



Full wwPDB X-ray Structure Validation Report ⓘ

May 29, 2020 – 03:37 am BST

PDB ID : 3OME
Title : Crystal structure of a probable ENOYL-COA Hydratase from Mycobacterium Smegmatis
Authors : Seattle Structural Genomics Center for Infectious Disease (SSGCID)
Deposited on : 2010-08-26
Resolution : 2.05 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	:	2.11
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.11

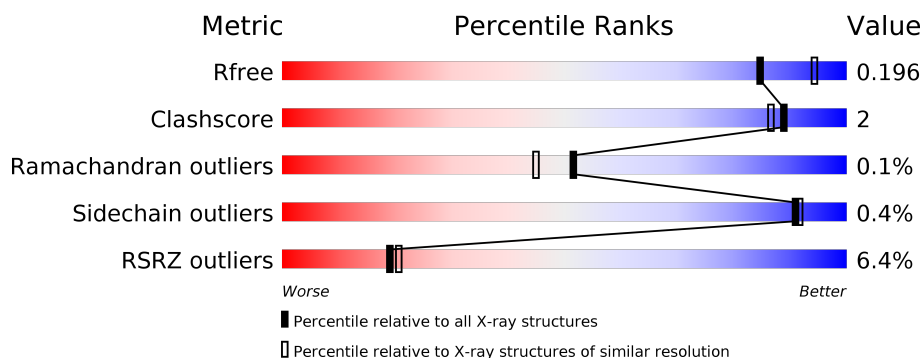
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.05 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	1692 (2.04-2.04)
Clashscore	141614	1773 (2.04-2.04)
Ramachandran outliers	138981	1752 (2.04-2.04)
Sidechain outliers	138945	1752 (2.04-2.04)
RSRZ outliers	127900	1672 (2.04-2.04)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. 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	282	<div> <div>4%</div> <div> <div></div> <div>82%</div> <div>5%</div> <div>14%</div> </div> </div>
1	B	282	<div> <div>5%</div> <div> <div></div> <div>81%</div> <div>5%</div> <div>14%</div> </div> </div>
1	C	282	<div> <div>6%</div> <div> <div></div> <div>82%</div> <div>5%</div> <div>12%</div> </div> </div>
1	D	282	<div> <div>6%</div> <div> <div></div> <div>81%</div> <div>5%</div> <div>13%</div> </div> </div>
1	E	282	<div> <div>6%</div> <div> <div></div> <div>80%</div> <div>6%</div> <div>13%</div> </div> </div>
1	F	282	<div> <div>6%</div> <div> <div></div> <div>82%</div> <div>6%</div> <div>12%</div> </div> </div>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 11909 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 Enoyl-CoA hydratase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	243	Total	C	N	O	S	0	1	0
			1867	1175	337	348	7			
1	B	242	Total	C	N	O	S	0	1	0
			1862	1173	334	348	7			
1	C	247	Total	C	N	O	S	0	1	0
			1894	1194	339	354	7			
1	D	244	Total	C	N	O	S	0	1	0
			1851	1167	331	347	6			
1	E	244	Total	C	N	O	S	0	1	0
			1883	1187	338	351	7			
1	F	247	Total	C	N	O	S	0	2	0
			1880	1190	335	349	6			

There are 132 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-21	MET	-	EXPRESSION TAG	UNP A0R1R4
A	-20	ALA	-	EXPRESSION TAG	UNP A0R1R4
A	-19	HIS	-	EXPRESSION TAG	UNP A0R1R4
A	-18	HIS	-	EXPRESSION TAG	UNP A0R1R4
A	-17	HIS	-	EXPRESSION TAG	UNP A0R1R4
A	-16	HIS	-	EXPRESSION TAG	UNP A0R1R4
A	-15	HIS	-	EXPRESSION TAG	UNP A0R1R4
A	-14	HIS	-	EXPRESSION TAG	UNP A0R1R4
A	-13	MET	-	EXPRESSION TAG	UNP A0R1R4
A	-12	GLY	-	EXPRESSION TAG	UNP A0R1R4
A	-11	THR	-	EXPRESSION TAG	UNP A0R1R4
A	-10	LEU	-	EXPRESSION TAG	UNP A0R1R4
A	-9	GLU	-	EXPRESSION TAG	UNP A0R1R4
A	-8	ALA	-	EXPRESSION TAG	UNP A0R1R4
A	-7	GLN	-	EXPRESSION TAG	UNP A0R1R4
A	-6	THR	-	EXPRESSION TAG	UNP A0R1R4
A	-5	GLN	-	EXPRESSION TAG	UNP A0R1R4

