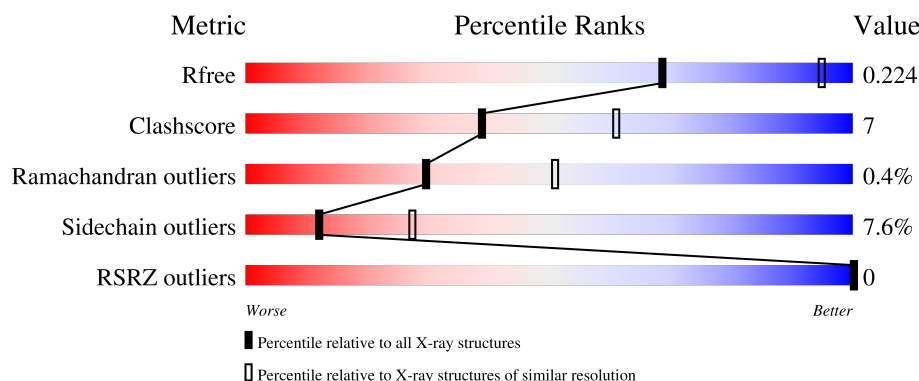
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.60 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	4008 (2.60-2.60)
Clashscore	190562	4347 (2.60-2.60)
Ramachandran outliers	187476	4277 (2.60-2.60)
Sidechain outliers	187428	4277 (2.60-2.60)
RSRZ outliers	180081	4008 (2.60-2.60)

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	216	 60% 15% • 23%
1	B	216	 60% 15% • 24%
1	C	216	 62% 12% • • 23%
1	D	216	 62% 12% • • 24%
1	E	216	 61% 12% • 24%

Continued on next page...

Continued from previous page...

Mol	Chain	Length	Quality of chain
1	F	216	
1	G	216	
1	H	216	
1	I	216	
1	J	216	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 12817 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 Peroxiredoxin TSA2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	167	Total	C	N	O	S	0	0	0
			1286	831	213	241	1			
1	B	165	Total	C	N	O	S	0	0	0
			1266	817	212	236	1			
1	C	166	Total	C	N	O	S	0	0	0
			1284	831	214	238	1			
1	D	165	Total	C	N	O	S	0	0	0
			1282	831	213	236	2			
1	E	164	Total	C	N	O	S	0	0	0
			1275	826	213	235	1			
1	F	165	Total	C	N	O	S	0	0	0
			1271	825	212	232	2			
1	G	165	Total	C	N	O	S	0	0	0
			1278	827	212	238	1			
1	H	167	Total	C	N	O	S	0	0	0
			1302	843	217	241	1			
1	I	163	Total	C	N	O	S	0	0	0
			1270	824	210	235	1			
1	J	164	Total	C	N	O	S	0	0	0
			1279	829	213	236	1			

There are 210 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-19	MET	-	initiating methionine	UNP Q04120
A	-18	GLY	-	expression tag	UNP Q04120
A	-17	SER	-	expression tag	UNP Q04120
A	-16	SER	-	expression tag	UNP Q04120
A	-15	HIS	-	expression tag	UNP Q04120
A	-14	HIS	-	expression tag	UNP Q04120
A	-13	HIS	-	expression tag	UNP Q04120
A	-12	HIS	-	expression tag	UNP Q04120
A	-11	HIS	-	expression tag	UNP Q04120

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
A	-10	HIS	-	expression tag	UNP Q04120
A	-9	SER	-	expression tag	UNP Q04120
A	-8	SER	-	expression tag	UNP Q04120
A	-7	GLY	-	expression tag	UNP Q04120
A	-6	LEU	-	expression tag	UNP Q04120
A	-5	VAL	-	expression tag	UNP Q04120
A	-4	PRO	-	expression tag	UNP Q04120
A	-3	ARG	-	expression tag	UNP Q04120
A	-2	GLY	-	expression tag	UNP Q04120
A	-1	SER	-	expression tag	UNP Q04120
A	0	HIS	-	expression tag	UNP Q04120
A	171	SER	CYS	engineered mutation	UNP Q04120
B	-19	MET	-	initiating methionine	UNP Q04120
B	-18	GLY	-	expression tag	UNP Q04120
B	-17	SER	-	expression tag	UNP Q04120
B	-16	SER	-	expression tag	UNP Q04120
B	-15	HIS	-	expression tag	UNP Q04120
B	-14	HIS	-	expression tag	UNP Q04120
B	-13	HIS	-	expression tag	UNP Q04120
B	-12	HIS	-	expression tag	UNP Q04120
B	-11	HIS	-	expression tag	UNP Q04120
B	-10	HIS	-	expression tag	UNP Q04120
B	-9	SER	-	expression tag	UNP Q04120
B	-8	SER	-	expression tag	UNP Q04120
B	-7	GLY	-	expression tag	UNP Q04120
B	-6	LEU	-	expression tag	UNP Q04120
B	-5	VAL	-	expression tag	UNP Q04120
B	-4	PRO	-	expression tag	UNP Q04120
B	-3	ARG	-	expression tag	UNP Q04120
B	-2	GLY	-	expression tag	UNP Q04120
B	-1	SER	-	expression tag	UNP Q04120
B	0	HIS	-	expression tag	UNP Q04120
B	171	SER	CYS	engineered mutation	UNP Q04120
C	-19	MET	-	initiating methionine	UNP Q04120
C	-18	GLY	-	expression tag	UNP Q04120
C	-17	SER	-	expression tag	UNP Q04120
C	-16	SER	-	expression tag	UNP Q04120
C	-15	HIS	-	expression tag	UNP Q04120
C	-14	HIS	-	expression tag	UNP Q04120
C	-13	HIS	-	expression tag	UNP Q04120
C	-12	HIS	-	expression tag	UNP Q04120
C	-11	HIS	-	expression tag	UNP Q04120

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
C	-10	HIS	-	expression tag	UNP Q04120
C	-9	SER	-	expression tag	UNP Q04120
C	-8	SER	-	expression tag	UNP Q04120
C	-7	GLY	-	expression tag	UNP Q04120
C	-6	LEU	-	expression tag	UNP Q04120
C	-5	VAL	-	expression tag	UNP Q04120
C	-4	PRO	-	expression tag	UNP Q04120
C	-3	ARG	-	expression tag	UNP Q04120
C	-2	GLY	-	expression tag	UNP Q04120
C	-1	SER	-	expression tag	UNP Q04120
C	0	HIS	-	expression tag	UNP Q04120
C	171	SER	CYS	engineered mutation	UNP Q04120
D	-19	MET	-	initiating methionine	UNP Q04120
D	-18	GLY	-	expression tag	UNP Q04120
D	-17	SER	-	expression tag	UNP Q04120
D	-16	SER	-	expression tag	UNP Q04120
D	-15	HIS	-	expression tag	UNP Q04120
D	-14	HIS	-	expression tag	UNP Q04120
D	-13	HIS	-	expression tag	UNP Q04120
D	-12	HIS	-	expression tag	UNP Q04120
D	-11	HIS	-	expression tag	UNP Q04120
D	-10	HIS	-	expression tag	UNP Q04120
D	-9	SER	-	expression tag	UNP Q04120
D	-8	SER	-	expression tag	UNP Q04120
D	-7	GLY	-	expression tag	UNP Q04120
D	-6	LEU	-	expression tag	UNP Q04120
D	-5	VAL	-	expression tag	UNP Q04120
D	-4	PRO	-	expression tag	UNP Q04120
D	-3	ARG	-	expression tag	UNP Q04120
D	-2	GLY	-	expression tag	UNP Q04120
D	-1	SER	-	expression tag	UNP Q04120
D	0	HIS	-	expression tag	UNP Q04120
D	171	SER	CYS	engineered mutation	UNP Q04120
E	-19	MET	-	initiating methionine	UNP Q04120
E	-18	GLY	-	expression tag	UNP Q04120
E	-17	SER	-	expression tag	UNP Q04120
E	-16	SER	-	expression tag	UNP Q04120
E	-15	HIS	-	expression tag	UNP Q04120
E	-14	HIS	-	expression tag	UNP Q04120
E	-13	HIS	-	expression tag	UNP Q04120
E	-12	HIS	-	expression tag	UNP Q04120
E	-11	HIS	-	expression tag	UNP Q04120

