



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 18, 2026 – 03:13 AM UTC

PDB ID : 7XKG / pdb_00007xkg
Title : Crystal structure of an intramolecular mesacyl-CoA transferase from the 3-hydroxypropionic acid cycle of *Roseiflexus castenholzii*
Authors : Min, Z.Z.; Fan, C.P.; Wu, W.P.; Xin, Y.Y.; Liu, M.H.; Zhang, X.; Wang, Z.G.; Xu, X.L.
Deposited on : 2022-04-19
Resolution : 2.50 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
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

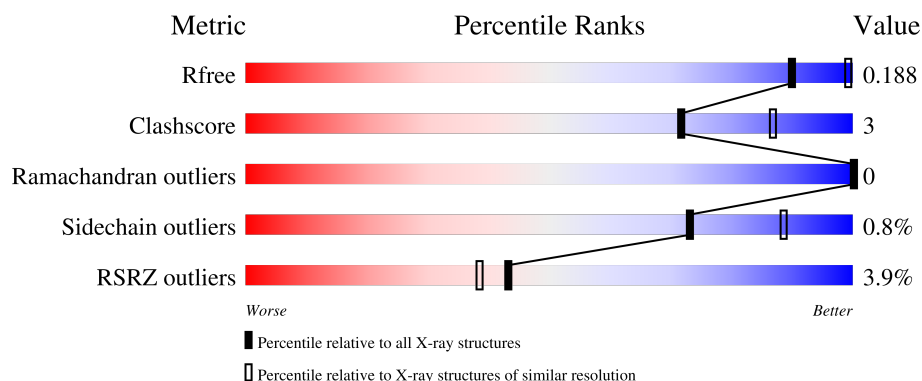
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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}	180053	5829 (2.50-2.50)
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)
RSRZ outliers	180081	5833 (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 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	429	3% 90% 5% 5%
1	B	429	3% 87% 8% 5%
1	C	429	4% 85% 10% 5%
1	D	429	6% 85% 9% 6%
1	E	429	3% 88% 7% 5%

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Mol	Chain	Length	Quality of chain
1	F	429	3% 88% 6% 5%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 20538 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 Acyl-CoA transferase/carnitine dehydratase-like protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	406	Total	C	N	O	S	0	0	0
			3134	1986	564	567	17			
1	B	406	Total	C	N	O	S	0	0	0
			3130	1984	564	565	17			
1	C	406	Total	C	N	O	S	0	0	0
			3134	1986	564	567	17			
1	D	405	Total	C	N	O	S	0	0	0
			3126	1981	563	566	16			
1	E	406	Total	C	N	O	S	0	0	0
			3134	1986	564	567	17			
1	F	406	Total	C	N	O	S	0	0	0
			3134	1986	564	567	17			

There are 126 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-20	MET	-	initiating methionine	UNP A7NHS9
A	-19	ARG	-	expression tag	UNP A7NHS9
A	-18	GLY	-	expression tag	UNP A7NHS9
A	-17	SER	-	expression tag	UNP A7NHS9
A	-16	HIS	-	expression tag	UNP A7NHS9
A	-15	HIS	-	expression tag	UNP A7NHS9
A	-14	HIS	-	expression tag	UNP A7NHS9
A	-13	HIS	-	expression tag	UNP A7NHS9
A	-12	HIS	-	expression tag	UNP A7NHS9
A	-11	HIS	-	expression tag	UNP A7NHS9
A	-10	GLY	-	expression tag	UNP A7NHS9
A	-9	MET	-	expression tag	UNP A7NHS9
A	-8	ALA	-	expression tag	UNP A7NHS9
A	-7	SER	-	expression tag	UNP A7NHS9
A	-6	GLU	-	expression tag	UNP A7NHS9
A	-5	LEU	-	expression tag	UNP A7NHS9
A	-4	ALA	-	expression tag	UNP A7NHS9

