



wwPDB X-ray Structure Validation Summary Report ⓘ

May 13, 2020 – 02:15 am BST

PDB ID : 4CP8
Title : Structure of the amidase domain of allophanate hydrolase from *Pseudomonas* sp strain ADP
Authors : Balotra, S.; Newman, J.; French, N.; French, L.; Peat, T.S.; Scott, C.
Deposited on : 2014-02-03
Resolution : 2.50 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

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

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

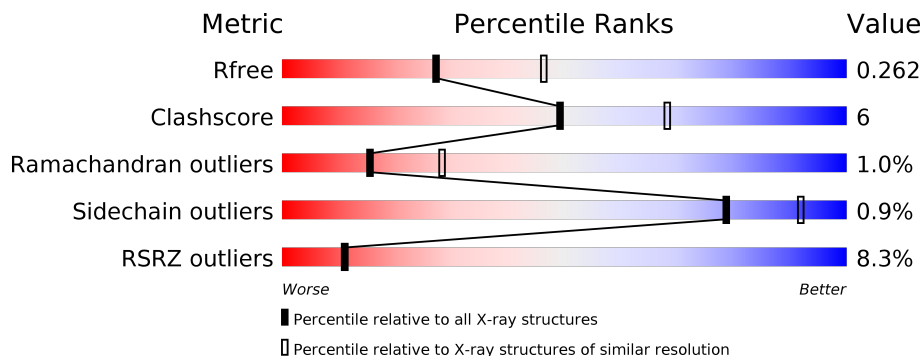
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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	4661 (2.50-2.50)
Clashscore	141614	5346 (2.50-2.50)
Ramachandran outliers	138981	5231 (2.50-2.50)
Sidechain outliers	138945	5233 (2.50-2.50)
RSRZ outliers	127900	4559 (2.50-2.50)

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	487	 2% 84% 7% • 8%
1	B	487	 2% 86% 6% 8%
1	C	487	 3% 86% 6% 8%
1	D	487	 2% 85% 6% • 8%
1	E	487	 18% 80% 11% • 8%
1	F	487	 19% 78% 12% • 9%

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	MLI	A	1466	-	-	X	-
2	MLI	B	1465	-	-	X	-
2	MLI	C	1466	-	-	X	-
2	MLI	D	1465	-	-	X	-

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 20462 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 ALLOPHANATE HYDROLASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	448	3375	2146	580	639	10	0	2	0
1	B	447	3345	2126	574	635	10	0	0	0
1	C	448	3355	2132	577	636	10	0	0	0
1	D	447	3356	2132	577	637	10	0	1	0
1	E	450	3373	2144	580	639	10	0	1	0
1	F	444	3318	2110	568	630	10	0	0	0

There are 120 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-19	MET	-	expression tag	UNP Q936X2
A	-18	GLY	-	expression tag	UNP Q936X2
A	-17	SER	-	expression tag	UNP Q936X2
A	-16	SER	-	expression tag	UNP Q936X2
A	-15	HIS	-	expression tag	UNP Q936X2
A	-14	HIS	-	expression tag	UNP Q936X2
A	-13	HIS	-	expression tag	UNP Q936X2
A	-12	HIS	-	expression tag	UNP Q936X2
A	-11	HIS	-	expression tag	UNP Q936X2
A	-10	HIS	-	expression tag	UNP Q936X2
A	-9	SER	-	expression tag	UNP Q936X2
A	-8	SER	-	expression tag	UNP Q936X2
A	-7	GLY	-	expression tag	UNP Q936X2
A	-6	LEU	-	expression tag	UNP Q936X2
A	-5	VAL	-	expression tag	UNP Q936X2
A	-4	PRO	-	expression tag	UNP Q936X2
A	-3	ARG	-	expression tag	UNP Q936X2