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
E	-10	HIS	-	expression tag	UNP Q04120
E	-9	SER	-	expression tag	UNP Q04120
E	-8	SER	-	expression tag	UNP Q04120
E	-7	GLY	-	expression tag	UNP Q04120
E	-6	LEU	-	expression tag	UNP Q04120
E	-5	VAL	-	expression tag	UNP Q04120
E	-4	PRO	-	expression tag	UNP Q04120
E	-3	ARG	-	expression tag	UNP Q04120
E	-2	GLY	-	expression tag	UNP Q04120
E	-1	SER	-	expression tag	UNP Q04120
E	0	HIS	-	expression tag	UNP Q04120
E	171	SER	CYS	engineered mutation	UNP Q04120
F	-19	MET	-	initiating methionine	UNP Q04120
F	-18	GLY	-	expression tag	UNP Q04120
F	-17	SER	-	expression tag	UNP Q04120
F	-16	SER	-	expression tag	UNP Q04120
F	-15	HIS	-	expression tag	UNP Q04120
F	-14	HIS	-	expression tag	UNP Q04120
F	-13	HIS	-	expression tag	UNP Q04120
F	-12	HIS	-	expression tag	UNP Q04120
F	-11	HIS	-	expression tag	UNP Q04120
F	-10	HIS	-	expression tag	UNP Q04120
F	-9	SER	-	expression tag	UNP Q04120
F	-8	SER	-	expression tag	UNP Q04120
F	-7	GLY	-	expression tag	UNP Q04120
F	-6	LEU	-	expression tag	UNP Q04120
F	-5	VAL	-	expression tag	UNP Q04120
F	-4	PRO	-	expression tag	UNP Q04120
F	-3	ARG	-	expression tag	UNP Q04120
F	-2	GLY	-	expression tag	UNP Q04120
F	-1	SER	-	expression tag	UNP Q04120
F	0	HIS	-	expression tag	UNP Q04120
F	171	SER	CYS	engineered mutation	UNP Q04120
G	-19	MET	-	initiating methionine	UNP Q04120
G	-18	GLY	-	expression tag	UNP Q04120
G	-17	SER	-	expression tag	UNP Q04120
G	-16	SER	-	expression tag	UNP Q04120
G	-15	HIS	-	expression tag	UNP Q04120
G	-14	HIS	-	expression tag	UNP Q04120
G	-13	HIS	-	expression tag	UNP Q04120
G	-12	HIS	-	expression tag	UNP Q04120
G	-11	HIS	-	expression tag	UNP Q04120

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
G	-10	HIS	-	expression tag	UNP Q04120
G	-9	SER	-	expression tag	UNP Q04120
G	-8	SER	-	expression tag	UNP Q04120
G	-7	GLY	-	expression tag	UNP Q04120
G	-6	LEU	-	expression tag	UNP Q04120
G	-5	VAL	-	expression tag	UNP Q04120
G	-4	PRO	-	expression tag	UNP Q04120
G	-3	ARG	-	expression tag	UNP Q04120
G	-2	GLY	-	expression tag	UNP Q04120
G	-1	SER	-	expression tag	UNP Q04120
G	0	HIS	-	expression tag	UNP Q04120
G	171	SER	CYS	engineered mutation	UNP Q04120
H	-19	MET	-	initiating methionine	UNP Q04120
H	-18	GLY	-	expression tag	UNP Q04120
H	-17	SER	-	expression tag	UNP Q04120
H	-16	SER	-	expression tag	UNP Q04120
H	-15	HIS	-	expression tag	UNP Q04120
H	-14	HIS	-	expression tag	UNP Q04120
H	-13	HIS	-	expression tag	UNP Q04120
H	-12	HIS	-	expression tag	UNP Q04120
H	-11	HIS	-	expression tag	UNP Q04120
H	-10	HIS	-	expression tag	UNP Q04120
H	-9	SER	-	expression tag	UNP Q04120
H	-8	SER	-	expression tag	UNP Q04120
H	-7	GLY	-	expression tag	UNP Q04120
H	-6	LEU	-	expression tag	UNP Q04120
H	-5	VAL	-	expression tag	UNP Q04120
H	-4	PRO	-	expression tag	UNP Q04120
H	-3	ARG	-	expression tag	UNP Q04120
H	-2	GLY	-	expression tag	UNP Q04120
H	-1	SER	-	expression tag	UNP Q04120
H	0	HIS	-	expression tag	UNP Q04120
H	171	SER	CYS	engineered mutation	UNP Q04120
I	-19	MET	-	initiating methionine	UNP Q04120
I	-18	GLY	-	expression tag	UNP Q04120
I	-17	SER	-	expression tag	UNP Q04120
I	-16	SER	-	expression tag	UNP Q04120
I	-15	HIS	-	expression tag	UNP Q04120
I	-14	HIS	-	expression tag	UNP Q04120
I	-13	HIS	-	expression tag	UNP Q04120
I	-12	HIS	-	expression tag	UNP Q04120
I	-11	HIS	-	expression tag	UNP Q04120

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
I	-10	HIS	-	expression tag	UNP Q04120
I	-9	SER	-	expression tag	UNP Q04120
I	-8	SER	-	expression tag	UNP Q04120
I	-7	GLY	-	expression tag	UNP Q04120
I	-6	LEU	-	expression tag	UNP Q04120
I	-5	VAL	-	expression tag	UNP Q04120
I	-4	PRO	-	expression tag	UNP Q04120
I	-3	ARG	-	expression tag	UNP Q04120
I	-2	GLY	-	expression tag	UNP Q04120
I	-1	SER	-	expression tag	UNP Q04120
I	0	HIS	-	expression tag	UNP Q04120
I	171	SER	CYS	engineered mutation	UNP Q04120
J	-19	MET	-	initiating methionine	UNP Q04120
J	-18	GLY	-	expression tag	UNP Q04120
J	-17	SER	-	expression tag	UNP Q04120
J	-16	SER	-	expression tag	UNP Q04120
J	-15	HIS	-	expression tag	UNP Q04120
J	-14	HIS	-	expression tag	UNP Q04120
J	-13	HIS	-	expression tag	UNP Q04120
J	-12	HIS	-	expression tag	UNP Q04120
J	-11	HIS	-	expression tag	UNP Q04120
J	-10	HIS	-	expression tag	UNP Q04120
J	-9	SER	-	expression tag	UNP Q04120
J	-8	SER	-	expression tag	UNP Q04120
J	-7	GLY	-	expression tag	UNP Q04120
J	-6	LEU	-	expression tag	UNP Q04120
J	-5	VAL	-	expression tag	UNP Q04120
J	-4	PRO	-	expression tag	UNP Q04120
J	-3	ARG	-	expression tag	UNP Q04120
J	-2	GLY	-	expression tag	UNP Q04120
J	-1	SER	-	expression tag	UNP Q04120
J	0	HIS	-	expression tag	UNP Q04120
J	171	SER	CYS	engineered mutation	UNP Q04120

- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	5	Total O 5 5	0	0
2	B	2	Total O 2 2	0	0
2	C	6	Total O 6 6	0	0

Continued on next page...