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Chain	Residue	Modelled	Actual	Comment	Reference
A	-4	GLY	-	EXPRESSION TAG	UNP A0R1R4
A	-3	PRO	-	EXPRESSION TAG	UNP A0R1R4
A	-2	GLY	-	EXPRESSION TAG	UNP A0R1R4
A	-1	SER	-	EXPRESSION TAG	UNP A0R1R4
A	1	VAL	-	EXPRESSION TAG	UNP A0R1R4
B	-21	MET	-	EXPRESSION TAG	UNP A0R1R4
B	-20	ALA	-	EXPRESSION TAG	UNP A0R1R4
B	-19	HIS	-	EXPRESSION TAG	UNP A0R1R4
B	-18	HIS	-	EXPRESSION TAG	UNP A0R1R4
B	-17	HIS	-	EXPRESSION TAG	UNP A0R1R4
B	-16	HIS	-	EXPRESSION TAG	UNP A0R1R4
B	-15	HIS	-	EXPRESSION TAG	UNP A0R1R4
B	-14	HIS	-	EXPRESSION TAG	UNP A0R1R4
B	-13	MET	-	EXPRESSION TAG	UNP A0R1R4
B	-12	GLY	-	EXPRESSION TAG	UNP A0R1R4
B	-11	THR	-	EXPRESSION TAG	UNP A0R1R4
B	-10	LEU	-	EXPRESSION TAG	UNP A0R1R4
B	-9	GLU	-	EXPRESSION TAG	UNP A0R1R4
B	-8	ALA	-	EXPRESSION TAG	UNP A0R1R4
B	-7	GLN	-	EXPRESSION TAG	UNP A0R1R4
B	-6	THR	-	EXPRESSION TAG	UNP A0R1R4
B	-5	GLN	-	EXPRESSION TAG	UNP A0R1R4
B	-4	GLY	-	EXPRESSION TAG	UNP A0R1R4
B	-3	PRO	-	EXPRESSION TAG	UNP A0R1R4
B	-2	GLY	-	EXPRESSION TAG	UNP A0R1R4
B	-1	SER	-	EXPRESSION TAG	UNP A0R1R4
B	1	VAL	-	EXPRESSION TAG	UNP A0R1R4
C	-21	MET	-	EXPRESSION TAG	UNP A0R1R4
C	-20	ALA	-	EXPRESSION TAG	UNP A0R1R4
C	-19	HIS	-	EXPRESSION TAG	UNP A0R1R4
C	-18	HIS	-	EXPRESSION TAG	UNP A0R1R4
C	-17	HIS	-	EXPRESSION TAG	UNP A0R1R4
C	-16	HIS	-	EXPRESSION TAG	UNP A0R1R4
C	-15	HIS	-	EXPRESSION TAG	UNP A0R1R4
C	-14	HIS	-	EXPRESSION TAG	UNP A0R1R4
C	-13	MET	-	EXPRESSION TAG	UNP A0R1R4
C	-12	GLY	-	EXPRESSION TAG	UNP A0R1R4
C	-11	THR	-	EXPRESSION TAG	UNP A0R1R4
C	-10	LEU	-	EXPRESSION TAG	UNP A0R1R4
C	-9	GLU	-	EXPRESSION TAG	UNP A0R1R4
C	-8	ALA	-	EXPRESSION TAG	UNP A0R1R4
C	-7	GLN	-	EXPRESSION TAG	UNP A0R1R4

