



## Full wwPDB X-ray Structure Validation Report ⓘ

May 29, 2020 – 02:36 am BST

PDB ID : 3GNS  
Title : Crystal Structure of the Staphylococcus aureus Enoyl-Acyl Carrier Protein Reductase (FabI) in apo form (one molecule in AU)  
Authors : Priyadarshi, A.; Hwang, K.Y.  
Deposited on : 2009-03-18  
Resolution : 2.71 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

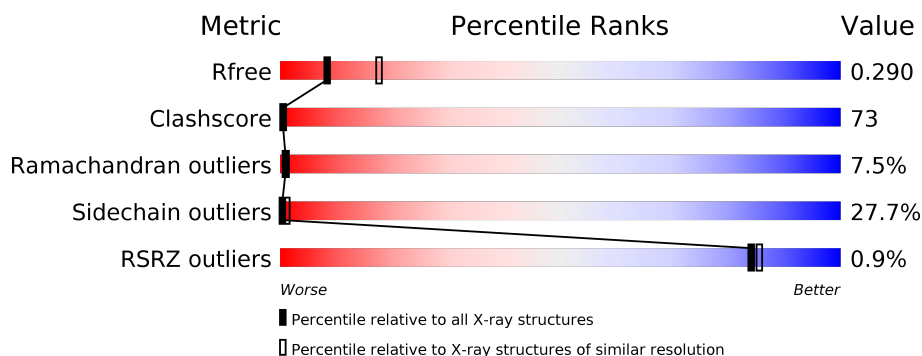
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.71 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	2808 (2.70-2.70)
Clashscore	141614	3122 (2.70-2.70)
Ramachandran outliers	138981	3069 (2.70-2.70)
Sidechain outliers	138945	3069 (2.70-2.70)
RSRZ outliers	127900	2737 (2.70-2.70)

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	260	

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 1716 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-[acyl-carrier-protein] reductase [NADH].

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	221	Total	C	N	O	S	0	0	0
			1704	1077	292	331	4			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-3	LEU	-	EXPRESSION TAG	UNP Q6GI75
A	-2	ALA	-	EXPRESSION TAG	UNP Q6GI75
A	-1	ALA	-	EXPRESSION TAG	UNP Q6GI75
A	0	ALA	-	EXPRESSION TAG	UNP Q6GI75

- Molecule 2 is SODIUM ION (three-letter code: NA) (formula: Na).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Na	0	0
			1	1		

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	11	Total	O	0	0
			11	11		

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.

- Chain A:**

Sequence logo showing the percentage of amino acids at each position for Chain A. The y-axis lists amino acid positions (T242 to S239). The x-axis represents the percentage of each amino acid at that position, with a color scale from green (17%) to red (15%).

Position	Percentage (%)
T242	17%
P180	17%
D181	17%
N182	17%
I183	17%
I184	17%
H185	17%
V186	17%
N187	17%
A188	17%
I189	17%
P192	17%
I193	17%
R194	17%
T195	17%
L196	17%
SER	17%
ALA	17%
LYS	17%
GLY	17%
VAL	17%
GLY	17%
GLY	17%
PHE	17%
N205	17%
T206	17%
I207	17%
L208	17%
K209	17%
E210	17%
I211	17%
E212	17%
E213	17%
R214	17%
A215	17%
P216	17%
I217	17%
R218	17%
R219	17%
N220	17%
V221	17%
D222	17%
Q223	17%
V224	17%
E225	17%
V226	17%
G227	17%
K228	17%
T229	17%
A230	17%
A231	17%
Y232	17%
L233	17%
L234	17%
S235	17%
D236	17%
S237	17%
S238	17%
S239	17%

## 4 Data and refinement statistics

Property	Value	Source
Space group	I 2 2 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	70.88Å 74.72Å 110.65Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	37.66 – 2.71 37.36 – 2.71	Depositor EDS
% Data completeness (in resolution range)	91.8 (37.66-2.71) 91.7 (37.36-2.71)	Depositor EDS
$R_{merge}$	0.08	Depositor
$R_{sym}$	0.28	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.32 (at 2.72Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, $R_{free}$	0.211 , 0.289 0.219 , 0.290	Depositor DCC
$R_{free}$ test set	347 reflections (4.55%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	46.9	Xtriage
Anisotropy	0.170	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.36 , 84.0	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.45$ , $\langle L^2 \rangle = 0.28$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.92	EDS
Total number of atoms	1716	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	41.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.25% 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

### 5.1 Standard geometry

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

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	2.17	66/1723 (3.8%)	1.96	53/2322 (2.3%)