Continued from previous page...

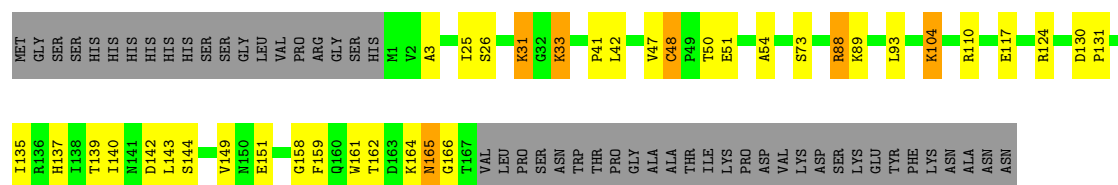
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	D	2	Total 2	O 2	0	0
2	E	1	Total 1	O 1	0	0
2	F	1	Total 1	O 1	0	0
2	G	2	Total 2	O 2	0	0
2	H	3	Total 3	O 3	0	0
2	J	2	Total 2	O 2	0	0

3 Residue-property plots

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.

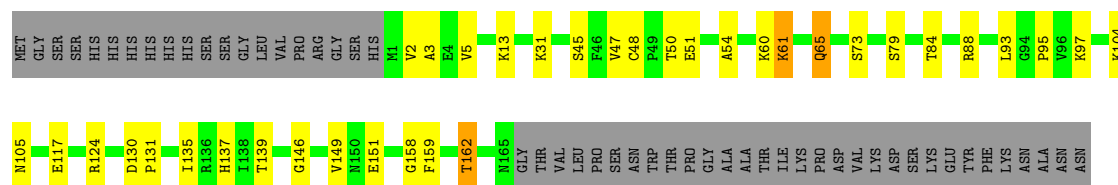
• Molecule 1: Peroxiredoxin TSA2

Chain A: 



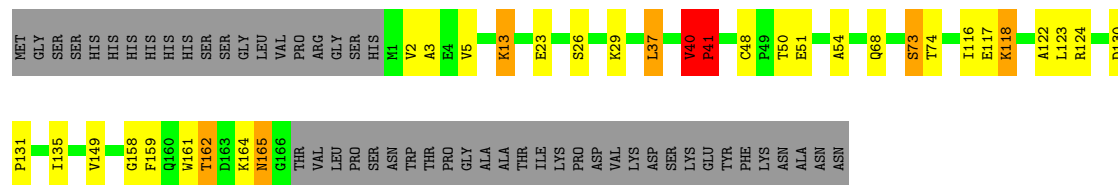
• Molecule 1: Peroxiredoxin TSA2

Chain B: 



• Molecule 1: Peroxiredoxin TSA2

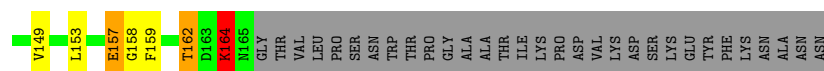
Chain C: 



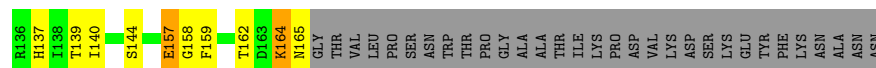
• Molecule 1: Peroxiredoxin TSA2

Chain D: 

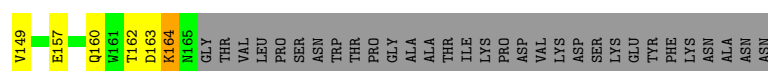




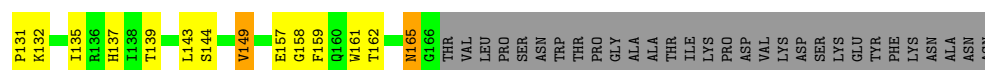
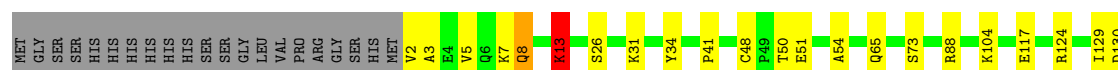
• Molecule 1: Peroxiredoxin TSA2



• Molecule 1: Peroxiredoxin TSA2



• Molecule 1: Peroxiredoxin TSA2

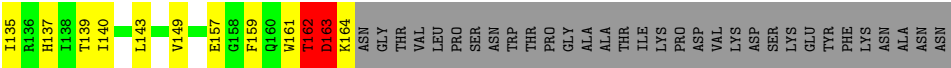


• Molecule 1: Peroxiredoxin TSA2

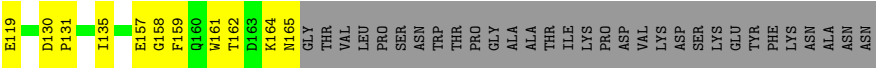


• Molecule 1: Peroxiredoxin TSA2





● Molecule 1: Peroxiredoxin TSA2



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	53.37Å 177.31Å 115.26Å 90.00° 103.28° 90.00°	Depositor
Resolution (Å)	47.40 – 2.60 47.40 – 2.60	Depositor EDS
% Data completeness (in resolution range)	97.7 (47.40-2.60) 97.8 (47.40-2.60)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.16 (at 2.61Å)	Xtriage
Refinement program	REFMAC 5.8.0158	Depositor
R, R_{free}	0.198 , 0.229 0.194 , 0.224	Depositor DCC
R_{free} test set	3018 reflections (4.61%)	wwPDB-VP
Wilson B-factor (Å ²)	76.0	Xtriage
Anisotropy	0.233	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 38.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.43$, $\langle L^2 \rangle = 0.26$	Xtriage
Estimated twinning fraction	0.210 for h,-k,-h-l	Xtriage
Reported twinning fraction	0.733 for H, K, L 0.267 for -H, -K, H+L	Depositor
Outliers	0 of 62339 reflections	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	12817	wwPDB-VP
Average B, all atoms (Å ²)	75.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.19% 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

Bond lengths and bond angles in the following residue types are not validated in this section: YCM

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# $ Z > 5$	RMSZ	# $ Z > 5$
1	A	1.15	0/1302	1.23	4/1767 (0.2%)
1	B	1.22	2/1281 (0.2%)	1.20	0/1738
1	C	1.28	3/1300 (0.2%)	1.28	7/1760 (0.4%)
1	D	1.19	1/1298 (0.1%)	1.20	7/1758 (0.4%)
1	E	1.27	4/1291 (0.3%)	1.31	13/1748 (0.7%)
1	F	1.20	5/1287 (0.4%)	1.22	11/1745 (0.6%)
1	G	1.11	0/1294	1.20	6/1754 (0.3%)
1	H	1.14	0/1318	1.22	6/1783 (0.3%)
1	I	1.20	3/1286 (0.2%)	1.46	20/1742 (1.1%)
1	J	1.18	1/1295 (0.1%)	1.26	5/1752 (0.3%)
All	All	1.19	19/12952 (0.1%)	1.26	79/17547 (0.5%)