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Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	GLY	-	expression tag	UNP Q936X2
A	-1	SER	-	expression tag	UNP Q936X2
A	0	HIS	-	expression tag	UNP Q936X2
B	-19	MET	-	expression tag	UNP Q936X2
B	-18	GLY	-	expression tag	UNP Q936X2
B	-17	SER	-	expression tag	UNP Q936X2
B	-16	SER	-	expression tag	UNP Q936X2
B	-15	HIS	-	expression tag	UNP Q936X2
B	-14	HIS	-	expression tag	UNP Q936X2
B	-13	HIS	-	expression tag	UNP Q936X2
B	-12	HIS	-	expression tag	UNP Q936X2
B	-11	HIS	-	expression tag	UNP Q936X2
B	-10	HIS	-	expression tag	UNP Q936X2
B	-9	SER	-	expression tag	UNP Q936X2
B	-8	SER	-	expression tag	UNP Q936X2
B	-7	GLY	-	expression tag	UNP Q936X2
B	-6	LEU	-	expression tag	UNP Q936X2
B	-5	VAL	-	expression tag	UNP Q936X2
B	-4	PRO	-	expression tag	UNP Q936X2
B	-3	ARG	-	expression tag	UNP Q936X2
B	-2	GLY	-	expression tag	UNP Q936X2
B	-1	SER	-	expression tag	UNP Q936X2
B	0	HIS	-	expression tag	UNP Q936X2
C	-19	MET	-	expression tag	UNP Q936X2
C	-18	GLY	-	expression tag	UNP Q936X2
C	-17	SER	-	expression tag	UNP Q936X2
C	-16	SER	-	expression tag	UNP Q936X2
C	-15	HIS	-	expression tag	UNP Q936X2
C	-14	HIS	-	expression tag	UNP Q936X2
C	-13	HIS	-	expression tag	UNP Q936X2
C	-12	HIS	-	expression tag	UNP Q936X2
C	-11	HIS	-	expression tag	UNP Q936X2
C	-10	HIS	-	expression tag	UNP Q936X2
C	-9	SER	-	expression tag	UNP Q936X2
C	-8	SER	-	expression tag	UNP Q936X2
C	-7	GLY	-	expression tag	UNP Q936X2
C	-6	LEU	-	expression tag	UNP Q936X2
C	-5	VAL	-	expression tag	UNP Q936X2
C	-4	PRO	-	expression tag	UNP Q936X2
C	-3	ARG	-	expression tag	UNP Q936X2
C	-2	GLY	-	expression tag	UNP Q936X2
C	-1	SER	-	expression tag	UNP Q936X2