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Chain	Residue	Modelled	Actual	Comment	Reference
A	-3	LEU	-	expression tag	UNP A7NHS9
A	-2	GLU	-	expression tag	UNP A7NHS9
A	-1	ALA	-	expression tag	UNP A7NHS9
A	0	LEU	-	expression tag	UNP A7NHS9
B	-20	MET	-	initiating methionine	UNP A7NHS9
B	-19	ARG	-	expression tag	UNP A7NHS9
B	-18	GLY	-	expression tag	UNP A7NHS9
B	-17	SER	-	expression tag	UNP A7NHS9
B	-16	HIS	-	expression tag	UNP A7NHS9
B	-15	HIS	-	expression tag	UNP A7NHS9
B	-14	HIS	-	expression tag	UNP A7NHS9
B	-13	HIS	-	expression tag	UNP A7NHS9
B	-12	HIS	-	expression tag	UNP A7NHS9
B	-11	HIS	-	expression tag	UNP A7NHS9
B	-10	GLY	-	expression tag	UNP A7NHS9
B	-9	MET	-	expression tag	UNP A7NHS9
B	-8	ALA	-	expression tag	UNP A7NHS9
B	-7	SER	-	expression tag	UNP A7NHS9
B	-6	GLU	-	expression tag	UNP A7NHS9
B	-5	LEU	-	expression tag	UNP A7NHS9
B	-4	ALA	-	expression tag	UNP A7NHS9
B	-3	LEU	-	expression tag	UNP A7NHS9
B	-2	GLU	-	expression tag	UNP A7NHS9
B	-1	ALA	-	expression tag	UNP A7NHS9
B	0	LEU	-	expression tag	UNP A7NHS9
C	-20	MET	-	initiating methionine	UNP A7NHS9
C	-19	ARG	-	expression tag	UNP A7NHS9
C	-18	GLY	-	expression tag	UNP A7NHS9
C	-17	SER	-	expression tag	UNP A7NHS9
C	-16	HIS	-	expression tag	UNP A7NHS9
C	-15	HIS	-	expression tag	UNP A7NHS9
C	-14	HIS	-	expression tag	UNP A7NHS9
C	-13	HIS	-	expression tag	UNP A7NHS9
C	-12	HIS	-	expression tag	UNP A7NHS9
C	-11	HIS	-	expression tag	UNP A7NHS9
C	-10	GLY	-	expression tag	UNP A7NHS9
C	-9	MET	-	expression tag	UNP A7NHS9
C	-8	ALA	-	expression tag	UNP A7NHS9
C	-7	SER	-	expression tag	UNP A7NHS9
C	-6	GLU	-	expression tag	UNP A7NHS9
C	-5	LEU	-	expression tag	UNP A7NHS9
C	-4	ALA	-	expression tag	UNP A7NHS9

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Chain	Residue	Modelled	Actual	Comment	Reference
C	-3	LEU	-	expression tag	UNP A7NHS9
C	-2	GLU	-	expression tag	UNP A7NHS9
C	-1	ALA	-	expression tag	UNP A7NHS9
C	0	LEU	-	expression tag	UNP A7NHS9
D	-20	MET	-	initiating methionine	UNP A7NHS9
D	-19	ARG	-	expression tag	UNP A7NHS9
D	-18	GLY	-	expression tag	UNP A7NHS9
D	-17	SER	-	expression tag	UNP A7NHS9
D	-16	HIS	-	expression tag	UNP A7NHS9
D	-15	HIS	-	expression tag	UNP A7NHS9
D	-14	HIS	-	expression tag	UNP A7NHS9
D	-13	HIS	-	expression tag	UNP A7NHS9
D	-12	HIS	-	expression tag	UNP A7NHS9
D	-11	HIS	-	expression tag	UNP A7NHS9
D	-10	GLY	-	expression tag	UNP A7NHS9
D	-9	MET	-	expression tag	UNP A7NHS9
D	-8	ALA	-	expression tag	UNP A7NHS9
D	-7	SER	-	expression tag	UNP A7NHS9
D	-6	GLU	-	expression tag	UNP A7NHS9
D	-5	LEU	-	expression tag	UNP A7NHS9
D	-4	ALA	-	expression tag	UNP A7NHS9
D	-3	LEU	-	expression tag	UNP A7NHS9
D	-2	GLU	-	expression tag	UNP A7NHS9
D	-1	ALA	-	expression tag	UNP A7NHS9
D	0	LEU	-	expression tag	UNP A7NHS9
E	-20	MET	-	initiating methionine	UNP A7NHS9
E	-19	ARG	-	expression tag	UNP A7NHS9
E	-18	GLY	-	expression tag	UNP A7NHS9
E	-17	SER	-	expression tag	UNP A7NHS9
E	-16	HIS	-	expression tag	UNP A7NHS9
E	-15	HIS	-	expression tag	UNP A7NHS9
E	-14	HIS	-	expression tag	UNP A7NHS9
E	-13	HIS	-	expression tag	UNP A7NHS9
E	-12	HIS	-	expression tag	UNP A7NHS9
E	-11	HIS	-	expression tag	UNP A7NHS9
E	-10	GLY	-	expression tag	UNP A7NHS9
E	-9	MET	-	expression tag	UNP A7NHS9
E	-8	ALA	-	expression tag	UNP A7NHS9
E	-7	SER	-	expression tag	UNP A7NHS9
E	-6	GLU	-	expression tag	UNP A7NHS9
E	-5	LEU	-	expression tag	UNP A7NHS9
E	-4	ALA	-	expression tag	UNP A7NHS9

