



# Full wwPDB X-ray Structure Validation Report ⓘ

May 21, 2020 – 06:34 pm BST

PDB ID : 6ACQ  
Title : Crystal structure of (S)-3-hydroxybutyryl-CoA dehydrogenase from *Clostridium acetobutylicum*, apo form  
Authors : Takenoya, M.; Taguchi, S.; Yajima, S.  
Deposited on : 2018-07-27  
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](mailto: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.

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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

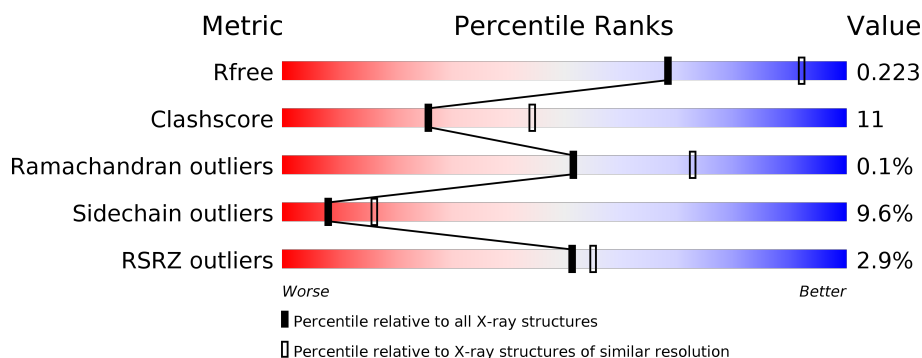
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.50 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  $\geq 3$ , 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	302	<div> <div>76%</div> <div>15%</div> <div>7%</div> </div>
1	B	302	<div>2%</div> <div>71%</div> <div>22%</div> <div>7%</div>
1	C	302	<div>%</div> <div>73%</div> <div>19%</div> <div>7%</div>
1	D	302	<div>13%</div> <div>64%</div> <div>25%</div> <div>7%</div>
1	E	302	<div>67%</div> <div>22%</div> <div>7%</div>
1	F	302	<div>78%</div> <div>17%</div> <div>5%</div>

## 2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 12991 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 3-hydroxybutyryl-CoA dehydrogenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	281	Total	C	N	O	S	0	0	0
			2132	1359	355	404	14			
1	B	289	Total	C	N	O	S	0	0	0
			2189	1394	368	413	14			
1	C	281	Total	C	N	O	S	0	0	0
			2132	1359	355	404	14			
1	D	281	Total	C	N	O	S	0	0	0
			2132	1359	355	404	14			
1	E	281	Total	C	N	O	S	0	0	0
			2132	1359	355	404	14			
1	F	290	Total	C	N	O	S	0	0	0
			2195	1397	369	415	14			

There are 120 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-19	MET	-	expression tag	UNP P52041
A	-18	GLY	-	expression tag	UNP P52041
A	-17	SER	-	expression tag	UNP P52041
A	-16	SER	-	expression tag	UNP P52041
A	-15	HIS	-	expression tag	UNP P52041
A	-14	HIS	-	expression tag	UNP P52041
A	-13	HIS	-	expression tag	UNP P52041
A	-12	HIS	-	expression tag	UNP P52041
A	-11	HIS	-	expression tag	UNP P52041
A	-10	HIS	-	expression tag	UNP P52041
A	-9	SER	-	expression tag	UNP P52041
A	-8	SER	-	expression tag	UNP P52041
A	-7	GLY	-	expression tag	UNP P52041
A	-6	LEU	-	expression tag	UNP P52041
A	-5	VAL	-	expression tag	UNP P52041
A	-4	PRO	-	expression tag	UNP P52041
A	-3	ARG	-	expression tag	UNP P52041

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Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	GLY	-	expression tag	UNP P52041
A	-1	SER	-	expression tag	UNP P52041
A	0	HIS	-	expression tag	UNP P52041
B	-19	MET	-	expression tag	UNP P52041
B	-18	GLY	-	expression tag	UNP P52041
B	-17	SER	-	expression tag	UNP P52041
B	-16	SER	-	expression tag	UNP P52041
B	-15	HIS	-	expression tag	UNP P52041
B	-14	HIS	-	expression tag	UNP P52041
B	-13	HIS	-	expression tag	UNP P52041
B	-12	HIS	-	expression tag	UNP P52041
B	-11	HIS	-	expression tag	UNP P52041
B	-10	HIS	-	expression tag	UNP P52041
B	-9	SER	-	expression tag	UNP P52041
B	-8	SER	-	expression tag	UNP P52041
B	-7	GLY	-	expression tag	UNP P52041
B	-6	LEU	-	expression tag	UNP P52041
B	-5	VAL	-	expression tag	UNP P52041
B	-4	PRO	-	expression tag	UNP P52041
B	-3	ARG	-	expression tag	UNP P52041
B	-2	GLY	-	expression tag	UNP P52041
B	-1	SER	-	expression tag	UNP P52041
B	0	HIS	-	expression tag	UNP P52041
C	-19	MET	-	expression tag	UNP P52041
C	-18	GLY	-	expression tag	UNP P52041
C	-17	SER	-	expression tag	UNP P52041
C	-16	SER	-	expression tag	UNP P52041
C	-15	HIS	-	expression tag	UNP P52041
C	-14	HIS	-	expression tag	UNP P52041
C	-13	HIS	-	expression tag	UNP P52041
C	-12	HIS	-	expression tag	UNP P52041
C	-11	HIS	-	expression tag	UNP P52041
C	-10	HIS	-	expression tag	UNP P52041
C	-9	SER	-	expression tag	UNP P52041
C	-8	SER	-	expression tag	UNP P52041
C	-7	GLY	-	expression tag	UNP P52041
C	-6	LEU	-	expression tag	UNP P52041
C	-5	VAL	-	expression tag	UNP P52041
C	-4	PRO	-	expression tag	UNP P52041
C	-3	ARG	-	expression tag	UNP P52041
C	-2	GLY	-	expression tag	UNP P52041
C	-1	SER	-	expression tag	UNP P52041