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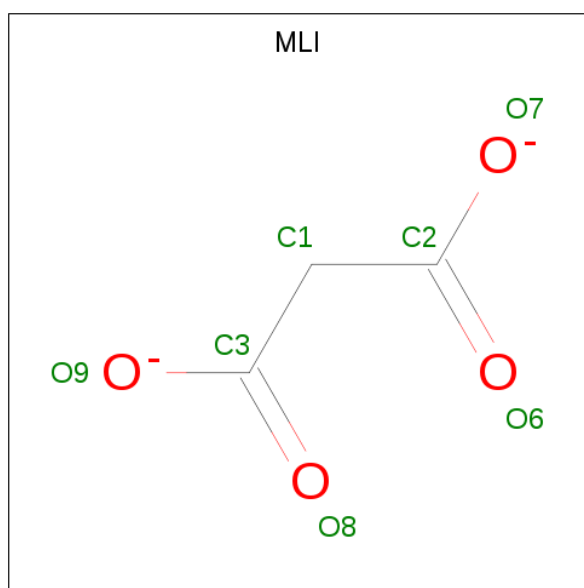
Chain	Residue	Modelled	Actual	Comment	Reference
C	0	HIS	-	expression tag	UNP Q936X2
D	-19	MET	-	expression tag	UNP Q936X2
D	-18	GLY	-	expression tag	UNP Q936X2
D	-17	SER	-	expression tag	UNP Q936X2
D	-16	SER	-	expression tag	UNP Q936X2
D	-15	HIS	-	expression tag	UNP Q936X2
D	-14	HIS	-	expression tag	UNP Q936X2
D	-13	HIS	-	expression tag	UNP Q936X2
D	-12	HIS	-	expression tag	UNP Q936X2
D	-11	HIS	-	expression tag	UNP Q936X2
D	-10	HIS	-	expression tag	UNP Q936X2
D	-9	SER	-	expression tag	UNP Q936X2
D	-8	SER	-	expression tag	UNP Q936X2
D	-7	GLY	-	expression tag	UNP Q936X2
D	-6	LEU	-	expression tag	UNP Q936X2
D	-5	VAL	-	expression tag	UNP Q936X2
D	-4	PRO	-	expression tag	UNP Q936X2
D	-3	ARG	-	expression tag	UNP Q936X2
D	-2	GLY	-	expression tag	UNP Q936X2
D	-1	SER	-	expression tag	UNP Q936X2
D	0	HIS	-	expression tag	UNP Q936X2
E	-19	MET	-	expression tag	UNP Q936X2
E	-18	GLY	-	expression tag	UNP Q936X2
E	-17	SER	-	expression tag	UNP Q936X2
E	-16	SER	-	expression tag	UNP Q936X2
E	-15	HIS	-	expression tag	UNP Q936X2
E	-14	HIS	-	expression tag	UNP Q936X2
E	-13	HIS	-	expression tag	UNP Q936X2
E	-12	HIS	-	expression tag	UNP Q936X2
E	-11	HIS	-	expression tag	UNP Q936X2
E	-10	HIS	-	expression tag	UNP Q936X2
E	-9	SER	-	expression tag	UNP Q936X2
E	-8	SER	-	expression tag	UNP Q936X2
E	-7	GLY	-	expression tag	UNP Q936X2
E	-6	LEU	-	expression tag	UNP Q936X2
E	-5	VAL	-	expression tag	UNP Q936X2
E	-4	PRO	-	expression tag	UNP Q936X2
E	-3	ARG	-	expression tag	UNP Q936X2
E	-2	GLY	-	expression tag	UNP Q936X2
E	-1	SER	-	expression tag	UNP Q936X2
E	0	HIS	-	expression tag	UNP Q936X2
F	-19	MET	-	expression tag	UNP Q936X2

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Chain	Residue	Modelled	Actual	Comment	Reference
F	-18	GLY	-	expression tag	UNP Q936X2
F	-17	SER	-	expression tag	UNP Q936X2
F	-16	SER	-	expression tag	UNP Q936X2
F	-15	HIS	-	expression tag	UNP Q936X2
F	-14	HIS	-	expression tag	UNP Q936X2
F	-13	HIS	-	expression tag	UNP Q936X2
F	-12	HIS	-	expression tag	UNP Q936X2
F	-11	HIS	-	expression tag	UNP Q936X2
F	-10	HIS	-	expression tag	UNP Q936X2
F	-9	SER	-	expression tag	UNP Q936X2
F	-8	SER	-	expression tag	UNP Q936X2
F	-7	GLY	-	expression tag	UNP Q936X2
F	-6	LEU	-	expression tag	UNP Q936X2
F	-5	VAL	-	expression tag	UNP Q936X2
F	-4	PRO	-	expression tag	UNP Q936X2
F	-3	ARG	-	expression tag	UNP Q936X2
F	-2	GLY	-	expression tag	UNP Q936X2
F	-1	SER	-	expression tag	UNP Q936X2
F	0	HIS	-	expression tag	UNP Q936X2

- Molecule 2 is MALONATE ION (three-letter code: MLI) (formula: C₃H₂O₄).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			7	3	4		
2	B	1	Total	C	O	0	0
			7	3	4		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	C	1	Total	C	O	0	0
			7	3	4		
2	D	1	Total	C	O	0	0
			7	3	4		