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Chain	Residue	Modelled	Actual	Comment	Reference
E	-3	LEU	-	expression tag	UNP A7NHS9
E	-2	GLU	-	expression tag	UNP A7NHS9
E	-1	ALA	-	expression tag	UNP A7NHS9
E	0	LEU	-	expression tag	UNP A7NHS9
F	-20	MET	-	initiating methionine	UNP A7NHS9
F	-19	ARG	-	expression tag	UNP A7NHS9
F	-18	GLY	-	expression tag	UNP A7NHS9
F	-17	SER	-	expression tag	UNP A7NHS9
F	-16	HIS	-	expression tag	UNP A7NHS9
F	-15	HIS	-	expression tag	UNP A7NHS9
F	-14	HIS	-	expression tag	UNP A7NHS9
F	-13	HIS	-	expression tag	UNP A7NHS9
F	-12	HIS	-	expression tag	UNP A7NHS9
F	-11	HIS	-	expression tag	UNP A7NHS9
F	-10	GLY	-	expression tag	UNP A7NHS9
F	-9	MET	-	expression tag	UNP A7NHS9
F	-8	ALA	-	expression tag	UNP A7NHS9
F	-7	SER	-	expression tag	UNP A7NHS9
F	-6	GLU	-	expression tag	UNP A7NHS9
F	-5	LEU	-	expression tag	UNP A7NHS9
F	-4	ALA	-	expression tag	UNP A7NHS9
F	-3	LEU	-	expression tag	UNP A7NHS9
F	-2	GLU	-	expression tag	UNP A7NHS9
F	-1	ALA	-	expression tag	UNP A7NHS9
F	0	LEU	-	expression tag	UNP A7NHS9

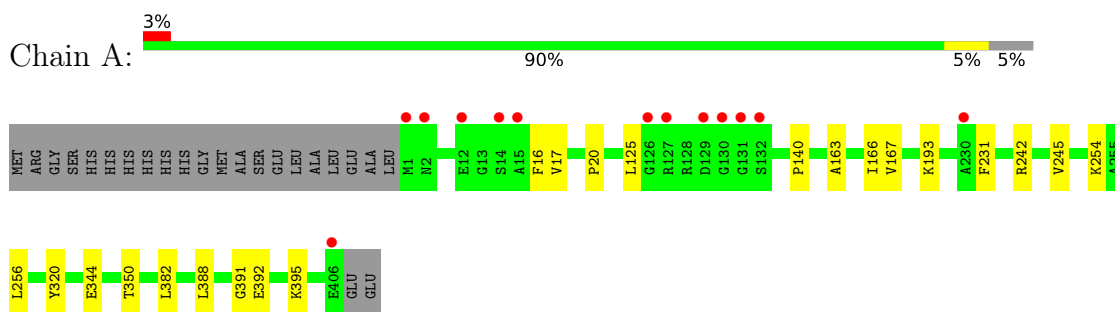
- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	303	Total O 303 303	0	0
2	B	305	Total O 305 305	0	0
2	C	295	Total O 295 295	0	0
2	D	285	Total O 285 285	0	0
2	E	278	Total O 278 278	0	0
2	F	280	Total O 280 280	0	0

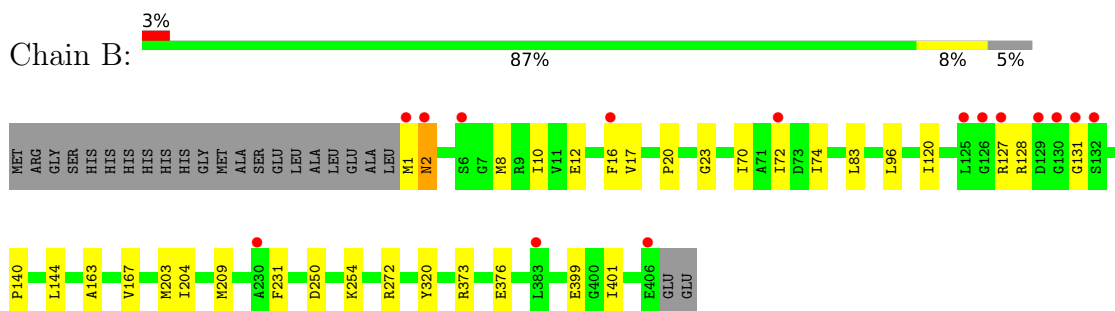
3 Residue-property plots [i](#)

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.