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Chain	Residue	Modelled	Actual	Comment	Reference
C	0	HIS	-	expression tag	UNP P52041
D	-19	MET	-	expression tag	UNP P52041
D	-18	GLY	-	expression tag	UNP P52041
D	-17	SER	-	expression tag	UNP P52041
D	-16	SER	-	expression tag	UNP P52041
D	-15	HIS	-	expression tag	UNP P52041
D	-14	HIS	-	expression tag	UNP P52041
D	-13	HIS	-	expression tag	UNP P52041
D	-12	HIS	-	expression tag	UNP P52041
D	-11	HIS	-	expression tag	UNP P52041
D	-10	HIS	-	expression tag	UNP P52041
D	-9	SER	-	expression tag	UNP P52041
D	-8	SER	-	expression tag	UNP P52041
D	-7	GLY	-	expression tag	UNP P52041
D	-6	LEU	-	expression tag	UNP P52041
D	-5	VAL	-	expression tag	UNP P52041
D	-4	PRO	-	expression tag	UNP P52041
D	-3	ARG	-	expression tag	UNP P52041
D	-2	GLY	-	expression tag	UNP P52041
D	-1	SER	-	expression tag	UNP P52041
D	0	HIS	-	expression tag	UNP P52041
E	-19	MET	-	expression tag	UNP P52041
E	-18	GLY	-	expression tag	UNP P52041
E	-17	SER	-	expression tag	UNP P52041
E	-16	SER	-	expression tag	UNP P52041
E	-15	HIS	-	expression tag	UNP P52041
E	-14	HIS	-	expression tag	UNP P52041
E	-13	HIS	-	expression tag	UNP P52041
E	-12	HIS	-	expression tag	UNP P52041
E	-11	HIS	-	expression tag	UNP P52041
E	-10	HIS	-	expression tag	UNP P52041
E	-9	SER	-	expression tag	UNP P52041
E	-8	SER	-	expression tag	UNP P52041
E	-7	GLY	-	expression tag	UNP P52041
E	-6	LEU	-	expression tag	UNP P52041
E	-5	VAL	-	expression tag	UNP P52041
E	-4	PRO	-	expression tag	UNP P52041
E	-3	ARG	-	expression tag	UNP P52041
E	-2	GLY	-	expression tag	UNP P52041
E	-1	SER	-	expression tag	UNP P52041
E	0	HIS	-	expression tag	UNP P52041
F	-19	MET	-	expression tag	UNP P52041

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Chain	Residue	Modelled	Actual	Comment	Reference
F	-18	GLY	-	expression tag	UNP P52041
F	-17	SER	-	expression tag	UNP P52041
F	-16	SER	-	expression tag	UNP P52041
F	-15	HIS	-	expression tag	UNP P52041
F	-14	HIS	-	expression tag	UNP P52041
F	-13	HIS	-	expression tag	UNP P52041
F	-12	HIS	-	expression tag	UNP P52041
F	-11	HIS	-	expression tag	UNP P52041
F	-10	HIS	-	expression tag	UNP P52041
F	-9	SER	-	expression tag	UNP P52041
F	-8	SER	-	expression tag	UNP P52041
F	-7	GLY	-	expression tag	UNP P52041
F	-6	LEU	-	expression tag	UNP P52041
F	-5	VAL	-	expression tag	UNP P52041
F	-4	PRO	-	expression tag	UNP P52041
F	-3	ARG	-	expression tag	UNP P52041
F	-2	GLY	-	expression tag	UNP P52041
F	-1	SER	-	expression tag	UNP P52041
F	0	HIS	-	expression tag	UNP P52041

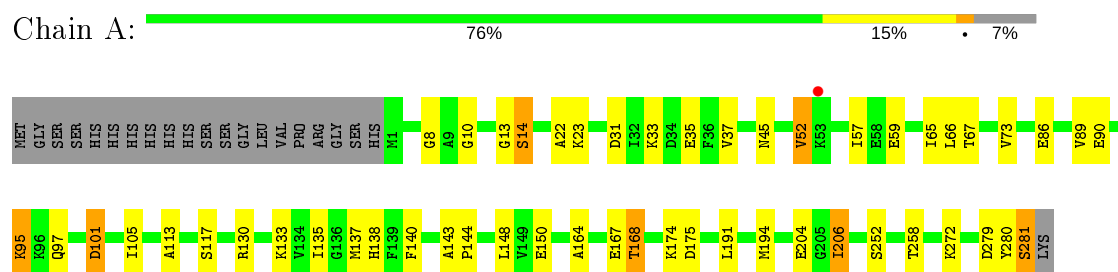
- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	23	Total O 23 23	0	0
2	B	8	Total O 8 8	0	0
2	C	16	Total O 16 16	0	0
2	D	4	Total O 4 4	0	0
2	E	8	Total O 8 8	0	0
2	F	20	Total O 20 20	0	0

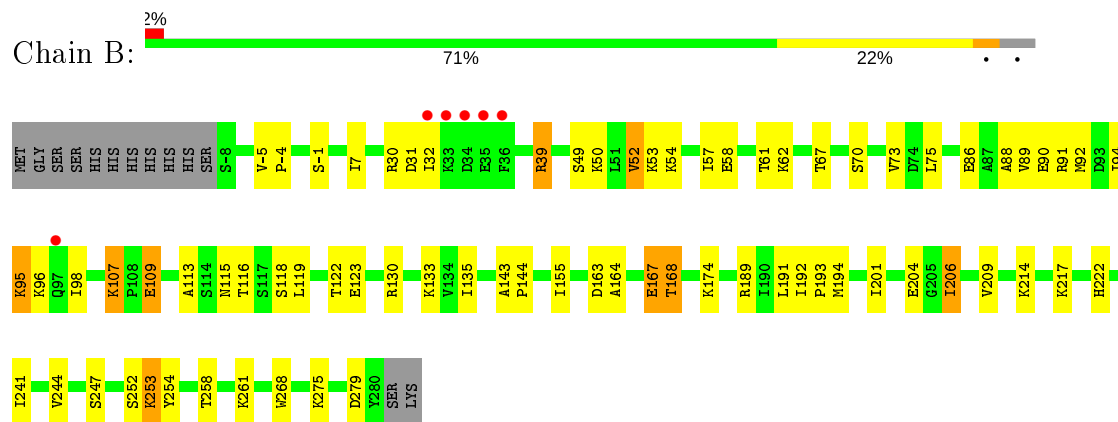
### 3 Residue-property plots [i](#)

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.