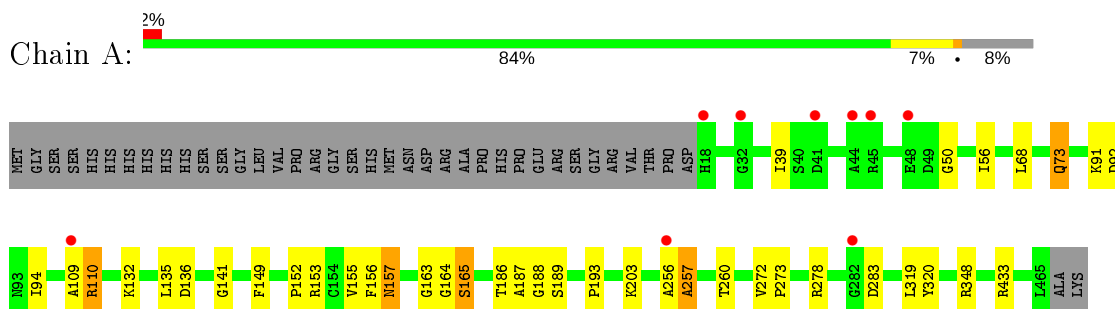
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	86	Total	O	0	0
			86	86		
3	B	71	Total	O	0	0
			71	71		
3	C	84	Total	O	0	0
			84	84		
3	D	51	Total	O	0	0
			51	51		
3	E	18	Total	O	0	0
			18	18		
3	F	2	Total	O	0	0
			2	2		

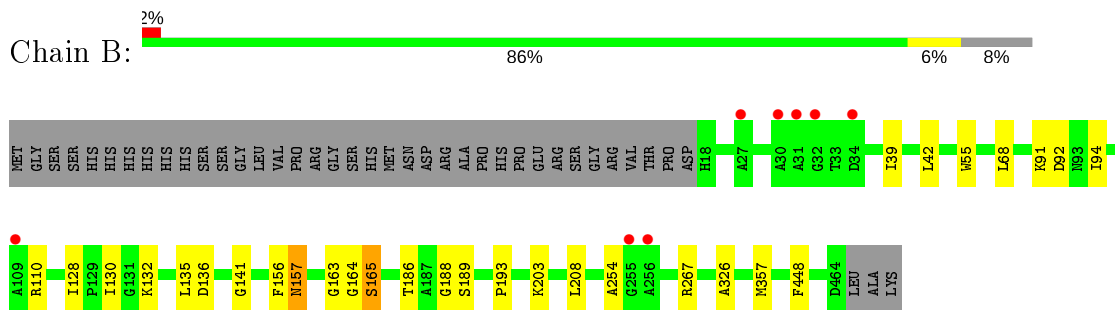
3 Residue-property plots

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.