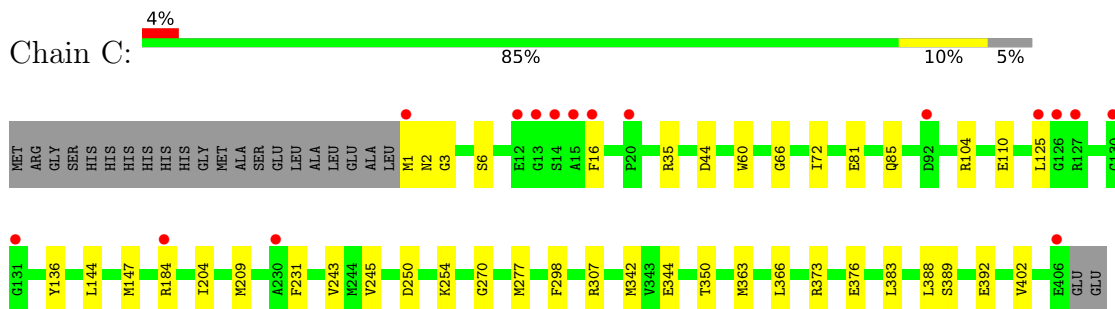
- Molecule 1: Acyl-CoA transferase/carnitine dehydratase-like protein



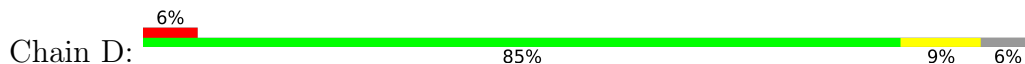
- Molecule 1: Acyl-CoA transferase/carnitine dehydratase-like protein

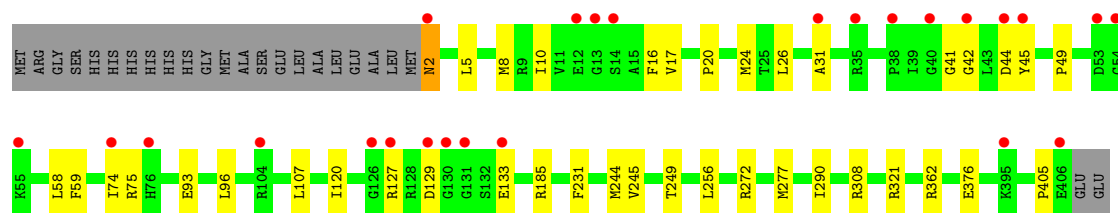


- Molecule 1: Acyl-CoA transferase/carnitine dehydratase-like protein

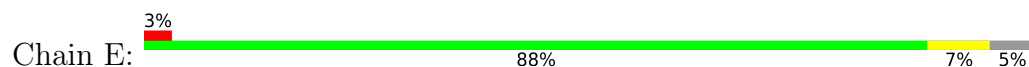


- Molecule 1: Acyl-CoA transferase/carnitine dehydratase-like protein

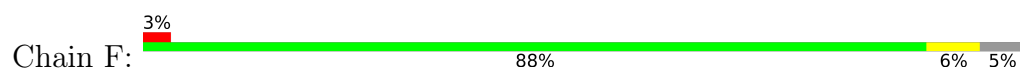




- Molecule 1: Acyl-CoA transferase/carnitine dehydratase-like protein



- Molecule 1: Acyl-CoA transferase/carnitine dehydratase-like protein



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	364.41Å 210.12Å 73.48Å 90.00° 95.81° 90.00°	Depositor
Resolution (Å)	44.42 – 2.50 44.42 – 2.50	Depositor EDS
% Data completeness (in resolution range)	99.2 (44.42-2.50) 99.3 (44.42-2.50)	Depositor EDS
R_{merge}	0.15	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.98 (at 2.51Å)	Xtriage
Refinement program	PHENIX 1.19.2_4158	Depositor
R, R_{free}	0.159 , 0.192 (Not available) , 0.188	Depositor DCC
R_{free} test set	9390 reflections (4.94%)	wwPDB-VP
Wilson B-factor (Å ²)	27.8	Xtriage
Anisotropy	0.317	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 38.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	20538	wwPDB-VP
Average B, all atoms (Å ²)	30.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.74% 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	0/3203	0.66	2/4338 (0.0%)
1	B	0.44	0/3199	0.65	2/4333 (0.0%)
1	C	0.42	0/3203	0.61	1/4338 (0.0%)
1	D	0.51	0/3195	0.72	3/4328 (0.1%)
1	E	0.51	0/3203	0.69	3/4338 (0.1%)
1	F	0.42	0/3203	0.61	0/4338
All	All	0.46	0/19206	0.66	11/26013 (0.0%)

There are no bond length outliers.