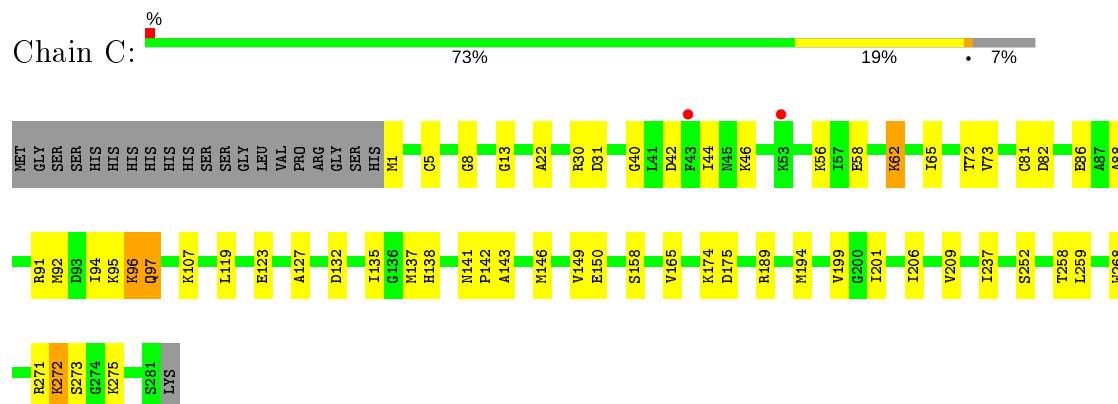
#### • Molecule 1: 3-hydroxybutyryl-CoA dehydrogenase



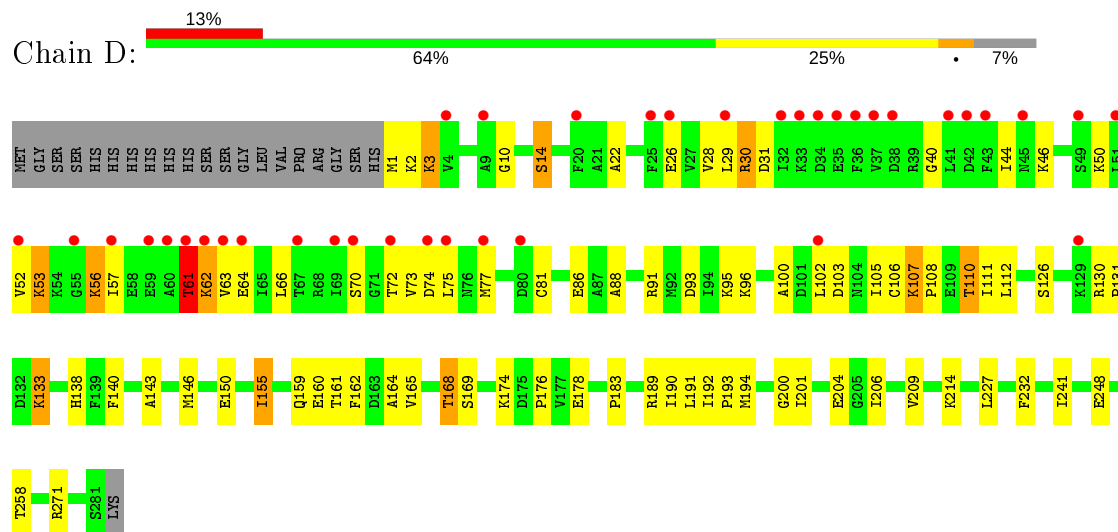
#### • Molecule 1: 3-hydroxybutyryl-CoA dehydrogenase



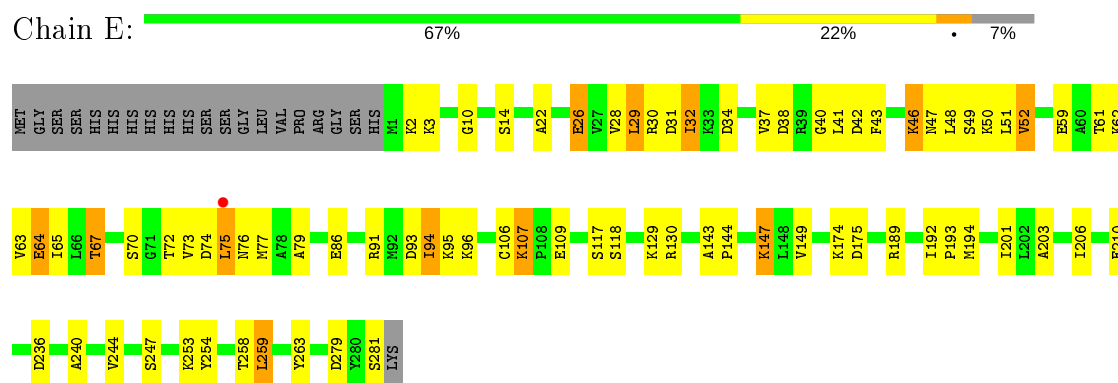
#### • Molecule 1: 3-hydroxybutyryl-CoA dehydrogenase



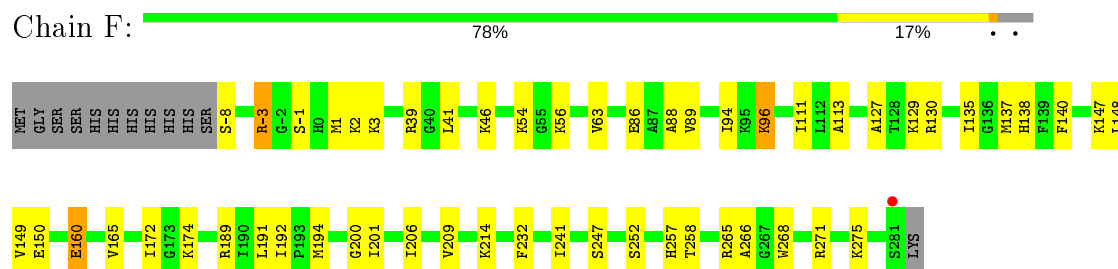
- Molecule 1: 3-hydroxybutyryl-CoA dehydrogenase



- Molecule 1: 3-hydroxybutyryl-CoA dehydrogenase



- Molecule 1: 3-hydroxybutyryl-CoA dehydrogenase



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	133.43Å 134.48Å 153.80Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	47.90 – 2.50 47.86 – 2.50	Depositor EDS
% Data completeness (in resolution range)	99.6 (47.90-2.50) 99.6 (47.86-2.50)	Depositor EDS
$R_{merge}$	0.09	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.39 (at 2.51Å)	Xtriage
Refinement program	REFMAC 5.8.0218	Depositor
R, $R_{free}$	0.199 , 0.222 0.203 , 0.223	Depositor DCC
$R_{free}$ test set	4780 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	48.4	Xtriage
Anisotropy	0.430	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.33 , 33.0	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.010 for k,h,-l	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	12991	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	54.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.18% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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	0.57	0/2162	0.71	0/2914
1	B	0.52	0/2221	0.69	0/2994
1	C	0.53	0/2162	0.69	0/2914
1	D	0.48	0/2162	0.69	0/2914
1	E	0.49	0/2162	0.69	0/2914
1	F	0.57	0/2227	0.70	0/3002
All	All	0.53	0/13096	0.69	0/17652