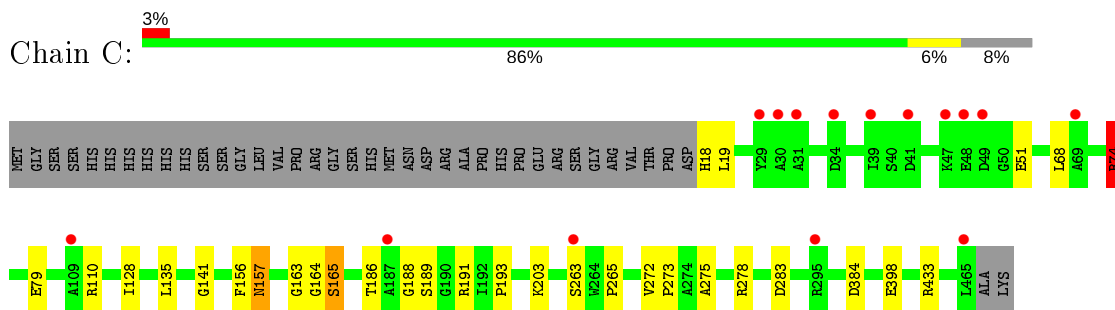
- Molecule 1: ALLOPHANATE HYDROLASE



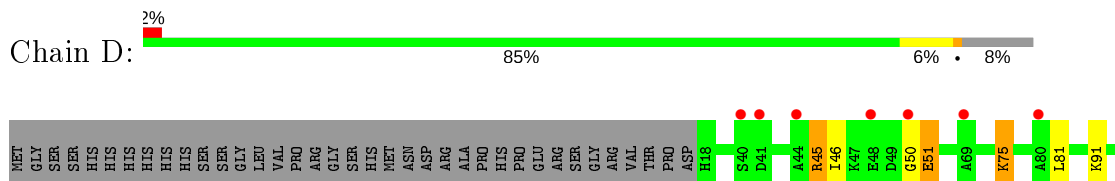
- Molecule 1: ALLOPHANATE HYDROLASE



- Molecule 1: ALLOPHANATE HYDROLASE

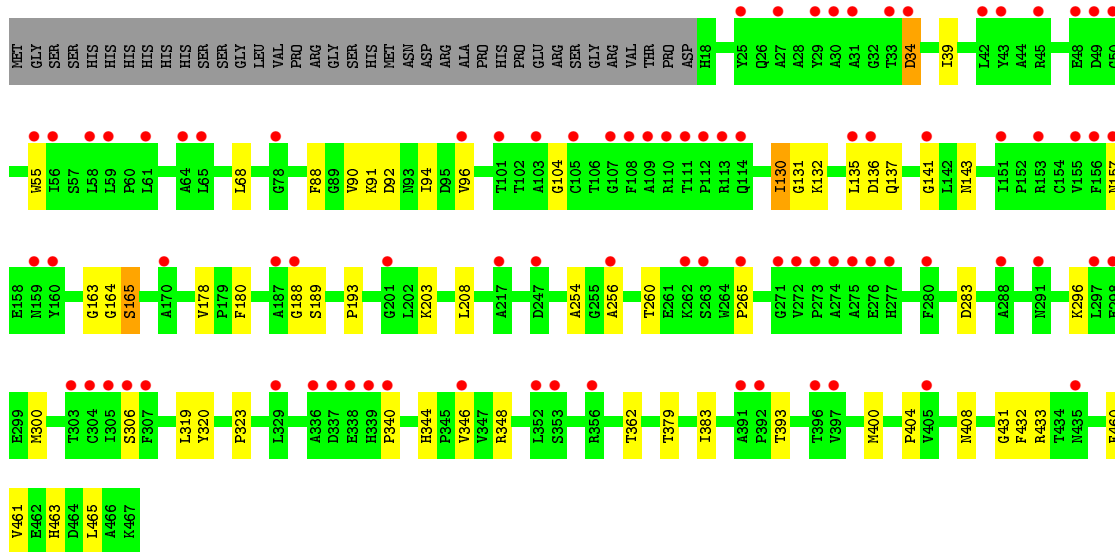
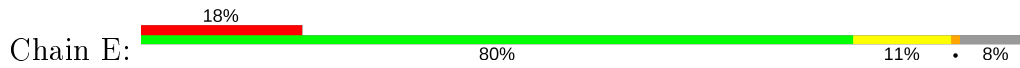


- Molecule 1: ALLOPHANATE HYDROLASE

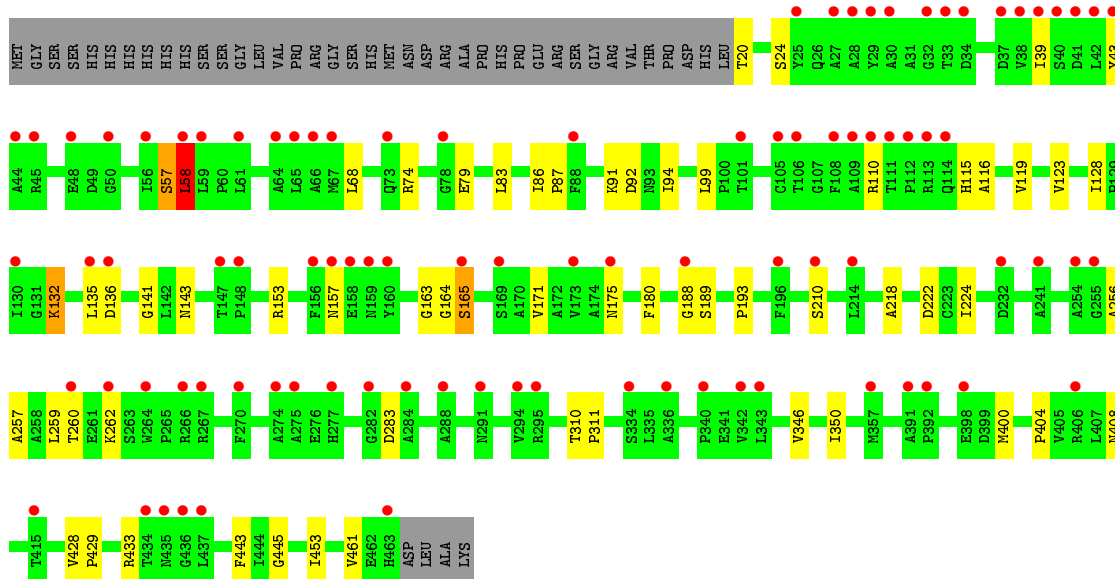
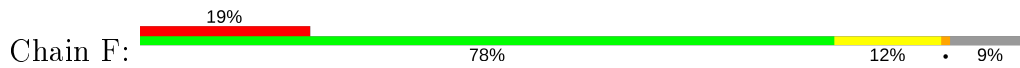