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	17	VAL	N-CA-C	11.32	121.98	111.45
1	E	17	VAL	N-CA-C	9.42	120.62	111.67
1	A	17	VAL	N-CA-C	9.25	120.05	111.45
1	B	231	PHE	N-CA-C	7.89	122.31	109.76
1	B	17	VAL	N-CA-C	7.79	119.06	111.67
1	C	231	PHE	N-CA-C	6.02	119.34	109.76
1	A	231	PHE	N-CA-C	5.86	119.46	110.20
1	E	131	GLY	N-CA-C	5.61	117.48	110.29
1	D	41	GLY	CA-C-O	-5.50	118.43	122.23
1	D	231	PHE	N-CA-C	5.20	118.09	109.46
1	E	336	GLU	N-CA-C	-5.03	105.79	111.28

There are no chirality outliers.

There are no planarity outliers.

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	3134	0	3136	14	0
1	B	3130	0	3132	27	0
1	C	3134	0	3136	30	0
1	D	3126	0	3124	28	0
1	E	3134	0	3136	19	0
1	F	3134	0	3136	20	0
2	A	303	0	0	0	0
2	B	305	0	0	3	0
2	C	295	0	0	5	0
2	D	285	0	0	1	0
2	E	278	0	0	0	0
2	F	280	0	0	0	0
All	All	20538	0	18800	122	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 3.