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	D	0	1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	D	61	THR	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	2132	0	2209	37	0
1	B	2189	0	2267	43	0
1	C	2132	0	2209	37	0
1	D	2132	0	2209	83	0
1	E	2132	0	2209	67	0
1	F	2195	0	2272	32	0
2	A	23	0	0	0	0
2	B	8	0	0	0	0
2	C	16	0	0	1	0
2	D	4	0	0	0	0
2	E	8	0	0	0	0
2	F	20	0	0	2	0
All	All	12991	0	13375	277	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 11.

All (277) 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:B:194:MET:CE	1:C:194:MET:SD	2.24	1.25
1:B:194:MET:HE2	1:C:194:MET:SD	1.79	1.19
1:D:194:MET:SD	1:E:194:MET:CE	2.34	1.16
1:D:102:LEU:CD2	1:D:112:LEU:HD22	1.82	1.10
1:D:102:LEU:HD23	1:D:112:LEU:HD22	1.35	1.05
1:D:194:MET:SD	1:E:194:MET:HE1	1.96	1.02
1:C:86:GLU:OE1	1:C:95:LYS:HE3	1.61	1.01
1:E:29:LEU:HD12	1:E:30:ARG:N	1.77	0.99
1:D:30:ARG:HD2	1:D:30:ARG:C	1.81	0.98
1:B:194:MET:HE1	1:C:194:MET:SD	2.05	0.96
1:E:46:LYS:HE2	1:E:50:LYS:HE3	1.45	0.95
1:E:29:LEU:HD12	1:E:29:LEU:C	1.85	0.94
1:D:30:ARG:HD2	1:D:31:ASP:N	1.82	0.93
1:D:102:LEU:CD2	1:D:112:LEU:CD2	2.47	0.92
1:E:41:LEU:H	1:E:41:LEU:HD12	1.38	0.89
1:D:86:GLU:OE1	1:D:95:LYS:NZ	2.06	0.87
1:D:102:LEU:HD21	1:D:112:LEU:CD2	2.05	0.86
1:A:89:VAL:O	1:A:95:LYS:NZ	2.08	0.86
1:D:10:GLY:O	1:D:14:SER:OG	1.94	0.85
1:C:40:GLY:O	1:C:44:ILE:HG13	1.79	0.83
1:E:86:GLU:OE2	1:E:95:LYS:HE3	1.80	0.81
1:E:37:VAL:O	1:E:41:LEU:HD12	1.81	0.80