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Chain	Residue	Modelled	Actual	Comment	Reference
C	-6	THR	-	EXPRESSION TAG	UNP A0R1R4
C	-5	GLN	-	EXPRESSION TAG	UNP A0R1R4
C	-4	GLY	-	EXPRESSION TAG	UNP A0R1R4
C	-3	PRO	-	EXPRESSION TAG	UNP A0R1R4
C	-2	GLY	-	EXPRESSION TAG	UNP A0R1R4
C	-1	SER	-	EXPRESSION TAG	UNP A0R1R4
C	1	VAL	-	EXPRESSION TAG	UNP A0R1R4
D	-21	MET	-	EXPRESSION TAG	UNP A0R1R4
D	-20	ALA	-	EXPRESSION TAG	UNP A0R1R4
D	-19	HIS	-	EXPRESSION TAG	UNP A0R1R4
D	-18	HIS	-	EXPRESSION TAG	UNP A0R1R4
D	-17	HIS	-	EXPRESSION TAG	UNP A0R1R4
D	-16	HIS	-	EXPRESSION TAG	UNP A0R1R4
D	-15	HIS	-	EXPRESSION TAG	UNP A0R1R4
D	-14	HIS	-	EXPRESSION TAG	UNP A0R1R4
D	-13	MET	-	EXPRESSION TAG	UNP A0R1R4
D	-12	GLY	-	EXPRESSION TAG	UNP A0R1R4
D	-11	THR	-	EXPRESSION TAG	UNP A0R1R4
D	-10	LEU	-	EXPRESSION TAG	UNP A0R1R4
D	-9	GLU	-	EXPRESSION TAG	UNP A0R1R4
D	-8	ALA	-	EXPRESSION TAG	UNP A0R1R4
D	-7	GLN	-	EXPRESSION TAG	UNP A0R1R4
D	-6	THR	-	EXPRESSION TAG	UNP A0R1R4
D	-5	GLN	-	EXPRESSION TAG	UNP A0R1R4
D	-4	GLY	-	EXPRESSION TAG	UNP A0R1R4
D	-3	PRO	-	EXPRESSION TAG	UNP A0R1R4
D	-2	GLY	-	EXPRESSION TAG	UNP A0R1R4
D	-1	SER	-	EXPRESSION TAG	UNP A0R1R4
D	1	VAL	-	EXPRESSION TAG	UNP A0R1R4
E	-21	MET	-	EXPRESSION TAG	UNP A0R1R4
E	-20	ALA	-	EXPRESSION TAG	UNP A0R1R4
E	-19	HIS	-	EXPRESSION TAG	UNP A0R1R4
E	-18	HIS	-	EXPRESSION TAG	UNP A0R1R4
E	-17	HIS	-	EXPRESSION TAG	UNP A0R1R4
E	-16	HIS	-	EXPRESSION TAG	UNP A0R1R4
E	-15	HIS	-	EXPRESSION TAG	UNP A0R1R4
E	-14	HIS	-	EXPRESSION TAG	UNP A0R1R4
E	-13	MET	-	EXPRESSION TAG	UNP A0R1R4
E	-12	GLY	-	EXPRESSION TAG	UNP A0R1R4
E	-11	THR	-	EXPRESSION TAG	UNP A0R1R4
E	-10	LEU	-	EXPRESSION TAG	UNP A0R1R4
E	-9	GLU	-	EXPRESSION TAG	UNP A0R1R4

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Chain	Residue	Modelled	Actual	Comment	Reference
E	-8	ALA	-	EXPRESSION TAG	UNP A0R1R4
E	-7	GLN	-	EXPRESSION TAG	UNP A0R1R4
E	-6	THR	-	EXPRESSION TAG	UNP A0R1R4
E	-5	GLN	-	EXPRESSION TAG	UNP A0R1R4
E	-4	GLY	-	EXPRESSION TAG	UNP A0R1R4
E	-3	PRO	-	EXPRESSION TAG	UNP A0R1R4
E	-2	GLY	-	EXPRESSION TAG	UNP A0R1R4
E	-1	SER	-	EXPRESSION TAG	UNP A0R1R4
E	1	VAL	-	EXPRESSION TAG	UNP A0R1R4
F	-21	MET	-	EXPRESSION TAG	UNP A0R1R4
F	-20	ALA	-	EXPRESSION TAG	UNP A0R1R4
F	-19	HIS	-	EXPRESSION TAG	UNP A0R1R4
F	-18	HIS	-	EXPRESSION TAG	UNP A0R1R4
F	-17	HIS	-	EXPRESSION TAG	UNP A0R1R4
F	-16	HIS	-	EXPRESSION TAG	UNP A0R1R4
F	-15	HIS	-	EXPRESSION TAG	UNP A0R1R4
F	-14	HIS	-	EXPRESSION TAG	UNP A0R1R4
F	-13	MET	-	EXPRESSION TAG	UNP A0R1R4
F	-12	GLY	-	EXPRESSION TAG	UNP A0R1R4
F	-11	THR	-	EXPRESSION TAG	UNP A0R1R4
F	-10	LEU	-	EXPRESSION TAG	UNP A0R1R4
F	-9	GLU	-	EXPRESSION TAG	UNP A0R1R4
F	-8	ALA	-	EXPRESSION TAG	UNP A0R1R4
F	-7	GLN	-	EXPRESSION TAG	UNP A0R1R4
F	-6	THR	-	EXPRESSION TAG	UNP A0R1R4
F	-5	GLN	-	EXPRESSION TAG	UNP A0R1R4
F	-4	GLY	-	EXPRESSION TAG	UNP A0R1R4
F	-3	PRO	-	EXPRESSION TAG	UNP A0R1R4
F	-2	GLY	-	EXPRESSION TAG	UNP A0R1R4
F	-1	SER	-	EXPRESSION TAG	UNP A0R1R4
F	1	VAL	-	EXPRESSION TAG	UNP A0R1R4