All (122) 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:277:MET:HE1	1:D:290:ILE:HD11	1.61	0.81
1:D:16:PHE:O	1:D:20:PRO:HG2	1.88	0.73
1:A:392:GLU:HA	1:A:395:LYS:HE2	1.73	0.71
1:D:376:GLU:HA	1:D:405:PRO:HG3	1.73	0.70
1:B:70:ILE:HG12	1:B:72:ILE:HG22	1.73	0.69
1:B:272:ARG:NH1	2:B:501:HOH:O	2.26	0.67
1:E:185:ARG:HH21	1:F:184:ARG:HH21	1.41	0.66
1:A:16:PHE:O	1:A:20:PRO:HG2	1.96	0.65
1:C:307:ARG:NH1	2:C:501:HOH:O	2.25	0.64
1:E:2:ASN:HB3	1:E:366:LEU:HD21	1.79	0.64
1:C:184:ARG:HH21	1:D:185:ARG:HH21	1.45	0.64
1:C:110:GLU:HG2	2:C:648:HOH:O	1.96	0.63
1:C:72:ILE:HG12	1:C:402:VAL:HG12	1.79	0.63
1:D:2:ASN:N	1:D:2:ASN:HD22	1.94	0.63
1:C:204:ILE:HB	1:C:209:MET:HE2	1.81	0.62
1:E:77:PRO:O	1:E:81:GLU:HG3	2.01	0.61
1:E:72:ILE:HG12	1:E:402:VAL:HG12	1.82	0.60
1:C:104:ARG:HG2	1:C:104:ARG:HH11	1.68	0.59
1:E:185:ARG:HH21	1:F:184:ARG:NH2	2.00	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:245:VAL:HG21	1:A:256:LEU:HD13	1.86	0.58
1:E:16:PHE:O	1:E:20:PRO:HG2	2.04	0.57
1:B:204:ILE:HB	1:B:209:MET:HE2	1.85	0.57
1:B:2:ASN:OD1	1:B:2:ASN:N	2.36	0.57
1:C:389:SER:OG	1:C:392:GLU:HG3	2.04	0.56
1:B:72:ILE:HG12	1:B:83:LEU:HD11	1.87	0.56
1:D:127:ARG:HB3	1:D:129:ASP:OD2	2.06	0.55
1:B:399:GLU:HB2	1:B:401:ILE:HD12	1.88	0.55
1:F:3:GLY:HA3	1:F:366:LEU:HD11	1.89	0.54
1:D:245:VAL:HG21	1:D:256:LEU:HD13	1.87	0.54
1:E:245:VAL:HG21	1:E:256:LEU:HD13	1.89	0.54
1:E:96:LEU:HD23	1:E:120:ILE:HB	1.88	0.54
1:C:1:MET:HG2	1:C:6:SER:HB3	1.90	0.54
1:C:344:GLU:HG2	1:C:350:THR:OG1	2.08	0.54
1:C:184:ARG:NH2	1:D:185:ARG:HH21	2.06	0.54
1:C:184:ARG:NH2	1:D:185:ARG:HE	2.06	0.53
1:C:81:GLU:O	1:C:85:GLN:HG3	2.09	0.53
1:B:250:ASP:HB3	1:B:254:LYS:NZ	2.24	0.53
1:C:147:MET:HE1	1:C:209:MET:HE3	1.90	0.53
1:A:166:ILE:HG23	1:B:203:MET:HE3	1.91	0.52
1:D:49:PRO:HD2	1:D:59:PHE:HB2	1.92	0.52
1:F:16:PHE:O	1:F:20:PRO:HG2	2.10	0.52
1:B:70:ILE:HG12	1:B:72:ILE:CG2	2.39	0.52
1:B:16:PHE:O	1:B:20:PRO:HG2	2.10	0.52
1:C:383:LEU:HD23	1:C:388:LEU:O	2.10	0.51
1:B:1:MET:HA	1:B:1:MET:HE2	1.92	0.51
1:B:373:ARG:HB2	1:B:376:GLU:HB2	1.93	0.51
1:B:8:MET:CE	1:B:10:ILE:HD11	2.42	0.50
1:A:242:ARG:HG2	1:A:320:TYR:CD1	2.47	0.50
1:B:128:ARG:HB2	2:B:783:HOH:O	2.11	0.49
1:C:250:ASP:O	1:C:254:LYS:HG3	2.12	0.49
1:E:166:ILE:HG23	1:F:203:MET:HE3	1.94	0.49
1:D:96:LEU:HD23	1:D:120:ILE:HB	1.95	0.47
1:D:107:LEU:HD23	1:D:107:LEU:HA	1.69	0.47
1:D:308:ARG:HG2	1:D:308:ARG:HH11	1.79	0.47
1:B:12:GLU:OE2	1:B:23:GLY:HA3	2.14	0.47
1:C:2:ASN:ND2	2:C:510:HOH:O	2.47	0.46
1:C:2:ASN:OD1	1:C:2:ASN:N	2.46	0.46
1:E:259:ALA:HB1	1:E:309:VAL:HG12	1.97	0.46
1:C:373:ARG:HB2	1:C:376:GLU:HB2	1.98	0.46
1:D:49:PRO:HB2	1:D:58:LEU:HD12	1.98	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:166:ILE:CG2	1:B:203:MET:HE3	2.46	0.46
1:B:127:ARG:HH21	1:B:131:GLY:HA3	1.80	0.45
1:E:363:MET:HE3	1:E:363:MET:HB3	1.87	0.45
1:F:12:GLU:HG2	1:F:14:SER:HB3	1.99	0.45
1:C:35:ARG:NH2	1:C:60:TRP:HE1	2.15	0.45
1:B:204:ILE:CB	1:B:209:MET:HE2	2.47	0.45
1:C:270:GLY:HA2	1:C:277:MET:HE3	1.97	0.45
1:F:194:LEU:HD12	1:F:194:LEU:HA	1.81	0.45
1:A:125:LEU:HD11	1:A:193:LYS:HG2	1.99	0.45