● Molecule 1: ALLOPHANATE HYDROLASE



● Molecule 1: ALLOPHANATE HYDROLASE



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	82.45Å 179.23Å 112.61Å 90.00° 106.63° 90.00°	Depositor
Resolution (Å)	40.00 – 2.50 39.59 – 2.50	Depositor EDS
% Data completeness (in resolution range)	100.0 (40.00-2.50) 100.0 (39.59-2.50)	Depositor EDS
R_{merge}	0.15	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.72 (at 2.51Å)	Xtrriage
Refinement program	REFMAC 5.7.0032	Depositor
R, R_{free}	0.224 , 0.259 0.226 , 0.262	Depositor DCC
R_{free} test set	5383 reflections (4.99%)	wwPDB-VP
Wilson B-factor (Å ²)	33.9	Xtrriage
Anisotropy	0.072	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 32.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	20462	wwPDB-VP
Average B, all atoms (Å ²)	44.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 11.37% 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](#)

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

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	0.50	0/3449	0.69	1/4696 (0.0%)
1	B	0.51	0/3418	0.67	0/4654
1	C	0.52	0/3429	0.68	1/4669 (0.0%)
1	D	0.50	0/3430	0.66	1/4670 (0.0%)
1	E	0.44	0/3447	0.67	1/4694 (0.0%)
1	F	0.44	0/3391	0.70	2/4618 (0.0%)
All	All	0.49	0/20564	0.68	6/28001 (0.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1
1	B	0	1
1	C	0	1
1	D	0	1
1	E	0	1
1	F	0	1
All	All	0	6

There are no bond length outliers.

The worst 5 of 6 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	74	ARG	NE-CZ-NH1	7.24	123.92	120.30
1	E	34	ASP	N-CA-CB	5.99	121.38	110.60
1	F	57	SER	CB-CA-C	-5.83	99.02	110.10
1	F	58	LEU	CA-CB-CG	5.17	127.19	115.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	348	ARG	NE-CZ-NH1	5.06	122.83	120.30

There are no chirality outliers.

5 of 6 planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	163	GLY	Peptide
1	B	163	GLY	Peptide
1	C	163	GLY	Peptide
1	D	163	GLY	Peptide
1	E	163	GLY	Peptide

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3375	0	3328	35	0
1	B	3345	0	3295	27	0
1	C	3355	0	3302	28	0
1	D	3356	0	3304	25	0
1	E	3373	0	3318	59	0
1	F	3318	0	3267	64	0
2	A	7	0	2	5	0
2	B	7	0	2	5	0
2	C	7	0	2	6	0
2	D	7	0	2	5	0
3	A	86	0	0	2	0
3	B	71	0	0	0	0
3	C	84	0	0	2	0
3	D	51	0	0	1	0
3	E	18	0	0	2	0
3	F	2	0	0	1	0
All	All	20462	0	19822	235	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 6.