- Molecule 2 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	D	1	Total Zn 1 1	0	0
2	E	1	Total Zn 1 1	0	0
2	B	1	Total Zn 1 1	0	0
2	C	1	Total Zn 1 1	0	0

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total 1	Zn 1	0	0
2	F	1	Total 1	Zn 1	0	0

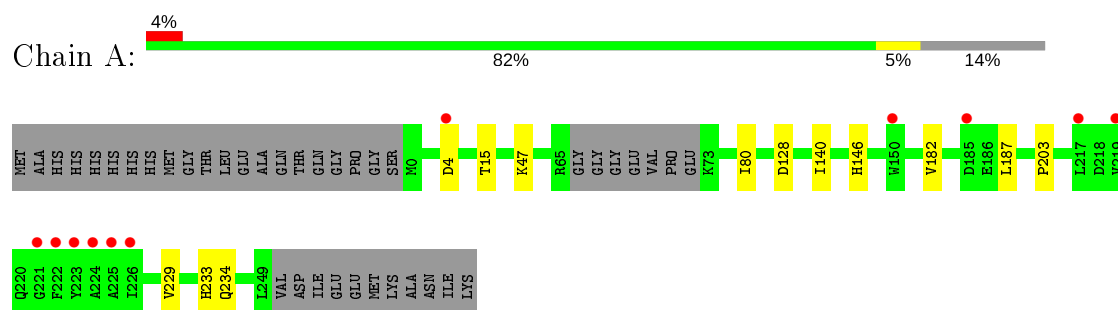
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	112	Total 112	O 112	0	0
3	B	115	Total 115	O 115	0	0
3	C	141	Total 141	O 141	0	0
3	D	95	Total 95	O 95	0	0
3	E	120	Total 120	O 120	0	0
3	F	83	Total 83	O 83	0	0

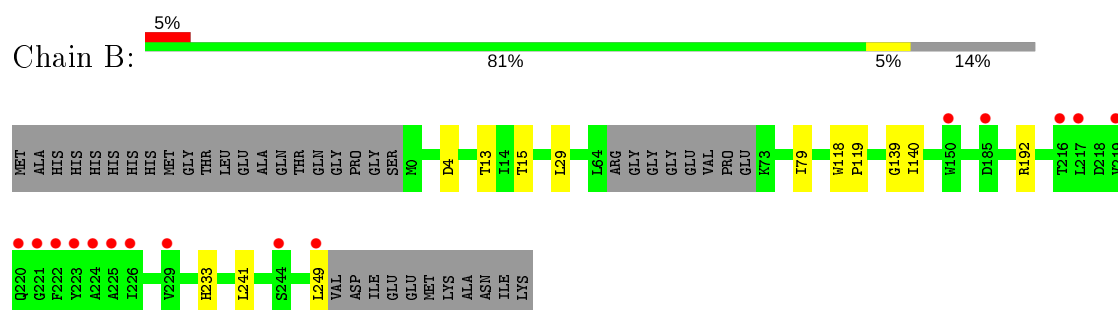
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