1:A:382:LEU:HB3	1:A:388:LEU:HD12	1.98	0.45
1:D:2:ASN:N	1:D:2:ASN:ND2	2.64	0.44
1:E:362:ARG:HG2	1:F:190:GLN:HB3	2.00	0.44
1:F:245:VAL:HG21	1:F:256:LEU:HD13	1.99	0.44
1:A:140:PRO:HB3	1:B:320:TYR:CD2	2.52	0.44
1:F:96:LEU:HD23	1:F:120:ILE:HB	2.00	0.44
1:F:86:LEU:HD23	1:F:86:LEU:HA	1.82	0.43
1:C:136:TYR:O	1:D:244:MET:HG2	2.18	0.43
1:B:272:ARG:NH2	2:B:508:HOH:O	2.45	0.43
1:B:250:ASP:HB3	1:B:254:LYS:HZ2	1.82	0.43
1:D:308:ARG:HG2	1:D:308:ARG:NH1	2.34	0.43
1:C:16:PHE:HZ	1:D:249:THR:HG22	1.83	0.43
1:E:2:ASN:HB3	1:E:366:LEU:CD2	2.48	0.43
1:A:163:ALA:O	1:A:167:VAL:HG23	2.19	0.42
1:B:144:LEU:HD13	1:B:144:LEU:HA	1.89	0.42
1:F:351:TYR:CZ	1:F:353:MET:HE2	2.53	0.42
1:D:24:MET:HE2	1:D:24:MET:HB3	1.81	0.42
1:C:184:ARG:CZ	1:D:185:ARG:HE	2.32	0.42
1:D:16:PHE:H	1:D:44:ASP:HB3	1.85	0.42
1:C:16:PHE:H	1:C:44:ASP:HB3	1.84	0.42
1:C:245:VAL:HG11	1:C:298:PHE:HZ	1.85	0.42
1:B:96:LEU:HD23	1:B:120:ILE:HB	2.02	0.42
1:E:140:PRO:HB3	1:F:320:TYR:CD2	2.55	0.42
1:E:320:TYR:CD2	1:F:140:PRO:HB3	2.55	0.42
1:F:163:ALA:O	1:F:167:VAL:HG23	2.19	0.42
1:F:83:LEU:HD23	1:F:83:LEU:HA	1.87	0.42
1:C:363:MET:HE3	1:C:363:MET:HB3	1.95	0.42
2:C:651:HOH:O	1:D:321:ARG:HD2	2.19	0.42
1:F:74:ILE:H	1:F:74:ILE:HG13	1.61	0.41
1:D:10:ILE:HD13	1:D:26:LEU:HD13	2.01	0.41
1:C:147:MET:HE1	1:C:209:MET:CE	2.50	0.41
1:D:5:LEU:HB3	1:D:31:ALA:HB2	2.01	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:8:MET:HA	1:D:93:GLU:HB3	2.02	0.41
1:D:42:GLY:H	1:D:45:TYR:HB2	1.86	0.41
1:A:320:TYR:CD2	1:B:140:PRO:HB3	2.56	0.41
1:A:391:GLY:O	1:A:395:LYS:HG3	2.21	0.41
1:F:277:MET:HE2	1:F:277:MET:HA	2.03	0.41
1:A:344:GLU:HG2	1:A:350:THR:OG1	2.20	0.41
1:B:83:LEU:HD23	1:B:83:LEU:HA	1.90	0.41
1:E:70:ILE:HG12	1:E:72:ILE:HG13	2.03	0.41
1:E:83:LEU:HD23	1:E:83:LEU:HA	1.90	0.41
1:E:140:PRO:HA	1:E:145:PRO:HG2	2.02	0.41
1:D:127:ARG:HG3	1:D:133:GLU:OE2	2.20	0.41
1:F:305:GLU:OE2	1:F:308:ARG:NH2	2.45	0.41
1:A:254:LYS:HE2	1:A:254:LYS:HB2	1.88	0.40
1:B:72:ILE:HD11	1:B:74:ILE:HG22	2.03	0.40
1:B:163:ALA:O	1:B:167:VAL:HG23	2.21	0.40
1:C:66:GLY:HA3	2:C:719:HOH:O	2.21	0.40
1:C:342:MET:HE3	1:C:350:THR:HG22	2.03	0.40
1:E:163:ALA:O	1:E:167:VAL:HG23	2.21	0.40
1:D:75:ARG:HD2	2:D:501:HOH:O	2.20	0.40
1:F:184:ARG:HB2	1:F:184:ARG:NH1	2.36	0.40
1:C:3:GLY:HA3	1:C:366:LEU:HD11	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	404/429 (94%)	398 (98%)	6 (2%)	0	100	100
1	B	404/429 (94%)	397 (98%)	7 (2%)	0	100	100
1	C	404/429 (94%)	398 (98%)	6 (2%)	0	100	100
1	D	403/429 (94%)	397 (98%)	6 (2%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	E	404/429 (94%)	395 (98%)	9 (2%)	0	100	100
1	F	404/429 (94%)	398 (98%)	6 (2%)	0	100	100
All	All	2423/2574 (94%)	2383 (98%)	40 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	250/339 (74%)	250 (100%)	0	100	100
1	B	29/339 (9%)	28 (97%)	1 (3%)	32	60
1	C	273/339 (80%)	270 (99%)	3 (1%)	65	84
1	D	320/339 (94%)	316 (99%)	4 (1%)	61	82
1	E	321/339 (95%)	318 (99%)	3 (1%)	70	87
1	F	321/339 (95%)	320 (100%)	1 (0%)	86	94
All	All	1514/2034 (74%)	1502 (99%)	12 (1%)	73	88