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	C	0	3
1	D	0	1
1	E	0	2
1	F	0	2
1	I	0	3
All	All	0	11

All (19) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	41	PRO	C-O	11.38	1.36	1.23
1	E	41	PRO	CA-C	-10.71	1.40	1.52
1	E	41	PRO	C-O	9.88	1.34	1.23

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	41	PRO	CA-C	-8.78	1.42	1.52
1	I	41	PRO	CA-CB	7.16	1.64	1.53
1	I	162	THR	CA-C	6.74	1.60	1.53
1	F	41	PRO	CA-C	6.72	1.60	1.52
1	F	41	PRO	C-N	-6.00	1.25	1.33
1	D	138	ILE	C-O	-5.85	1.18	1.24
1	C	164	LYS	C-O	-5.84	1.16	1.24
1	B	97	LYS	C-O	-5.61	1.16	1.24
1	B	146	GLY	N-CA	5.55	1.50	1.44
1	F	42	LEU	N-CA	-5.48	1.39	1.46
1	E	40	VAL	CA-CB	5.47	1.61	1.54
1	J	68	GLN	CD-OE1	5.44	1.33	1.23
1	E	42	LEU	C-O	-5.35	1.17	1.23
1	I	164	LYS	N-CA	5.32	1.56	1.46
1	F	42	LEU	C-O	-5.19	1.17	1.24
1	F	3	ALA	CA-C	-5.06	1.46	1.52

All (79) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	I	40	VAL	CA-C-N	15.46	139.17	119.84
1	I	40	VAL	C-N-CA	15.46	139.17	119.84
1	I	40	VAL	N-CA-C	13.55	138.15	108.88
1	A	33	LYS	CA-CB-CG	11.07	136.24	114.10
1	E	41	PRO	CA-C-N	9.83	139.24	121.85
1	E	41	PRO	C-N-CA	9.83	139.24	121.85
1	I	42	LEU	CA-CB-CG	9.66	150.10	116.30
1	C	40	VAL	N-CA-C	9.38	129.15	108.88
1	E	40	VAL	N-CA-C	8.85	127.99	108.88
1	C	41	PRO	CA-C-N	8.73	137.31	121.85
1	C	41	PRO	C-N-CA	8.73	137.31	121.85
1	I	40	VAL	O-C-N	-7.99	112.00	121.10
1	A	31	LYS	CA-CB-CG	7.91	129.93	114.10
1	I	163	ASP	N-CA-C	7.74	127.28	110.80
1	F	31	LYS	CD-CE-NZ	-7.32	88.49	111.90
1	I	104	LYS	CG-CD-CE	7.30	128.10	111.30
1	A	104	LYS	CG-CD-CE	7.23	127.94	111.30
1	H	104	LYS	CG-CD-CE	7.11	127.65	111.30
1	J	13	LYS	CB-CG-CD	6.84	127.04	111.30
1	J	164	LYS	N-CA-C	6.79	119.21	110.65
1	J	110	ARG	NE-CZ-NH1	-6.73	114.77	121.50
1	H	162	THR	OG1-CB-CG2	6.65	122.61	109.30

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	104	LYS	CG-CD-CE	6.65	126.59	111.30
1	I	40	VAL	CB-CA-C	-6.34	99.81	111.36
1	I	42	LEU	N-CA-C	6.29	119.75	109.24
1	F	42	LEU	N-CA-CB	-6.27	99.89	110.49
1	F	164	LYS	O-C-N	-6.24	114.29	122.59
1	I	13	LYS	CD-CE-NZ	6.16	131.62	111.90
1	G	104	LYS	CG-CD-CE	6.15	125.44	111.30
1	F	110	ARG	NE-CZ-NH1	-6.07	115.43	121.50
1	A	144	SER	N-CA-C	6.00	119.70	112.38
1	I	41	PRO	N-CD-CG	6.00	112.20	103.20
1	H	144	SER	N-CA-C	5.96	119.64	112.38
1	I	164	LYS	N-CA-C	5.94	127.63	111.00
1	G	13	LYS	CB-CG-CD	5.94	124.96	111.30
1	E	104	LYS	N-CA-C	5.90	118.46	111.33
1	D	4	GLU	CA-CB-CG	5.88	125.86	114.10
1	G	88	ARG	CG-CD-NE	5.84	124.85	112.00
1	I	132	LYS	CG-CD-CE	5.83	124.70	111.30
1	D	33	LYS	CD-CE-NZ	5.78	130.40	111.90
1	F	42	LEU	N-CA-C	5.78	123.11	110.80
1	E	40	VAL	CB-CA-C	-5.74	100.92	111.36
1	E	40	VAL	CA-C-O	-5.69	112.33	119.95
1	D	13	LYS	CB-CG-CD	5.64	124.27	111.30
1	I	162	THR	CA-CB-OG1	-5.63	101.15	109.60
1	I	162	THR	N-CA-C	-5.62	98.08	108.24
1	D	164	LYS	N-CA-C	5.61	119.11	112.72
1	E	41	PRO	N-CA-CB	5.60	108.28	103.36
1	E	144	SER	N-CA-C	5.60	119.21	112.38
1	H	162	THR	CA-CB-CG2	5.52	119.89	110.50
1	F	45	SER	N-CA-C	-5.49	103.28	110.53
1	C	40	VAL	CB-CA-C	-5.48	101.38	111.36
1	C	37	LEU	CB-CG-CD2	5.44	127.02	110.70
1	E	42	LEU	N-CA-C	5.40	117.46	108.99
1	H	149	VAL	CB-CA-C	-5.38	103.83	112.16
1	F	164	LYS	CA-C-N	5.35	131.33	121.70
1	F	164	LYS	C-N-CA	5.35	131.33	121.70
1	I	163	ASP	CA-C-N	5.30	131.23	121.70
1	I	163	ASP	C-N-CA	5.30	131.23	121.70
1	H	45	SER	N-CA-C	-5.28	103.56	110.53
1	G	149	VAL	CB-CA-C	-5.28	103.98	112.16
1	F	144	SER	N-CA-C	5.28	118.82	112.38
1	F	42	LEU	CB-CG-CD2	5.27	126.51	110.70
1	D	157	GLU	CA-CB-CG	5.26	124.61	114.10

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	F	162	THR	N-CA-C	5.24	117.74	111.71
1	I	40	VAL	C-N-CD	-5.22	103.60	125.00
1	C	123	LEU	N-CA-C	-5.17	104.05	110.41
1	G	13	LYS	CD-CE-NZ	-5.16	95.38	111.90
1	E	10	PRO	CA-C-N	-5.13	114.95	120.03
1	E	10	PRO	C-N-CA	-5.13	114.95	120.03
1	J	42	LEU	N-CA-C	5.11	117.75	109.06
1	G	104	LYS	N-CA-C	5.09	116.83	111.28
1	I	162	THR	N-CA-CB	5.08	117.95	110.49
1	I	164	LYS	CA-C-O	5.07	129.42	120.80
1	C	40	VAL	CA-C-O	-5.05	113.18	119.95
1	D	164	LYS	CA-C-O	-5.04	113.78	119.78
1	J	61	LYS	CG-CD-CE	5.01	122.82	111.30
1	E	40	VAL	CA-C-N	-5.01	114.64	119.90
1	E	40	VAL	C-N-CA	-5.01	114.64	119.90

There are no chirality outliers.