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:164:ALA:O	1:A:168:THR:CG2	2.30	0.79
1:D:53:LYS:HG3	1:D:53:LYS:O	1.83	0.78
1:B:143:ALA:O	1:B:174:LYS:NZ	2.18	0.77
1:D:86:GLU:OE2	1:D:88:ALA:HB3	1.86	0.75
1:D:232:PHE:HD1	1:D:271:ARG:NH1	1.85	0.75
1:A:86:GLU:OE1	1:A:95:LYS:HE3	1.88	0.73
1:D:81:CYS:O	1:D:110:THR:HG23	1.87	0.73
1:A:86:GLU:CD	1:A:95:LYS:HE3	2.08	0.72
1:E:10:GLY:O	1:E:14:SER:OG	2.08	0.72
1:D:232:PHE:CD1	1:D:271:ARG:NH1	2.58	0.72
1:A:164:ALA:O	1:A:168:THR:HG22	1.90	0.71
1:F:-1:SER:O	1:F:2:LYS:NZ	2.24	0.70
1:D:30:ARG:CD	1:D:31:ASP:N	2.54	0.70
1:D:81:CYS:O	1:D:110:THR:CG2	2.39	0.70
1:B:95:LYS:HE2	1:B:119:LEU:HD12	1.72	0.70
1:D:107:LYS:O	1:D:110:THR:OG1	2.09	0.70
1:B:107:LYS:HD2	1:B:109:GLU:HG2	1.75	0.69
1:F:-3:ARG:HH11	1:F:-3:ARG:HG2	1.58	0.69
1:D:164:ALA:O	1:D:168:THR:HG22	1.92	0.69
1:E:29:LEU:C	1:E:29:LEU:CD1	2.59	0.69
1:D:194:MET:SD	1:E:194:MET:HE2	2.31	0.68
1:E:41:LEU:N	1:E:41:LEU:HD12	2.07	0.68
1:B:163:ASP:O	1:B:167:GLU:HG3	1.93	0.68
1:D:102:LEU:HD12	1:D:106:CYS:SG	2.34	0.68
1:E:37:VAL:O	1:E:41:LEU:CD1	2.41	0.68
1:C:8:GLY:O	1:C:13:GLY:HA3	1.94	0.68
1:A:22:ALA:HA	1:A:65:ILE:HG12	1.77	0.67
1:D:73:VAL:O	1:D:73:VAL:HG23	1.92	0.67
1:E:143:ALA:O	1:E:174:LYS:NZ	2.24	0.66
1:D:102:LEU:CD1	1:D:106:CYS:SG	2.84	0.66
1:D:232:PHE:HD1	1:D:271:ARG:HH11	1.43	0.65
1:C:86:GLU:OE2	1:C:88:ALA:N	2.27	0.65
1:E:203:ALA:HB2	1:E:259:LEU:HD12	1.79	0.65
1:A:52:VAL:HG12	1:A:57:ILE:O	1.97	0.64
1:B:217:LYS:HG2	1:B:222:HIS:O	1.97	0.64
1:E:72:THR:HG22	1:E:73:VAL:N	2.12	0.64
1:A:52:VAL:HG11	1:A:59:GLU:HA	1.80	0.63
1:B:194:MET:HE2	1:C:194:MET:CG	2.27	0.63
1:D:3:LYS:HD2	1:D:81:CYS:SG	2.38	0.63
1:A:59:GLU:HG3	1:A:59:GLU:O	1.99	0.63
1:B:49:SER:OG	1:B:62:LYS:NZ	2.31	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:201:ILE:HG23	1:E:206:ILE:HB	1.80	0.62
1:D:52:VAL:HG12	1:D:57:ILE:HG13	1.80	0.61
1:A:194:MET:SD	1:F:194:MET:SD	2.98	0.61
1:F:232:PHE:CD1	1:F:271:ARG:NH1	2.68	0.61
1:D:130:ARG:N	1:D:131:PRO:CD	2.64	0.61
1:D:159:GLN:O	1:D:159:GLN:HG3	1.99	0.61
1:D:194:MET:SD	1:E:194:MET:HE3	2.38	0.61
1:A:164:ALA:O	1:A:168:THR:HG23	2.00	0.61
1:E:59:GLU:O	1:E:63:VAL:HG23	2.01	0.61
1:E:63:VAL:O	1:E:67:THR:HG23	2.01	0.61
1:D:30:ARG:CD	1:D:30:ARG:C	2.64	0.60
1:E:3:LYS:HG3	1:E:26:GLU:HG2	1.82	0.60
1:D:100:ALA:O	1:D:103:ASP:HB3	2.02	0.60
1:D:30:ARG:HD2	1:D:31:ASP:CA	2.32	0.60
1:E:107:LYS:HD2	1:E:109:GLU:HG2	1.84	0.59
1:F:1:MET:HE3	1:F:111:ILE:HD12	1.83	0.59
1:F:1:MET:CE	1:F:111:ILE:HD12	2.32	0.59
1:B:122:THR:HG23	1:B:155:ILE:HG22	1.82	0.59
1:D:201:ILE:HG23	1:D:206:ILE:HB	1.84	0.59
1:D:111:ILE:HD13	1:D:161:THR:HG23	1.84	0.59
1:A:143:ALA:O	1:A:174:LYS:NZ	2.30	0.59
1:A:280:TYR:O	1:A:281:SER:C	2.41	0.58
1:E:49:SER:HA	1:E:52:VAL:HG23	1.84	0.58
1:E:91:ARG:HD2	1:E:94:ILE:HD13	1.86	0.58
1:D:165:VAL:HA	1:D:168:THR:HG23	1.85	0.58
1:E:86:GLU:CD	1:E:95:LYS:HE3	2.24	0.58
1:E:72:THR:HG22	1:E:74:ASP:N	2.19	0.58
1:D:169:SER:HB2	1:D:176:PRO:HG3	1.86	0.57
1:A:86:GLU:OE2	1:A:95:LYS:HE2	2.05	0.57
1:B:164:ALA:O	1:B:168:THR:HG23	2.03	0.57
1:D:214:LYS:HE3	1:E:175:ASP:OD2	2.05	0.56
1:C:272:LYS:HG2	2:C:308:HOH:O	2.05	0.56
1:B:90:GLU:OE1	1:B:116:THR:OG1	2.16	0.55
1:D:61:THR:HA	1:D:64:GLU:HB3	1.89	0.55
1:D:73:VAL:O	1:D:73:VAL:CG2	2.54	0.55
1:E:46:LYS:HE2	1:E:50:LYS:CE	2.29	0.55
1:E:259:LEU:HD21	1:E:263:TYR:HE1	1.72	0.54
1:C:95:LYS:HD2	1:C:119:LEU:HD12	1.89	0.54
1:C:95:LYS:HD2	1:C:119:LEU:CD1	2.37	0.54
1:A:8:GLY:O	1:A:13:GLY:HA3	2.08	0.54
1:A:86:GLU:OE2	1:A:95:LYS:CE	2.56	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:109:GLU:HA	1:B:133:LYS:HE2	1.91	0.53
1:E:72:THR:HG22	1:E:74:ASP:H	1.74	0.53
1:C:135:ILE:HG21	1:C:165:VAL:HG21	1.89	0.53
1:E:40:GLY:O	1:E:43:PHE:HB3	2.07	0.53
1:E:49:SER:CA	1:E:52:VAL:HG23	2.39	0.53
1:D:91:ARG:HG2	1:D:93:ASP:OD1	2.09	0.53