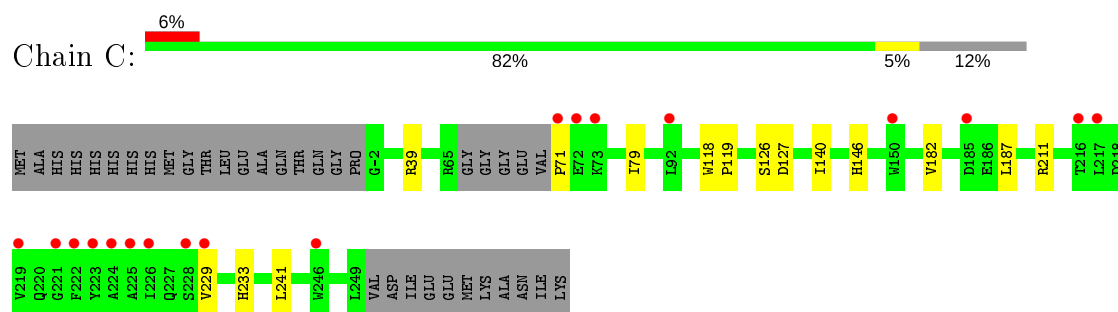
• Molecule 1: Enoyl-CoA hydratase



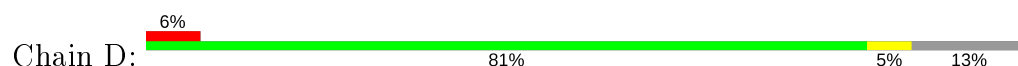
• Molecule 1: Enoyl-CoA hydratase

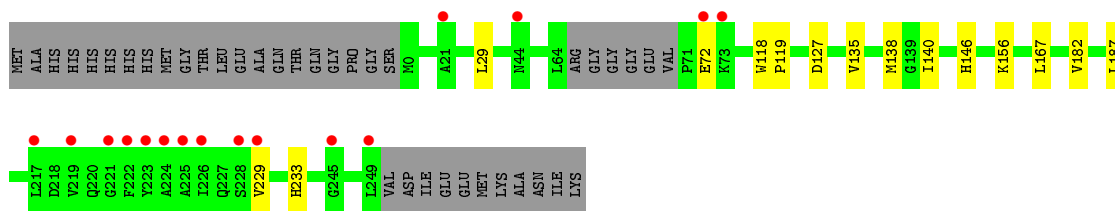


• Molecule 1: Enoyl-CoA hydratase

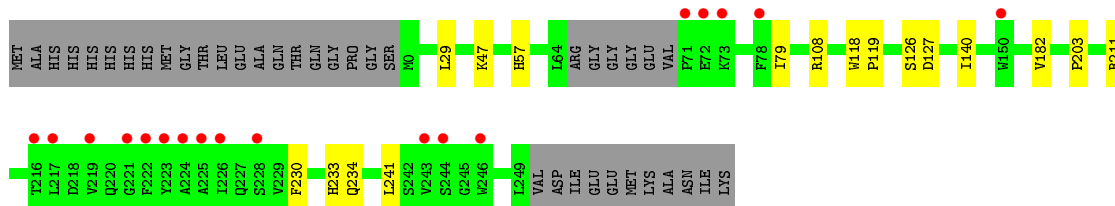
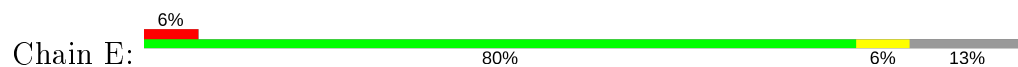


• Molecule 1: Enoyl-CoA hydratase

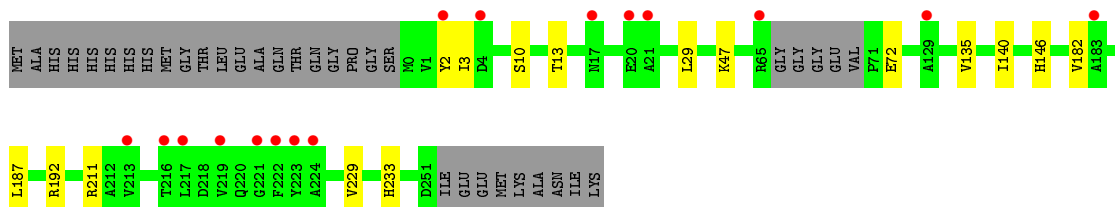
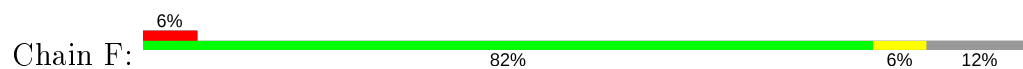