All (11) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	C	40	VAL	Peptide
1	C	41	PRO	Mainchain,Peptide
1	D	164	LYS	Mainchain
1	E	41	PRO	Mainchain,Peptide
1	F	164	LYS	Peptide
1	F	41	PRO	Peptide
1	I	163	ASP	Peptide
1	I	40	VAL	Mainchain,Peptide

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	1286	0	1281	27	0
1	B	1266	0	1262	17	0
1	C	1284	0	1292	21	0
1	D	1282	0	1296	16	0

Continued on next page...

Continued from previous page...

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	E	1275	0	1286	21	4
1	F	1271	0	1279	13	0
1	G	1278	0	1280	22	0
1	H	1302	0	1325	19	0
1	I	1270	0	1282	12	4
1	J	1279	0	1295	17	0
2	A	5	0	0	8	0
2	B	2	0	0	0	0
2	C	6	0	0	0	0
2	D	2	0	0	0	0
2	E	1	0	0	0	0
2	F	1	0	0	1	0
2	G	2	0	0	6	0
2	H	3	0	0	1	0
2	J	2	0	0	0	0
All	All	12817	0	12878	175	4

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 7.

All (175) 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:A:25:ILE:HG22	2:A:201:HOH:O	1.46	1.12
1:A:142:ASP:HB3	2:A:203:HOH:O	1.57	1.04
1:E:65:GLN:HG3	1:E:157:GLU:OE1	1.61	0.98
1:H:89:LYS:NZ	1:H:90:ASP:OD2	2.02	0.93
1:A:165:ASN:N	1:A:165:ASN:HD22	1.70	0.89
1:C:41:PRO:HD2	1:C:122:ALA:HB3	1.56	0.87
1:E:41:PRO:HD2	1:E:122:ALA:HB3	1.57	0.86
1:J:162:THR:O	1:J:165:ASN:ND2	2.09	0.84
1:I:41:PRO:HG3	1:I:143:LEU:HD22	1.62	0.82
1:B:84:THR:HG23	1:B:95:PRO:HA	1.59	0.81
1:B:45:SER:OG	1:B:47:VAL:HG12	1.81	0.80
1:F:45:SER:OG	1:F:47:VAL:HG22	1.81	0.78
1:A:164:LYS:C	1:A:165:ASN:HD22	1.90	0.77
1:E:41:PRO:CD	1:E:122:ALA:HB3	2.13	0.77
1:A:25:ILE:C	2:A:201:HOH:O	2.29	0.75
1:C:116:ILE:HG22	1:C:118:LYS:HD2	1.68	0.75
1:D:41:PRO:HG3	1:D:143:LEU:HD22	1.68	0.75
1:C:41:PRO:CD	1:C:122:ALA:HB3	2.16	0.75

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:34:TYR:CE2	2:G:201:HOH:O	2.40	0.74
1:I:159:PHE:O	1:I:163:ASP:HA	1.87	0.73
1:F:41:PRO:HG3	1:F:143:LEU:HD22	1.71	0.72
1:A:26:SER:N	2:A:201:HOH:O	2.23	0.71
1:A:26:SER:CA	2:A:201:HOH:O	2.38	0.71
1:G:144:SER:HB2	1:H:162:THR:HG21	1.73	0.71
1:F:163:ASP:CB	2:F:201:HOH:O	2.39	0.70
1:G:34:TYR:CD2	2:G:201:HOH:O	2.46	0.68
1:A:137:HIS:HE2	1:A:139:THR:HG1	1.38	0.68
1:A:41:PRO:HG3	1:A:143:LEU:HD22	1.75	0.68
1:E:65:GLN:HG3	1:E:157:GLU:CD	2.19	0.68
1:C:41:PRO:HG3	1:C:73:SER:HA	1.76	0.68
1:E:76:SER:HB3	1:E:104:LYS:HD2	1.77	0.66
1:A:26:SER:C	2:A:201:HOH:O	2.39	0.65
1:A:161:TRP:O	1:A:165:ASN:ND2	2.29	0.65
1:E:41:PRO:HG2	1:E:122:ALA:H	1.61	0.65
1:A:165:ASN:N	1:A:165:ASN:ND2	2.41	0.65
1:E:164:LYS:HG3	1:E:165:ASN:ND2	2.12	0.64
1:G:161:TRP:O	1:G:165:ASN:ND2	2.31	0.64
1:C:161:TRP:O	1:C:165:ASN:ND2	2.30	0.64
1:J:8:GLN:HE21	1:J:8:GLN:HA	1.64	0.62
1:E:65:GLN:CG	1:E:157:GLU:OE1	2.43	0.61
1:H:60:LYS:NZ	1:H:64:ASP:OD1	2.35	0.60
1:B:158:GLY:O	1:B:162:THR:HG22	2.02	0.60
1:E:158:GLY:O	1:E:162:THR:HG22	2.02	0.60
1:C:158:GLY:O	1:C:162:THR:HG22	2.02	0.60
1:G:129:ILE:C	2:G:201:HOH:O	2.44	0.60
1:E:140:ILE:HG22	1:F:5:VAL:HG21	1.83	0.60
1:D:48:YCM:NZ2	1:D:49:PRO:HD2	2.17	0.59
1:G:41:PRO:HG3	1:G:143:LEU:HD22	1.83	0.59
1:G:158:GLY:O	1:G:162:THR:HG22	2.02	0.59
1:A:158:GLY:O	1:A:162:THR:HG22	2.02	0.58
1:G:130:ASP:CA	2:G:201:HOH:O	2.51	0.58
1:D:158:GLY:O	1:D:162:THR:HG22	2.02	0.58
1:J:8:GLN:HA	1:J:8:GLN:NE2	2.20	0.57
1:J:158:GLY:O	1:J:162:THR:HG22	2.03	0.57
1:A:164:LYS:C	1:A:165:ASN:ND2	2.62	0.57
1:G:13:LYS:HG2	1:G:26:SER:HB3	1.88	0.56
1:F:40:VAL:HG22	1:F:41:PRO:O	2.06	0.56
1:G:130:ASP:N	2:G:201:HOH:O	2.37	0.56
1:F:31:LYS:HD2	1:F:131:PRO:O	2.09	0.54