1:D:143:ALA:O	1:D:174:LYS:NZ	2.33	0.53
1:E:48:LEU:HA	1:E:51:LEU:HD12	1.90	0.53
1:B:113:ALA:HA	1:B:135:ILE:O	2.10	0.52
1:B:50:LYS:O	1:B:54:LYS:HG3	2.09	0.52
1:D:30:ARG:CG	1:D:30:ARG:HH11	2.22	0.52
1:C:96:LYS:HG2	1:C:127:ALA:HB2	1.92	0.52
1:F:86:GLU:OE1	1:F:88:ALA:N	2.31	0.52
1:E:259:LEU:CD2	1:E:263:TYR:CE1	2.92	0.52
1:A:31:ASP:O	1:A:73:VAL:HA	2.10	0.52
1:E:41:LEU:CD1	1:E:41:LEU:H	2.17	0.52
1:C:30:ARG:HA	1:C:72:THR:O	2.09	0.52
1:C:22:ALA:HA	1:C:65:ILE:HG12	1.92	0.51
1:D:102:LEU:HD11	1:D:106:CYS:SG	2.50	0.51
1:D:138:HIS:HB3	1:D:150:GLU:HB2	1.92	0.51
1:E:30:ARG:HD3	1:E:31:ASP:O	2.09	0.51
1:E:130:ARG:HG3	1:E:130:ARG:O	2.10	0.51
1:E:49:SER:HA	1:E:52:VAL:CG2	2.41	0.51
1:A:194:MET:HG3	1:F:194:MET:HG3	1.92	0.51
1:B:214:LYS:HE3	1:C:175:ASP:OD1	2.10	0.51
1:C:201:ILE:HG23	1:C:206:ILE:HB	1.93	0.51
1:F:192:ILE:HG13	1:F:241:ILE:HG21	1.93	0.51
1:F:113:ALA:HA	1:F:135:ILE:O	2.11	0.51
1:D:192:ILE:HB	1:D:193:PRO:HD3	1.92	0.50
1:E:91:ARG:HD2	1:E:94:ILE:CD1	2.41	0.50
1:C:97:GLN:HA	1:C:97:GLN:NE2	2.26	0.50
1:B:107:LYS:HD2	1:B:109:GLU:CG	2.40	0.50
1:E:192:ILE:HB	1:E:193:PRO:HD3	1.93	0.50
1:A:86:GLU:CD	1:A:95:LYS:CE	2.80	0.49
1:D:62:LYS:O	1:D:66:LEU:HD12	2.13	0.49
1:A:140:PHE:CE2	1:A:148:LEU:HD23	2.48	0.49
1:B:123:GLU:HG3	1:F:265:ARG:HG2	1.93	0.49
1:D:169:SER:CB	1:D:176:PRO:HG3	2.43	0.49
1:C:30:ARG:HD3	1:C:31:ASP:N	2.27	0.49
1:F:-3:ARG:HH11	1:F:-3:ARG:CG	2.25	0.49
1:E:259:LEU:CD2	1:E:263:TYR:HE1	2.25	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:191:LEU:O	1:A:194:MET:HB2	2.13	0.48
1:B:52:VAL:HA	1:B:57:ILE:O	2.12	0.48
1:F:140:PHE:CE1	1:F:148:LEU:HD23	2.48	0.48
1:D:52:VAL:CG1	1:D:57:ILE:HG13	2.44	0.48
1:D:130:ARG:N	1:D:131:PRO:HD2	2.29	0.48
1:D:192:ILE:HG13	1:D:241:ILE:HG21	1.95	0.48
1:D:40:GLY:O	1:D:44:ILE:HG13	2.14	0.48
1:D:52:VAL:HG12	1:D:57:ILE:CG1	2.43	0.48
1:F:172:ILE:HG13	1:F:174:LYS:HG2	1.95	0.48
1:F:138:HIS:HB3	1:F:150:GLU:HB2	1.96	0.48
1:B:194:MET:CE	1:C:194:MET:CG	2.89	0.48
1:E:49:SER:OG	1:E:62:LYS:NZ	2.47	0.48
1:C:31:ASP:O	1:C:73:VAL:HA	2.14	0.47
1:B:143:ALA:N	1:B:144:PRO:CD	2.77	0.47
1:A:89:VAL:CG2	1:A:89:VAL:O	2.62	0.47
1:B:86:GLU:OE2	1:B:88:ALA:N	2.46	0.47
1:F:1:MET:HE2	2:F:303:HOH:O	2.13	0.47
1:F:147:LYS:HB3	1:F:147:LYS:HE2	1.55	0.47
1:A:143:ALA:N	1:A:144:PRO:CD	2.78	0.47
1:D:74:ASP:HB3	1:D:77:MET:CG	2.44	0.47
1:E:253:LYS:HD3	1:E:254:TYR:CZ	2.50	0.47
1:E:28:VAL:HG12	1:E:29:LEU:N	2.30	0.47
1:B:268:TRP:CZ3	1:D:155:ILE:HD13	2.50	0.47
1:D:102:LEU:HG	1:D:112:LEU:HD21	1.96	0.47
1:D:133:LYS:HB3	1:D:133:LYS:HE3	1.53	0.47
1:D:30:ARG:HD3	1:D:31:ASP:O	2.15	0.47
1:E:61:THR:O	1:E:64:GLU:N	2.48	0.46
1:B:253:LYS:HD3	1:B:254:TYR:CZ	2.49	0.46
1:D:133:LYS:HA	1:D:161:THR:HG21	1.96	0.46
1:D:72:THR:OG1	1:D:73:VAL:N	2.46	0.46
1:D:74:ASP:HB3	1:D:77:MET:HG2	1.96	0.46
1:D:93:ASP:N	1:D:93:ASP:OD1	2.48	0.46
1:D:140:PHE:O	1:D:146:MET:HB2	2.15	0.46
1:E:38:ASP:O	1:E:42:ASP:HB3	2.14	0.46
1:B:201:ILE:HG23	1:B:206:ILE:HB	1.98	0.46
1:E:37:VAL:HG12	1:E:41:LEU:HD11	1.97	0.46
1:B:30:ARG:HD3	1:B:31:ASP:N	2.31	0.46
1:C:143:ALA:O	1:C:174:LYS:NZ	2.37	0.46
1:E:79:ALA:HA	1:E:106:CYS:HA	1.98	0.46
1:B:122:THR:CG2	1:B:155:ILE:HG22	2.45	0.45
1:F:201:ILE:HG23	1:F:206:ILE:HB	1.97	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:268:TRP:CD2	1:F:275:LYS:HD3	2.50	0.45
1:B:155:ILE:HG21	1:F:266:ALA:HB2	1.98	0.45
1:A:86:GLU:OE1	1:A:95:LYS:CE	2.62	0.45
1:B:115:ASN:CG	1:B:115:ASN:O	2.55	0.45
1:A:113:ALA:HA	1:A:135:ILE:O	2.16	0.45
1:A:138:HIS:HB3	1:A:150:GLU:HB2	1.97	0.45
1:D:164:ALA:O	1:D:168:THR:CG2	2.61	0.45
1:F:135:ILE:HG21	1:F:165:VAL:HG21	1.99	0.45
1:D:102:LEU:HD21	1:D:112:LEU:HD23	1.95	0.45
1:A:175:ASP:CG	1:F:214:LYS:HE3	2.36	0.45
1:F:96:LYS:HG2	1:F:127:ALA:HB2	1.98	0.45
1:B:-5:VAL:HG13	1:B:-4:PRO:HD2	1.99	0.45