The worst 5 of 235 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:E:94:ILE:O	1:E:132:LYS:NZ	1.58	1.34
1:F:115:HIS:HB3	1:F:119:VAL:HG13	1.39	1.02
1:F:86:ILE:HD12	1:F:180:PHE:HE1	1.27	0.95
1:E:340:PRO:HB2	1:E:348:ARG:HE	1.35	0.92
1:D:189:SER:HG	2:D:1465:MLI:C2	1.83	0.90

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	448/487 (92%)	429 (96%)	14 (3%)	5 (1%)	14	26
1	B	445/487 (91%)	427 (96%)	14 (3%)	4 (1%)	17	31
1	C	446/487 (92%)	426 (96%)	16 (4%)	4 (1%)	17	31
1	D	446/487 (92%)	429 (96%)	12 (3%)	5 (1%)	14	26
1	E	449/487 (92%)	431 (96%)	14 (3%)	4 (1%)	17	31
1	F	442/487 (91%)	421 (95%)	16 (4%)	5 (1%)	14	26
All	All	2676/2922 (92%)	2563 (96%)	86 (3%)	27 (1%)	15	28

5 of 27 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	165	SER
1	B	165	SER
1	C	165	SER
1	D	51	GLU
1	D	165	SER

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	344/375 (92%)	340 (99%)	4 (1%)	71	88
1	B	340/375 (91%)	339 (100%)	1 (0%)	92	97
1	C	341/375 (91%)	337 (99%)	4 (1%)	71	88
1	D	342/375 (91%)	337 (98%)	5 (2%)	65	85
1	E	342/375 (91%)	340 (99%)	2 (1%)	86	95
1	F	337/375 (90%)	333 (99%)	4 (1%)	71	88
All	All	2046/2250 (91%)	2026 (99%)	20 (1%)	78	90

5 of 20 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	D	45	ARG
1	D	75	LYS
1	F	58	LEU
1	C	203	LYS
1	C	398	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	C	72	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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

4 ligands 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	$\# Z > 2$
2	MLI	C	1466	-	0,6,6	0.00	-	0,7,7	0.00	-
2	MLI	B	1465	-	0,6,6	0.00	-	0,7,7	0.00	-
2	MLI	A	1466	-	0,6,6	0.00	-	0,7,7	0.00	-
2	MLI	D	1465	-	0,6,6	0.00	-	0,7,7	0.00	-

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
2	MLI	C	1466	-	-	0/0/4/4	-
2	MLI	B	1465	-	-	0/0/4/4	-
2	MLI	A	1466	-	-	0/0/4/4	-
2	MLI	D	1465	-	-	0/0/4/4	-

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.

4 monomers are involved in 21 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	1466	MLI	6	0
2	B	1465	MLI	5	0
2	A	1466	MLI	5	0
2	D	1465	MLI	5	0

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	448/487 (91%)	-0.00	9 (2%) 65 68	16, 26, 59, 96	0
1	B	447/487 (91%)	-0.04	8 (1%) 68 71	16, 28, 59, 84	0
1	C	448/487 (91%)	0.07	15 (3%) 46 50	14, 25, 61, 81	0
1	D	447/487 (91%)	0.12	11 (2%) 57 61	15, 28, 58, 95	0
1	E	450/487 (92%)	1.22	86 (19%) 1 1	35, 63, 90, 115	0
1	F	444/487 (91%)	1.34	93 (20%) 1 0	43, 74, 100, 117	0
All	All	2684/2922 (91%)	0.45	222 (8%) 11 11	14, 35, 87, 117	0

The worst 5 of 222 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	50	GLY	6.5
1	F	78	GLY	5.8
1	E	108	PHE	5.8
1	F	434	THR	5.6
1	F	436	GLY	5.4

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

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

6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

6.4 Ligands [i](#)

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	MLI	A	1466	7/7	0.92	0.18	24,28,30,31	0
2	MLI	B	1465	7/7	0.93	0.17	29,30,36,40	0
2	MLI	C	1466	7/7	0.95	0.17	21,26,28,33	0
2	MLI	D	1465	7/7	0.95	0.15	25,31,35,36	0

6.5 Other polymers [i](#)

There are no such residues in this entry.