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:158:GLY:O	1:H:162:THR:HG23	2.07	0.53
1:A:140:ILE:HG22	1:B:5:VAL:HG21	1.93	0.51
1:E:41:PRO:HG3	1:E:73:SER:HA	1.92	0.51
1:A:142:ASP:C	2:A:203:HOH:O	2.53	0.51
1:A:54:ALA:HB1	1:A:149:VAL:HG21	1.92	0.51
1:A:110:ARG:NH1	1:A:117:GLU:OE1	2.38	0.50
1:C:13:LYS:HG2	1:C:26:SER:HB3	1.92	0.50
1:E:159:PHE:HA	1:E:162:THR:CG2	2.42	0.50
1:B:54:ALA:HB1	1:B:149:VAL:HG21	1.94	0.50
1:J:40:VAL:CG2	1:J:41:PRO:HD2	2.42	0.50
1:F:2:VAL:HG12	1:F:3:ALA:H	1.77	0.49
1:H:54:ALA:HB1	1:H:149:VAL:HG21	1.92	0.49
1:G:54:ALA:HB1	1:G:149:VAL:HG21	1.94	0.49
1:A:159:PHE:HA	1:A:162:THR:CG2	2.43	0.49
1:J:13:LYS:HG2	1:J:26:SER:HB3	1.94	0.49
1:G:159:PHE:HA	1:G:162:THR:CG2	2.43	0.49
1:J:40:VAL:CG2	1:J:41:PRO:CD	2.90	0.49
1:C:5:VAL:HG21	1:D:140:ILE:HG22	1.95	0.49
1:J:159:PHE:HA	1:J:162:THR:CG2	2.42	0.49
1:E:65:GLN:HE21	1:E:157:GLU:CD	2.20	0.48
1:H:161:TRP:HZ3	1:H:167:THR:O	1.96	0.48
1:I:54:ALA:HB1	1:I:149:VAL:HG21	1.94	0.48
1:J:116:ILE:HG21	1:J:119:GLU:HG3	1.96	0.48
1:C:2:VAL:HG12	1:C:3:ALA:H	1.77	0.48
1:D:54:ALA:HB1	1:D:149:VAL:HG21	1.94	0.48
1:D:159:PHE:HA	1:D:162:THR:CG2	2.43	0.48
1:A:48:YCM:OZ1	1:A:88:ARG:HD3	2.13	0.47
2:A:203:HOH:O	1:B:159:PHE:HE1	1.95	0.47
1:C:159:PHE:HA	1:C:162:THR:CG2	2.43	0.47
1:B:159:PHE:HA	1:B:162:THR:CG2	2.45	0.47
1:G:130:ASP:HA	2:G:201:HOH:O	2.14	0.47
1:E:3:ALA:HB1	1:E:135:ILE:CD1	2.45	0.47
1:G:51:GLU:OE1	1:G:124:ARG:NH1	2.48	0.47
1:C:2:VAL:HG12	1:C:3:ALA:N	2.30	0.46
1:C:51:GLU:OE1	1:C:124:ARG:NH1	2.48	0.46
1:E:137:HIS:NE2	1:E:139:THR:OG1	2.47	0.46
1:C:54:ALA:HB1	1:C:149:VAL:HG21	1.97	0.46
1:E:159:PHE:HA	1:E:162:THR:HG22	1.98	0.46
1:F:130:ASP:HB2	1:F:131:PRO:HD2	1.98	0.46
1:J:159:PHE:HA	1:J:162:THR:HG22	1.98	0.46
1:B:61:LYS:HE2	1:B:65:GLN:NE2	2.31	0.46

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:41:PRO:HG2	1:C:122:ALA:H	1.80	0.46
1:J:3:ALA:HB1	1:J:135:ILE:CD1	2.46	0.46
1:G:162:THR:HG21	1:H:144:SER:HB2	1.98	0.45
1:B:51:GLU:OE1	1:B:124:ARG:NH1	2.49	0.45
1:J:40:VAL:HG23	1:J:41:PRO:CD	2.46	0.45
1:A:159:PHE:HA	1:A:162:THR:HG22	1.99	0.45
1:C:130:ASP:HB2	1:C:131:PRO:HD2	1.99	0.45
1:H:3:ALA:HB1	1:H:135:ILE:CD1	2.47	0.45
1:D:130:ASP:HB2	1:D:131:PRO:HD2	1.99	0.45
1:J:8:GLN:HE21	1:J:8:GLN:CA	2.25	0.45
1:H:41:PRO:HG3	1:H:143:LEU:HD22	1.99	0.45
1:F:3:ALA:HB1	1:F:135:ILE:CD1	2.46	0.45
1:F:54:ALA:HB1	1:F:149:VAL:HG21	1.98	0.45
1:A:130:ASP:HB2	1:A:131:PRO:HD2	2.00	0.44
1:H:51:GLU:OE1	1:H:124:ARG:NH1	2.50	0.44
1:I:51:GLU:OE1	1:I:124:ARG:NH1	2.50	0.44
1:H:89:LYS:NZ	1:H:90:ASP:CG	2.75	0.44
1:H:130:ASP:HB2	1:H:131:PRO:HD2	1.99	0.44
1:A:51:GLU:OE1	1:A:124:ARG:NH1	2.50	0.44
1:F:51:GLU:OE1	1:F:124:ARG:NH1	2.49	0.44
1:H:137:HIS:NE2	1:H:139:THR:OG1	2.47	0.44
1:I:65:GLN:NE2	1:I:157:GLU:OE2	2.50	0.44
1:I:130:ASP:HB2	1:I:131:PRO:HD2	2.00	0.44
1:F:2:VAL:HG12	1:F:3:ALA:N	2.32	0.44
1:J:40:VAL:HG22	1:J:41:PRO:HD2	1.98	0.44
1:G:159:PHE:HA	1:G:162:THR:HG22	1.99	0.44
1:I:140:ILE:HG22	1:J:5:VAL:HG21	2.00	0.44
1:G:5:VAL:HG21	1:H:140:ILE:HG22	2.00	0.44
1:C:23:GLU:OE2	1:C:29:LYS:NZ	2.51	0.43
1:D:51:GLU:OE1	1:D:124:ARG:NH1	2.51	0.43
1:D:104:LYS:HB2	1:D:104:LYS:HE2	1.85	0.43
1:D:2:VAL:HG12	1:D:3:ALA:H	1.83	0.43
1:I:159:PHE:O	1:I:162:THR:O	2.36	0.43
1:A:3:ALA:HB1	1:A:135:ILE:CD1	2.47	0.43
1:G:130:ASP:HB2	1:G:131:PRO:HD2	2.01	0.43
1:H:110:ARG:NH2	2:H:202:HOH:O	2.39	0.43
1:J:130:ASP:HB2	1:J:131:PRO:HD2	2.00	0.43
1:C:159:PHE:HA	1:C:162:THR:HG22	1.99	0.43
1:A:42:LEU:HD13	1:J:78:TYR:CD2	2.54	0.43
1:I:47:VAL:HG12	1:I:48:YCM:N	2.34	0.43
1:C:3:ALA:HB1	1:C:135:ILE:CD1	2.49	0.43