• Molecule 1: Enoyl-CoA hydratase



• Molecule 1: Enoyl-CoA hydratase



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	145.22Å 86.65Å 140.05Å 90.00° 97.93° 90.00°	Depositor
Resolution (Å)	50.00 – 2.05 48.97 – 2.05	Depositor EDS
% Data completeness (in resolution range)	99.8 (50.00-2.05) 99.8 (48.97-2.05)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.45 (at 2.05Å)	Xtriage
Refinement program	REFMAC 5.5.0109	Depositor
R, R_{free}	0.145 , 0.174 0.172 , 0.196	Depositor DCC
R_{free} test set	5373 reflections (4.99%)	wwPDB-VP
Wilson B-factor (Å ²)	29.6	Xtriage
Anisotropy	0.461	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 41.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	11909	wwPDB-VP
Average B, all atoms (Å ²)	35.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.07% 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: ZN

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	0.72	0/1906	0.73	1/2591 (0.0%)
1	B	0.72	0/1901	0.70	0/2585
1	C	0.75	0/1936	0.74	2/2635 (0.1%)
1	D	0.63	0/1891	0.65	0/2576
1	E	0.73	0/1925	0.70	1/2619 (0.0%)
1	F	0.65	0/1925	0.70	1/2624 (0.0%)
All	All	0.70	0/11484	0.70	5/15630 (0.0%)

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	F	211	ARG	NE-CZ-NH2	6.93	123.77	120.30
1	A	128	ASP	CB-CG-OD2	6.90	124.51	118.30
1	E	211	ARG	NE-CZ-NH2	6.44	123.52	120.30
1	C	211	ARG	NE-CZ-NH2	6.26	123.43	120.30
1	C	39	ARG	NE-CZ-NH2	-5.04	117.78	120.30