1:E:91:ARG:CG	1:E:93:ASP:OD1	2.65	0.45
1:B:192:ILE:HG13	1:B:241:ILE:HG21	1.99	0.45
1:E:30:ARG:HH12	1:E:32:ILE:HD12	1.81	0.44
1:E:72:THR:CG2	1:E:73:VAL:N	2.78	0.44
1:E:49:SER:O	1:E:52:VAL:HG23	2.17	0.44
1:B:-1:SER:HB2	1:B:168:THR:HA	1.98	0.44
1:F:160:GLU:HG3	2:F:315:HOH:O	2.16	0.44
1:B:92:MET:HG3	1:B:96:LYS:HD2	1.99	0.44
1:C:132:ASP:HB2	1:C:158:SER:HB3	1.99	0.44
1:D:3:LYS:HG2	1:D:26:GLU:HB3	2.00	0.44
1:E:30:ARG:CZ	1:E:75:LEU:HD23	2.48	0.44
1:A:194:MET:HG3	1:F:194:MET:CG	2.48	0.44
1:D:52:VAL:CG1	1:D:57:ILE:CG1	2.96	0.44
1:E:86:GLU:OE2	1:E:95:LYS:CE	2.60	0.44
1:F:-3:ARG:CG	1:F:-3:ARG:NH1	2.81	0.44
1:B:191:LEU:O	1:B:194:MET:HB3	2.18	0.43
1:B:189:ARG:HG2	1:C:201:ILE:HD11	1.99	0.43
1:D:30:ARG:HG2	1:D:30:ARG:HH11	1.83	0.43
1:B:192:ILE:N	1:B:193:PRO:CD	2.81	0.43
1:C:268:TRP:CD2	1:C:275:LYS:HD3	2.53	0.43
1:C:5:CYS:HB2	1:C:81:CYS:SG	2.59	0.43
1:D:105:ILE:O	1:D:105:ILE:HG22	2.19	0.43
1:A:101:ASP:O	1:A:105:ILE:HG13	2.18	0.43
1:F:130:ARG:HG3	1:F:130:ARG:O	2.19	0.43
1:A:130:ARG:O	1:A:130:ARG:HG3	2.19	0.43
1:E:240:ALA:O	1:E:244:VAL:HG23	2.19	0.43
1:A:33:LYS:O	1:A:37:VAL:HG23	2.19	0.43
1:A:89:VAL:O	1:A:89:VAL:HG23	2.18	0.43
1:D:209:VAL:HB	1:D:227:LEU:HD13	1.99	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:160:GLU:H	1:F:160:GLU:HG3	1.58	0.43
1:B:91:ARG:HD2	1:B:94:ILE:CD1	2.48	0.43
1:C:86:GLU:OE2	1:C:88:ALA:HB3	2.18	0.43
1:D:200:GLY:O	1:D:204:GLU:HG3	2.19	0.42
1:A:10:GLY:O	1:A:14:SER:OG	2.36	0.42
1:B:39:ARG:HB3	1:B:39:ARG:HE	1.48	0.42
1:C:141:ASN:OD1	1:C:142:PRO:CA	2.67	0.42
1:C:62:LYS:O	1:C:62:LYS:HG3	2.19	0.42
1:F:200:GLY:HA2	1:F:257:HIS:CD2	2.54	0.42
1:D:190:ILE:CG2	1:E:194:MET:HG3	2.50	0.42
1:E:72:THR:HG21	1:E:77:MET:CG	2.49	0.42
1:D:162:PHE:CE1	1:D:178:GLU:HG3	2.54	0.42
1:D:86:GLU:OE2	1:D:88:ALA:CB	2.64	0.42
1:A:206:ILE:N	1:A:206:ILE:CD1	2.83	0.42
1:D:56:LYS:HA	1:D:56:LYS:HD2	1.58	0.42
1:C:42:ASP:O	1:C:46:LYS:HG3	2.20	0.41
1:D:30:ARG:CG	1:D:30:ARG:NH1	2.83	0.41
1:E:30:ARG:NH2	1:E:75:LEU:HD23	2.33	0.41
1:F:129:LYS:HA	1:F:129:LYS:HD3	1.78	0.41
1:A:45:ASN:ND2	1:A:66:LEU:HD11	2.35	0.41
1:D:105:ILE:O	1:D:105:ILE:CG2	2.67	0.41
1:C:97:GLN:CA	1:C:97:GLN:HE21	2.32	0.41
1:C:97:GLN:CA	1:C:97:GLN:NE2	2.84	0.41
1:D:102:LEU:HD21	1:D:112:LEU:HD22	1.72	0.41
1:E:22:ALA:HA	1:E:65:ILE:HG12	2.02	0.41
1:A:204:GLU:CB	1:A:206:ILE:HD13	2.50	0.41
1:E:143:ALA:N	1:E:144:PRO:HD2	2.35	0.41
1:B:261:LYS:NZ	1:D:248:GLU:O	2.49	0.41
1:C:138:HIS:HB3	1:C:150:GLU:HB2	2.03	0.41
1:D:30:ARG:CD	1:D:31:ASP:O	2.69	0.41
1:C:92:MET:HE3	1:C:123:GLU:CD	2.41	0.41
1:C:141:ASN:OD1	1:C:142:PRO:N	2.53	0.41
1:F:191:LEU:O	1:F:194:MET:HB2	2.21	0.41
1:C:1:MET:HE2	1:C:82:ASP:C	2.41	0.41
1:E:74:ASP:OD1	1:E:76:ASN:ND2	2.35	0.41
1:E:72:THR:HG21	1:E:77:MET:HG3	2.03	0.41
1:D:30:ARG:HD3	1:D:73:VAL:HA	2.03	0.40
1:A:175:ASP:OD1	1:F:214:LYS:HE3	2.20	0.40
1:B:31:ASP:O	1:B:73:VAL:HA	2.21	0.40
1:D:28:VAL:HG12	1:D:29:LEU:N	2.36	0.40
1:E:236:ASP:OD1	1:E:236:ASP:N	2.54	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:203:ALA:CB	1:E:259:LEU:HD12	2.49	0.40
1:D:191:LEU:O	1:D:194:MET:HB3	2.21	0.40
1:E:37:VAL:CG1	1:E:41:LEU:HD11	2.50	0.40
1:C:142:PRO:HD2	1:C:146:MET:HG2	2.02	0.40
1:B:118:SER:HB2	1:B:244:VAL:HG21	2.04	0.40
1:B:89:VAL:HG22	1:B:90:GLU:H	1.87	0.40
1:D:183:PRO:O	1:D:189:ARG:HD3	2.22	0.40
1:E:118:SER:HB2	1:E:244:VAL:HG21	2.03	0.40
1:E:147:LYS:HD3	1:E:175:ASP:HB2	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	279/302 (92%)	269 (96%)	10 (4%)	0	100	100
1	B	287/302 (95%)	280 (98%)	7 (2%)	0	100	100
1	C	279/302 (92%)	267 (96%)	12 (4%)	0	100	100
1	D	279/302 (92%)	261 (94%)	16 (6%)	2 (1%)	22	39
1	E	279/302 (92%)	263 (94%)	16 (6%)	0	100	100
1	F	288/302 (95%)	283 (98%)	5 (2%)	0	100	100
All	All	1691/1812 (93%)	1623 (96%)	66 (4%)	2 (0%)	51	73