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:116:ILE:HG22	1:C:118:LYS:CD	2.45	0.43
1:I:42:LEU:HD13	1:I:121:ILE:HG21	2.00	0.43
1:C:41:PRO:HG3	1:C:73:SER:CA	2.46	0.43
1:D:159:PHE:HA	1:D:162:THR:HG22	1.99	0.43
1:I:3:ALA:HB1	1:I:135:ILE:CD1	2.49	0.43
1:D:3:ALA:HB1	1:D:135:ILE:CD1	2.48	0.42
1:G:3:ALA:HB1	1:G:135:ILE:CD1	2.48	0.42
1:E:65:GLN:NE2	1:E:157:GLU:CD	2.77	0.42
1:G:2:VAL:HG12	1:G:3:ALA:N	2.35	0.42
1:G:137:HIS:NE2	1:G:139:THR:OG1	2.46	0.42
1:B:137:HIS:NE2	1:B:139:THR:OG1	2.46	0.42
1:B:130:ASP:HB2	1:B:131:PRO:HD2	2.01	0.42
1:G:8:GLN:NE2	1:G:132:LYS:O	2.52	0.42
1:I:137:HIS:NE2	1:I:139:THR:OG1	2.47	0.42
1:E:130:ASP:HB2	1:E:131:PRO:HD2	2.01	0.42
1:H:2:VAL:HG12	1:H:3:ALA:N	2.34	0.42
1:H:54:ALA:HB1	1:H:149:VAL:CG2	2.50	0.42
1:B:2:VAL:HG12	1:B:3:ALA:H	1.85	0.41
1:B:159:PHE:HA	1:B:162:THR:HG22	2.01	0.41
1:A:48:YCM:OZ1	1:A:88:ARG:CD	2.69	0.41
1:H:2:VAL:HG12	1:H:3:ALA:H	1.85	0.41
1:F:65:GLN:O	1:F:160:GLN:NE2	2.50	0.41
1:E:51:GLU:OE1	1:E:124:ARG:NH1	2.52	0.41
1:B:3:ALA:HB1	1:B:135:ILE:CD1	2.50	0.41
1:D:2:VAL:HG12	1:D:3:ALA:N	2.36	0.41
1:D:42:LEU:HD13	1:E:78:TYR:CD2	2.56	0.41
1:D:153:LEU:O	1:D:157:GLU:HG3	2.21	0.41
1:B:2:VAL:HG12	1:B:3:ALA:N	2.36	0.41
1:D:105:ASN:OD1	1:D:105:ASN:C	2.64	0.41
1:E:41:PRO:HG3	1:E:74:THR:H	1.86	0.41
1:H:89:LYS:HE2	1:H:89:LYS:HB2	1.95	0.40
1:C:41:PRO:HG3	1:C:74:THR:H	1.85	0.40
1:A:151:GLU:HG3	1:B:151:GLU:HG3	2.03	0.40
1:B:105:ASN:OD1	1:B:105:ASN:C	2.64	0.40

All (4) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:48:YCM:OZ1	1:I:161:TRP:CG[1_556]	1.93	0.27

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:48:YCM:OZ1	1:I:161:TRP:CD2[1_556]	1.95	0.25
1:E:48:YCM:NZ2	1:I:161:TRP:CZ2[1_556]	2.07	0.13
1:E:48:YCM:NZ2	1:I:161:TRP:CE2[1_556]	2.18	0.02

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	164/216 (76%)	151 (92%)	12 (7%)	1 (1%)	21	42
1	B	162/216 (75%)	152 (94%)	10 (6%)	0	100	100
1	C	163/216 (76%)	151 (93%)	11 (7%)	1 (1%)	21	42
1	D	162/216 (75%)	151 (93%)	11 (7%)	0	100	100
1	E	161/216 (74%)	151 (94%)	9 (6%)	1 (1%)	21	42
1	F	162/216 (75%)	150 (93%)	11 (7%)	1 (1%)	21	42
1	G	162/216 (75%)	152 (94%)	10 (6%)	0	100	100
1	H	164/216 (76%)	152 (93%)	11 (7%)	1 (1%)	21	42
1	I	160/216 (74%)	148 (92%)	11 (7%)	1 (1%)	21	42
1	J	161/216 (74%)	151 (94%)	10 (6%)	0	100	100
All	All	1621/2160 (75%)	1509 (93%)	106 (6%)	6 (0%)	30	51

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	F	42	LEU
1	I	163	ASP
1	C	40	VAL
1	E	40	VAL
1	H	49	PRO
1	A	166	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	134/181 (74%)	124 (92%)	10 (8%)	12	28
1	B	131/181 (72%)	118 (90%)	13 (10%)	7	16
1	C	134/181 (74%)	125 (93%)	9 (7%)	15	33
1	D	135/181 (75%)	126 (93%)	9 (7%)	15	33
1	E	134/181 (74%)	125 (93%)	9 (7%)	15	33
1	F	132/181 (73%)	127 (96%)	5 (4%)	29	56
1	G	134/181 (74%)	124 (92%)	10 (8%)	12	28
1	H	138/181 (76%)	125 (91%)	13 (9%)	8	18
1	I	134/181 (74%)	122 (91%)	12 (9%)	9	20
1	J	135/181 (75%)	123 (91%)	12 (9%)	9	20
All	All	1341/1810 (74%)	1239 (92%)	102 (8%)	12	27

All (102) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	31	LYS
1	A	33	LYS
1	A	47	VAL
1	A	50	THR
1	A	73	SER
1	A	88	ARG
1	A	89	LYS
1	A	93	LEU
1	A	104	LYS
1	A	165	ASN
1	B	13	LYS
1	B	31	LYS
1	B	50	THR
1	B	60	LYS
1	B	61	LYS
1	B	65	GLN

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	B	73	SER
1	B	79	SER
1	B	88	ARG
1	B	93	LEU
1	B	104	LYS
1	B	117	GLU
1	B	162	THR
1	C	13	LYS
1	C	37	LEU
1	C	50	THR
1	C	68	GLN
1	C	73	SER
1	C	117	GLU
1	C	118	LYS
1	C	162	THR
1	C	165	ASN
1	D	1	MET
1	D	13	LYS
1	D	50	THR
1	D	73	SER
1	D	88	ARG
1	D	93	LEU
1	D	104	LYS
1	D	162	THR
1	D	164	LYS
1	E	47	VAL
1	E	50	THR
1	E	65	GLN
1	E	73	SER
1	E	88	ARG
1	E	89	LYS
1	E	93	LEU
1	E	157	GLU
1	E	164	LYS
1	F	50	THR
1	F	60	LYS
1	F	65	GLN
1	F	73	SER
1	F	157	GLU
1	G	7	LYS
1	G	8	GLN
1	G	13	LYS

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	G	31	LYS
1	G	50	THR
1	G	65	GLN
1	G	73	SER
1	G	117	GLU
1	G	157	GLU
1	G	165	ASN
1	H	7	LYS
1	H	21	ILE
1	H	31	LYS
1	H	50	THR
1	H	65	GLN
1	H	73	SER
1	H	89	LYS
1	H	97	LYS
1	H	104	LYS
1	H	117	GLU
1	H	157	GLU
1	H	162	THR
1	H	164	LYS
1	I	13	LYS
1	I	33	LYS
1	I	42	LEU
1	I	50	THR
1	I	60	LYS
1	I	65	GLN
1	I	73	SER
1	I	89	LYS
1	I	104	LYS
1	I	117	GLU
1	I	132	LYS
1	I	162	THR
1	J	13	LYS
1	J	47	VAL
1	J	50	THR
1	J	60	LYS
1	J	61	LYS
1	J	73	SER
1	J	89	LYS
1	J	93	LEU
1	J	97	LYS
1	J	118	LYS

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	J	157	GLU
1	J	161	TRP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (17) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	8	GLN
1	A	160	GLN
1	A	165	ASN
1	B	8	GLN
1	B	65	GLN
1	C	68	GLN
1	C	165	ASN
1	D	8	GLN
1	E	8	GLN
1	E	165	ASN
1	F	8	GLN
1	G	6	GLN
1	G	8	GLN
1	G	165	ASN
1	H	8	GLN
1	I	65	GLN
1	J	8	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