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	1867	0	1832	6	0
1	B	1862	0	1828	7	0
1	C	1894	0	1845	6	0
1	D	1851	0	1796	8	0
1	E	1883	0	1841	8	0
1	F	1880	0	1829	8	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
2	E	1	0	0	0	0
2	F	1	0	0	0	0
3	A	112	0	0	0	0
3	B	115	0	0	0	0
3	C	141	0	0	0	0
3	D	95	0	0	0	0
3	E	120	0	0	0	0
3	F	83	0	0	0	0
All	All	11909	0	10971	43	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 (43) 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:D:135:VAL:HG23	1:D:138:MET:CE	2.23	0.68
1:D:135:VAL:HG23	1:D:138:MET:HE2	1.78	0.64
1:F:10:SER:HB3	1:F:47:LYS:HG3	1.86	0.58
1:C:182:VAL:HG21	1:C:187:LEU:HA	1.88	0.55
1:B:140:ILE:HG21	1:B:233:HIS:CE1	2.41	0.55
1:D:140:ILE:HG21	1:D:233:HIS:CE1	2.44	0.53
1:A:47:LYS:HE3	1:A:203:PRO:HG3	1.91	0.52
1:C:79:ILE:HD12	1:C:241:LEU:HD11	1.92	0.51
1:C:146:HIS:HB2	1:C:229:VAL:HG11	1.95	0.48
1:B:118:TRP:N	1:B:119:PRO:HD2	2.28	0.48
1:F:13:THR:HG1	1:F:192[B]:ARG:HH12	1.56	0.47
1:B:4:ASP:HB2	1:B:15:THR:HB	1.97	0.47
1:D:146:HIS:HB2	1:D:229:VAL:HG11	1.96	0.47
1:F:13:THR:OG1	1:F:192[B]:ARG:NH1	2.38	0.47
1:E:140:ILE:HG21	1:E:233:HIS:CE1	2.50	0.46
1:D:118:TRP:N	1:D:119:PRO:HD2	2.30	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:146:HIS:HB2	1:A:229:VAL:HG11	1.96	0.46
1:E:126:SER:HA	1:E:182:VAL:O	2.15	0.46
1:F:182:VAL:HG21	1:F:187:LEU:HA	1.98	0.45
1:D:182:VAL:HG21	1:D:187:LEU:HA	1.99	0.45
1:A:4:ASP:HB2	1:A:15:THR:HB	1.99	0.44
1:A:80:ILE:HG12	1:A:234:GLN:HG3	2.00	0.44
1:F:146:HIS:HB2	1:F:229:VAL:HG11	1.98	0.44
1:B:29:LEU:HD23	1:B:29:LEU:C	2.38	0.43
1:F:29:LEU:HD23	1:F:29:LEU:C	2.38	0.43
1:F:140:ILE:HG21	1:F:233:HIS:CE1	2.52	0.43
1:C:118:TRP:N	1:C:119:PRO:HD2	2.33	0.43
1:E:118:TRP:N	1:E:119:PRO:HD2	2.32	0.43
1:E:57:HIS:CE1	1:E:108:ARG:HG3	2.54	0.43
1:F:2:TYR:O	1:F:3:ILE:HD13	2.19	0.42
1:D:156:LYS:HD2	1:D:167:LEU:HD21	2.02	0.42
1:A:182:VAL:HG21	1:A:187:LEU:HA	2.03	0.41
1:D:29:LEU:HD23	1:D:29:LEU:C	2.41	0.41
1:E:79:ILE:HD12	1:E:241:LEU:HD11	2.02	0.41
1:C:140:ILE:HG21	1:C:233:HIS:CE1	2.56	0.41
1:B:13:THR:OG1	1:B:192[B]:ARG:NH1	2.48	0.41
1:C:126:SER:HA	1:C:182:VAL:O	2.20	0.41
1:E:230:PHE:O	1:E:234:GLN:HG2	2.21	0.41
1:B:79:ILE:HD12	1:B:241:LEU:HD11	2.03	0.41
1:B:139:GLY:CA	1:B:249:LEU:HD12	2.51	0.41
1:E:29:LEU:C	1:E:29:LEU:HD23	2.41	0.41
1:A:140:ILE:HG21	1:A:233:HIS:CE1	2.56	0.40
1:E:47:LYS:HE3	1:E:203:PRO:HG3	2.02	0.40

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	240/282 (85%)	233 (97%)	7 (3%)	0	100	100
1	B	239/282 (85%)	232 (97%)	7 (3%)	0	100	100
1	C	244/282 (86%)	235 (96%)	9 (4%)	0	100	100
1	D	241/282 (86%)	234 (97%)	6 (2%)	1 (0%)	34	24
1	E	241/282 (86%)	233 (97%)	8 (3%)	0	100	100
1	F	245/282 (87%)	236 (96%)	8 (3%)	1 (0%)	34	24
All	All	1450/1692 (86%)	1403 (97%)	45 (3%)	2 (0%)	51	45

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	72	GLU
1	F	72	GLU

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	189/225 (84%)	189 (100%)	0	100	100
1	B	189/225 (84%)	189 (100%)	0	100	100
1	C	191/225 (85%)	189 (99%)	2 (1%)	76	75
1	D	185/225 (82%)	184 (100%)	1 (0%)	88	89
1	E	191/225 (85%)	190 (100%)	1 (0%)	88	89
1	F	188/225 (84%)	187 (100%)	1 (0%)	88	89
All	All	1133/1350 (84%)	1128 (100%)	5 (0%)	91	91

All (5) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	71	PRO
1	C	127	ASP
1	D	127	ASP

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Mol	Chain	Res	Type
1	E	127	ASP
1	F	135	VAL

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

Of 6 ligands modelled in this entry, 6 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 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	243/282 (86%)	0.03	11 (4%) 33 35	16, 32, 49, 63	0
1	B	242/282 (85%)	0.10	15 (6%) 20 22	18, 33, 54, 63	0
1	C	247/282 (87%)	0.06	18 (7%) 15 16	15, 28, 48, 64	0
1	D	244/282 (86%)	0.13	16 (6%) 18 19	17, 41, 61, 74	0
1	E	244/282 (86%)	0.14	18 (7%) 14 16	17, 32, 52, 77	0
1	F	247/282 (87%)	0.24	16 (6%) 18 20	19, 43, 64, 72	0
All	All	1467/1692 (86%)	0.12	94 (6%) 19 21	15, 34, 59, 77	0