All (12) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	2	ASN
1	C	125	LEU
1	C	144	LEU
1	C	243	VAL
1	D	2	ASN
1	D	74	ILE
1	D	272	ARG
1	D	362	ARG
1	E	72	ILE
1	E	127	ARG
1	E	128	ARG
1	F	144	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	158	ASN
1	A	282	ASN
1	C	170	GLN
1	E	123	ASN
1	E	313	HIS
1	F	158	ASN
1	F	170	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 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 ⓘ

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	406/429 (94%)	-0.33	13 (3%)	50	46	19, 26, 40, 77	0
1	B	406/429 (94%)	-0.34	15 (3%)	45	40	19, 26, 43, 83	0
1	C	406/429 (94%)	-0.27	16 (3%)	43	38	19, 28, 45, 77	0
1	D	405/429 (94%)	-0.06	25 (6%)	26	23	19, 28, 44, 76	0
1	E	406/429 (94%)	-0.19	14 (3%)	48	43	20, 29, 49, 94	0
1	F	406/429 (94%)	-0.23	12 (2%)	52	48	20, 29, 51, 85	0
All	All	2435/2574 (94%)	-0.24	95 (3%)	43	38	19, 27, 46, 94	0

All (95) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	130	GLY	8.2
1	E	131	GLY	8.0
1	E	127	ARG	6.9
1	B	131	GLY	6.8
1	E	1	MET	6.6
1	A	129	ASP	6.1
1	A	127	ARG	5.9
1	D	131	GLY	5.8
1	E	2	ASN	5.5
1	A	130	GLY	5.4
1	B	126	GLY	5.4
1	C	1	MET	5.1
1	F	76	HIS	5.1
1	A	132	SER	5.0
1	B	16	PHE	5.0
1	F	1	MET	5.0
1	F	130	GLY	4.9
1	B	127	ARG	4.8
1	A	131	GLY	4.7

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Mol	Chain	Res	Type	RSRZ
1	D	40	GLY	4.7
1	D	127	ARG	4.7
1	C	130	GLY	4.5
1	C	131	GLY	4.4
1	B	130	GLY	4.2
1	C	12	GLU	4.1
1	A	1	MET	4.1
1	D	2	ASN	4.0
1	A	406	GLU	4.0
1	D	42	GLY	4.0
1	D	130	GLY	4.0
1	F	14	SER	3.9
1	F	406	GLU	3.9
1	D	126	GLY	3.9
1	B	129	ASP	3.9
1	C	406	GLU	3.8
1	E	129	ASP	3.8
1	C	15	ALA	3.8
1	C	14	SER	3.7
1	A	126	GLY	3.7
1	D	12	GLU	3.6
1	F	127	ARG	3.6
1	C	16	PHE	3.6
1	F	129	ASP	3.6
1	F	131	GLY	3.6
1	D	14	SER	3.5
1	E	406	GLU	3.4
1	D	31	ALA	3.4
1	B	1	MET	3.4
1	D	129	ASP	3.3
1	E	126	GLY	3.3
1	B	406	GLU	3.3
1	C	127	ARG	3.1
1	D	54	GLY	3.0
1	E	133	GLU	3.0
1	E	128	ARG	3.0
1	B	2	ASN	2.9
1	A	12	GLU	2.9
1	B	6	SER	2.9
1	C	126	GLY	2.9
1	C	20	PRO	2.9
1	D	44	ASP	2.9

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Mol	Chain	Res	Type	RSRZ
1	D	53	ASP	2.9
1	E	132	SER	2.8
1	E	76	HIS	2.8
1	B	72	ILE	2.8
1	D	45	TYR	2.8
1	A	14	SER	2.6
1	D	38	PRO	2.6
1	B	230	ALA	2.6
1	B	132	SER	2.5
1	E	185	ARG	2.5
1	B	125	LEU	2.5
1	D	35	ARG	2.5
1	D	133	GLU	2.4
1	F	92	ASP	2.4
1	A	2	ASN	2.4
1	D	104	ARG	2.4
1	F	12	GLU	2.3
1	F	2	ASN	2.3
1	D	76	HIS	2.3
1	B	383	LEU	2.3
1	C	13	GLY	2.2
1	F	15	ALA	2.2
1	C	92	ASP	2.2
1	C	125	LEU	2.1
1	D	13	GLY	2.1
1	D	406	GLU	2.1
1	D	395	LYS	2.1
1	C	230	ALA	2.1
1	E	314	ARG	2.1
1	D	55	LYS	2.1
1	A	230	ALA	2.1
1	D	74	ILE	2.1
1	A	15	ALA	2.0
1	C	184	ARG	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 [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.