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	22	ALA
1	D	108	PRO

### 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	231/249 (93%)	211 (91%)	20 (9%)	10	20
1	B	237/249 (95%)	211 (89%)	26 (11%)	6	12
1	C	231/249 (93%)	211 (91%)	20 (9%)	10	20
1	D	231/249 (93%)	208 (90%)	23 (10%)	7	15
1	E	231/249 (93%)	205 (89%)	26 (11%)	6	11
1	F	238/249 (96%)	218 (92%)	20 (8%)	11	21
All	All	1399/1494 (94%)	1264 (90%)	135 (10%)	8	16

All (135) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	14	SER
1	A	23	LYS
1	A	35	GLU
1	A	52	VAL
1	A	67	THR
1	A	90	GLU
1	A	95	LYS
1	A	97	GLN
1	A	101	ASP
1	A	117	SER
1	A	133	LYS
1	A	137	MET
1	A	167	GLU
1	A	168	THR
1	A	206	ILE
1	A	252	SER
1	A	258	THR
1	A	272	LYS
1	A	279	ASP
1	A	281	SER
1	B	7	ILE
1	B	32	ILE

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Mol	Chain	Res	Type
1	B	39	ARG
1	B	52	VAL
1	B	53	LYS
1	B	58	GLU
1	B	61	THR
1	B	67	THR
1	B	70	SER
1	B	75	LEU
1	B	95	LYS
1	B	98	ILE
1	B	107	LYS
1	B	109	GLU
1	B	130	ARG
1	B	167	GLU
1	B	168	THR
1	B	204	GLU
1	B	206	ILE
1	B	209	VAL
1	B	247	SER
1	B	252	SER
1	B	253	LYS
1	B	258	THR
1	B	275	LYS
1	B	279	ASP
1	C	56	LYS
1	C	58	GLU
1	C	62	LYS
1	C	91	ARG
1	C	94	ILE
1	C	96	LYS
1	C	97	GLN
1	C	107	LYS
1	C	137	MET
1	C	149	VAL
1	C	189	ARG
1	C	199	VAL
1	C	209	VAL
1	C	237	ILE
1	C	252	SER
1	C	258	THR
1	C	259	LEU
1	C	271	ARG

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Mol	Chain	Res	Type
1	C	272	LYS
1	C	273	SER
1	D	1	MET
1	D	2	LYS
1	D	3	LYS
1	D	14	SER
1	D	30	ARG
1	D	46	LYS
1	D	50	LYS
1	D	53	LYS
1	D	56	LYS
1	D	61	THR
1	D	62	LYS
1	D	63	VAL
1	D	70	SER
1	D	75	LEU
1	D	96	LYS
1	D	107	LYS
1	D	110	THR
1	D	126	SER
1	D	133	LYS
1	D	155	ILE
1	D	160	GLU
1	D	168	THR
1	D	258	THR
1	E	2	LYS
1	E	26	GLU
1	E	29	LEU
1	E	32	ILE
1	E	34	ASP
1	E	46	LYS
1	E	47	ASN
1	E	52	VAL
1	E	64	GLU
1	E	67	THR
1	E	70	SER
1	E	75	LEU
1	E	94	ILE
1	E	96	LYS
1	E	107	LYS
1	E	117	SER
1	E	129	LYS

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Mol	Chain	Res	Type
1	E	147	LYS
1	E	149	VAL
1	E	189	ARG
1	E	210	GLU
1	E	247	SER
1	E	258	THR
1	E	259	LEU
1	E	279	ASP
1	E	281	SER
1	F	-8	SER
1	F	-3	ARG
1	F	3	LYS
1	F	39	ARG
1	F	41	LEU
1	F	46	LYS
1	F	54	LYS
1	F	56	LYS
1	F	63	VAL
1	F	89	VAL
1	F	94	ILE
1	F	96	LYS
1	F	137	MET
1	F	149	VAL
1	F	160	GLU
1	F	189	ARG
1	F	209	VAL
1	F	247	SER
1	F	252	SER
1	F	258	THR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	45	ASN
1	C	47	ASN
1	C	97	GLN
1	D	97	GLN
1	E	45	ASN

### 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](#)

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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	281/302 (93%)	-0.15	1 (0%) 92 93	28, 41, 71, 92	0
1	B	289/302 (95%)	0.08	6 (2%) 63 66	33, 51, 77, 101	0
1	C	281/302 (93%)	-0.10	2 (0%) 87 89	30, 49, 74, 94	0
1	D	281/302 (93%)	0.59	38 (13%) 3 2	33, 67, 108, 125	0
1	E	281/302 (93%)	0.03	1 (0%) 92 93	33, 58, 94, 111	0
1	F	290/302 (96%)	-0.15	1 (0%) 94 94	30, 41, 65, 82	0
All	All	1703/1812 (93%)	0.05	49 (2%) 51 55	28, 48, 92, 125	0

All (49) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	61	THR	4.9
1	D	34	ASP	3.8
1	D	33	LYS	3.7
1	D	63	VAL	3.7
1	D	60	ALA	3.7
1	D	52	VAL	3.5
1	D	42	ASP	3.5
1	B	36	PHE	3.5
1	D	45	ASN	3.5
1	D	69	ILE	3.4
1	D	74	ASP	3.4
1	D	36	PHE	3.4
1	D	32	ILE	3.3
1	D	25	PHE	3.3
1	D	26	GLU	3.3
1	B	33	LYS	3.3
1	D	29	LEU	3.2
1	C	43	PHE	3.2
1	D	80	ASP	3.2

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Mol	Chain	Res	Type	RSRZ
1	D	43	PHE	3.2
1	D	75	LEU	3.1
1	D	9	ALA	3.1
1	D	38	ASP	3.0
1	D	67	THR	3.0
1	D	35	GLU	3.0
1	B	32	ILE	2.9
1	A	53	LYS	2.9
1	D	77	MET	2.9
1	D	37	VAL	2.9
1	D	102	LEU	2.8
1	D	4	VAL	2.7
1	D	49	SER	2.6
1	D	41	LEU	2.6
1	D	129	LYS	2.6
1	D	51	LEU	2.6
1	D	57	ILE	2.5
1	D	55	GLY	2.4
1	D	70	SER	2.4
1	B	35	GLU	2.3
1	E	75	LEU	2.3
1	B	34	ASP	2.3
1	D	64	GLU	2.2
1	B	97	GLN	2.2
1	D	72	THR	2.2
1	D	20	PHE	2.1
1	F	281	SER	2.0
1	D	59	GLU	2.0
1	D	62	LYS	2.0
1	C	53	LYS	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](#)

There are no ligands in this entry.

## 6.5 Other polymers [i](#)

There are no such residues in this entry.