All (94) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	246	TRP	6.2
1	E	219	VAL	4.7
1	E	72	GLU	4.7
1	A	219	VAL	4.5
1	E	71	PRO	4.2
1	E	217	LEU	4.0
1	B	219	VAL	4.0
1	F	219	VAL	3.9
1	C	219	VAL	3.8
1	C	246	TRP	3.6
1	E	224	ALA	3.6
1	C	72	GLU	3.4
1	E	223	TYR	3.3
1	D	73	LYS	3.3
1	C	224	ALA	3.3
1	F	65	ARG	3.3
1	C	223	TYR	3.3
1	D	217	LEU	3.2
1	D	219	VAL	3.2

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Mol	Chain	Res	Type	RSRZ
1	B	221	GLY	3.1
1	C	222	PHE	3.1
1	E	73	LYS	3.0
1	E	222	PHE	3.0
1	C	226	ILE	3.0
1	E	78	PHE	3.0
1	B	223	TYR	3.0
1	B	224	ALA	3.0
1	B	217	LEU	2.9
1	B	185	ASP	2.9
1	C	150	TRP	2.9
1	A	217	LEU	2.9
1	A	222	PHE	2.8
1	C	217	LEU	2.8
1	B	226	ILE	2.8
1	E	221	GLY	2.8
1	A	221	GLY	2.8
1	C	92	LEU	2.8
1	E	243	VAL	2.7
1	D	72	GLU	2.7
1	D	225	ALA	2.7
1	F	2	TYR	2.7
1	D	222	PHE	2.6
1	B	225	ALA	2.6
1	A	224	ALA	2.6
1	B	216	THR	2.6
1	C	73	LYS	2.6
1	B	249	LEU	2.6
1	F	217	LEU	2.6
1	C	225	ALA	2.6
1	E	225	ALA	2.6
1	E	150	TRP	2.6
1	F	183	ALA	2.5
1	F	224	ALA	2.5
1	A	226	ILE	2.5
1	C	221	GLY	2.5
1	D	223	TYR	2.5
1	D	224	ALA	2.5
1	D	221	GLY	2.5
1	C	229	VAL	2.4
1	B	244	SER	2.4
1	D	226	ILE	2.4

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Mol	Chain	Res	Type	RSRZ
1	F	221	GLY	2.4
1	F	21	ALA	2.4
1	A	150	TRP	2.4
1	B	150	TRP	2.3
1	C	216	THR	2.3
1	A	225	ALA	2.3
1	D	229	VAL	2.3
1	D	228	SER	2.3
1	F	223	TYR	2.3
1	B	229	VAL	2.3
1	D	245	GLY	2.3
1	E	244	SER	2.3
1	B	222	PHE	2.2
1	F	17	ASN	2.2
1	D	249	LEU	2.2
1	C	71	PRO	2.2
1	A	223	TYR	2.2
1	E	216	THR	2.1
1	F	213	VAL	2.1
1	C	228	SER	2.1
1	F	222	PHE	2.1
1	F	20	GLU	2.1
1	E	228	SER	2.1
1	B	220	GLN	2.1
1	A	185	ASP	2.0
1	E	226	ILE	2.0
1	D	21	ALA	2.0
1	C	185	ASP	2.0
1	D	44	ASN	2.0
1	F	129	ALA	2.0
1	F	216	THR	2.0
1	A	4	ASP	2.0
1	F	4	ASP	2.0

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

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

6.4 Ligands [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
2	ZN	D	300	1/1	0.99	0.06	37,37,37,37	0
2	ZN	B	300	1/1	0.99	0.05	41,41,41,41	0
2	ZN	F	300	1/1	1.00	0.06	41,41,41,41	0
2	ZN	C	300	1/1	1.00	0.07	29,29,29,29	0
2	ZN	A	300	1/1	1.00	0.04	36,36,36,36	0
2	ZN	E	300	1/1	1.00	0.06	40,40,40,40	0

6.5 Other polymers [i](#)

There are no such residues in this entry.