10 non-standard protein/DNA/RNA residues are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. 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 $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
1	YCM	F	48	1	7,9,10	0.97	0	5,10,12	1.48	1 (20%)
1	YCM	G	48	1	7,9,10	1.26	1 (14%)	5,10,12	1.70	2 (40%)
1	YCM	I	48	1	7,9,10	1.49	1 (14%)	5,10,12	2.64	2 (40%)
1	YCM	E	48	1	7,9,10	1.41	1 (14%)	5,10,12	3.34	2 (40%)
1	YCM	C	48	1	7,9,10	0.59	0	5,10,12	1.36	1 (20%)
1	YCM	J	48	1	7,9,10	0.96	0	5,10,12	1.23	0
1	YCM	B	48	1	7,9,10	1.52	0	5,10,12	2.44	3 (60%)
1	YCM	H	48	1	7,9,10	1.14	1 (14%)	5,10,12	2.09	1 (20%)
1	YCM	D	48	1	7,9,10	0.70	0	5,10,12	1.31	0
1	YCM	A	48	1	7,9,10	2.28	2 (28%)	5,10,12	3.09	1 (20%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
1	YCM	F	48	1	-	3/6/8/10	-
1	YCM	G	48	1	-	3/6/8/10	-
1	YCM	I	48	1	-	3/6/8/10	-
1	YCM	E	48	1	-	5/6/8/10	-
1	YCM	C	48	1	-	3/6/8/10	-
1	YCM	J	48	1	-	3/6/8/10	-
1	YCM	B	48	1	-	2/6/8/10	-
1	YCM	H	48	1	-	2/6/8/10	-
1	YCM	D	48	1	-	5/6/8/10	-
1	YCM	A	48	1	-	3/6/8/10	-

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	48	YCM	CB-SG	-4.56	1.63	1.81
1	I	48	YCM	CB-CA	-3.27	1.45	1.53
1	E	48	YCM	CD-SG	-3.23	1.73	1.81
1	A	48	YCM	CB-CA	-3.09	1.45	1.53
1	G	48	YCM	CB-SG	-2.57	1.71	1.81
1	H	48	YCM	CB-SG	-2.51	1.71	1.81

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	48	YCM	CE-CD-SG	-6.76	92.46	113.81
1	E	48	YCM	CE-CD-SG	-6.47	93.37	113.81
1	I	48	YCM	CA-CB-SG	-4.99	95.11	113.51
1	B	48	YCM	CB-CA-C	-3.75	100.62	110.80
1	H	48	YCM	CE-CD-SG	-3.14	103.90	113.81
1	E	48	YCM	CA-CB-SG	3.10	124.93	113.51
1	B	48	YCM	CE-CD-SG	-3.00	104.33	113.81
1	I	48	YCM	CB-CA-C	2.71	118.16	110.80
1	B	48	YCM	OZ1-CE-NZ2	2.37	128.87	122.53
1	C	48	YCM	CE-CD-SG	2.32	121.14	113.81
1	G	48	YCM	CA-CB-SG	-2.30	105.05	113.51
1	F	48	YCM	CB-CA-C	-2.11	105.08	110.80
1	G	48	YCM	CE-CD-SG	-2.06	107.31	113.81

There are no chirality outliers.

All (32) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	48	YCM	CE-CD-SG-CB
1	A	48	YCM	SG-CD-CE-OZ1
1	A	48	YCM	SG-CD-CE-NZ2
1	B	48	YCM	CE-CD-SG-CB
1	B	48	YCM	SG-CD-CE-NZ2
1	C	48	YCM	N-CA-CB-SG
1	C	48	YCM	C-CA-CB-SG
1	C	48	YCM	CE-CD-SG-CB
1	D	48	YCM	N-CA-CB-SG
1	D	48	YCM	C-CA-CB-SG
1	D	48	YCM	CA-CB-SG-CD
1	D	48	YCM	SG-CD-CE-OZ1
1	D	48	YCM	SG-CD-CE-NZ2
1	E	48	YCM	N-CA-CB-SG
1	E	48	YCM	C-CA-CB-SG
1	E	48	YCM	CA-CB-SG-CD
1	F	48	YCM	N-CA-CB-SG
1	F	48	YCM	C-CA-CB-SG
1	F	48	YCM	CA-CB-SG-CD
1	G	48	YCM	SG-CD-CE-NZ2
1	H	48	YCM	CE-CD-SG-CB
1	I	48	YCM	SG-CD-CE-OZ1
1	I	48	YCM	SG-CD-CE-NZ2

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms
1	J	48	YCM	N-CA-CB-SG
1	J	48	YCM	C-CA-CB-SG
1	J	48	YCM	CE-CD-SG-CB
1	E	48	YCM	CE-CD-SG-CB
1	G	48	YCM	CE-CD-SG-CB
1	G	48	YCM	SG-CD-CE-OZ1
1	I	48	YCM	CE-CD-SG-CB
1	H	48	YCM	SG-CD-CE-NZ2
1	E	48	YCM	SG-CD-CE-NZ2

There are no ring outliers.

4 monomers are involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	I	48	YCM	1	0
1	E	48	YCM	0	4
1	D	48	YCM	1	0
1	A	48	YCM	2	0

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	166/216 (76%)	-1.14	0 100 100	59, 75, 97, 114	0
1	B	164/216 (75%)	-1.12	0 100 100	59, 73, 87, 109	0
1	C	165/216 (76%)	-1.10	0 100 100	58, 73, 90, 100	0
1	D	164/216 (75%)	-1.18	0 100 100	57, 69, 82, 94	0
1	E	163/216 (75%)	-1.15	0 100 100	57, 70, 85, 99	0
1	F	164/216 (75%)	-1.15	0 100 100	61, 76, 95, 109	0
1	G	164/216 (75%)	-1.13	0 100 100	63, 75, 85, 92	0
1	H	166/216 (76%)	-1.09	0 100 100	62, 74, 90, 102	0
1	I	162/216 (75%)	-1.07	0 100 100	64, 79, 96, 103	0
1	J	163/216 (75%)	-1.13	0 100 100	59, 77, 97, 118	0
All	All	1641/2160 (75%)	-1.12	0 100 100	57, 74, 92, 118	0

There are no RSRZ outliers to report.

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled ‘Q< 0.9’ lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
1	YCM	B	48	10/11	0.92	0.07	91,99,103,107	0
1	YCM	J	48	10/11	0.93	0.07	89,98,103,105	0
1	YCM	C	48	10/11	0.94	0.06	72,97,101,105	0
1	YCM	H	48	10/11	0.95	0.07	76,90,95,99	0
1	YCM	I	48	10/11	0.95	0.06	78,103,110,111	0
1	YCM	F	48	10/11	0.95	0.06	87,94,109,110	0

Continued on next page...

Continued from previous page...

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
1	YCM	D	48	10/11	0.96	0.06	79,86,90,97	0
1	YCM	G	48	10/11	0.97	0.05	81,88,91,92	0
1	YCM	A	48	10/11	0.97	0.05	79,87,100,104	0
1	YCM	E	48	10/11	0.99	0.04	79,88,112,120	0

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.