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	1	10

All (66) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	72	GLU	CG-CD	9.30	1.65	1.51
1	A	71	GLU	CG-CD	8.92	1.65	1.51
1	A	53	GLU	CG-CD	8.82	1.65	1.51
1	A	173	TYR	CB-CG	8.76	1.64	1.51
1	A	131	GLU	CD-OE2	8.65	1.35	1.25
1	A	46	LYS	CD-CE	8.26	1.71	1.51
1	A	173	TYR	CG-CD1	8.02	1.49	1.39
1	A	173	TYR	CG-CD2	8.02	1.49	1.39
1	A	26	LYS	CD-CE	7.89	1.71	1.51
1	A	24	VAL	CB-CG2	7.06	1.67	1.52
1	A	213	GLU	CB-CG	7.04	1.65	1.52
1	A	131	GLU	CG-CD	6.98	1.62	1.51
1	A	53	GLU	CD-OE1	6.96	1.33	1.25
1	A	148	LEU	CG-CD2	6.90	1.77	1.51
1	A	183	ILE	CA-CB	-6.90	1.39	1.54
1	A	138	GLU	CG-CD	6.62	1.61	1.51
1	A	171	VAL	CB-CG2	-6.49	1.39	1.52
1	A	224	VAL	CB-CG2	-6.45	1.39	1.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	187	ALA	N-CA	6.31	1.58	1.46
1	A	148	LEU	CG-CD1	6.29	1.75	1.51
1	A	37	PHE	CE2-CZ	-6.21	1.25	1.37
1	A	182	ASN	CB-CG	-6.14	1.36	1.51
1	A	72	GLU	CD-OE2	6.10	1.32	1.25
1	A	47	GLU	CB-CG	6.08	1.63	1.52
1	A	90	VAL	CB-CG2	-6.01	1.40	1.52
1	A	22	PHE	CE2-CZ	5.99	1.48	1.37
1	A	195	THR	CB-CG2	5.99	1.72	1.52
1	A	5	GLU	CD-OE1	5.96	1.32	1.25
1	A	14	ILE	N-CA	-5.92	1.34	1.46
1	A	213	GLU	CG-CD	5.92	1.60	1.51
1	A	94	ILE	N-CA	5.83	1.58	1.46
1	A	59	GLU	CG-CD	5.83	1.60	1.51
1	A	189	SER	C-O	5.81	1.34	1.23
1	A	47	GLU	CG-CD	5.78	1.60	1.51
1	A	63	TYR	CD1-CE1	5.76	1.48	1.39
1	A	138	GLU	CD-OE1	5.71	1.31	1.25
1	A	239	SER	CA-CB	-5.69	1.44	1.52
1	A	72	GLU	CB-CG	5.68	1.62	1.52
1	A	225	GLU	CG-CD	5.68	1.60	1.51
1	A	217	LEU	C-O	5.68	1.34	1.23
1	A	72	GLU	CD-OE1	5.67	1.31	1.25
1	A	34	LYS	N-CA	5.67	1.57	1.46
1	A	253	HIS	C-O	5.59	1.33	1.23
1	A	239	SER	CB-OG	-5.59	1.34	1.42
1	A	60	ALA	N-CA	-5.55	1.35	1.46
1	A	67	VAL	N-CA	5.49	1.57	1.46
1	A	53	GLU	CB-CG	5.49	1.62	1.52
1	A	52	LEU	CG-CD2	5.46	1.72	1.51
1	A	91	TYR	CD1-CE1	5.46	1.47	1.39
1	A	38	THR	C-O	5.45	1.33	1.23
1	A	225	GLU	C-O	5.45	1.33	1.23
1	A	142	ILE	CA-CB	-5.41	1.42	1.54
1	A	71	GLU	CB-CG	5.40	1.62	1.52
1	A	94	ILE	CA-CB	5.39	1.67	1.54
1	A	256	LYS	CB-CG	5.33	1.67	1.52
1	A	5	GLU	CG-CD	5.24	1.59	1.51
1	A	218	LYS	CE-NZ	5.19	1.62	1.49
1	A	62	LEU	CG-CD1	5.18	1.71	1.51
1	A	138	GLU	CD-OE2	5.17	1.31	1.25
1	A	78	GLU	CD-OE2	5.13	1.31	1.25

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	22	PHE	CD1-CE1	5.12	1.49	1.39
1	A	230	ALA	CA-CB	-5.11	1.41	1.52
1	A	-1	ALA	C-O	-5.10	1.13	1.23
1	A	27	VAL	CB-CG2	-5.08	1.42	1.52
1	A	213	GLU	CD-OE2	5.08	1.31	1.25
1	A	185	VAL	CA-CB	-5.04	1.44	1.54

All (53) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	164	LYS	O-C-N	-13.84	100.56	122.70
1	A	237	LEU	CA-CB-CG	12.30	143.59	115.30
1	A	176	LEU	CA-CB-CG	-11.15	89.65	115.30
1	A	163	ALA	CB-CA-C	-10.68	94.08	110.10
1	A	249	ASP	CB-CG-OD1	-9.49	109.76	118.30
1	A	236	ASP	CB-CG-OD1	9.17	126.55	118.30
1	A	233	LEU	CB-CG-CD1	8.99	126.28	111.00
1	A	164	LYS	CA-C-N	8.75	136.46	117.20
1	A	148	LEU	CB-CG-CD2	8.70	125.80	111.00
1	A	148	LEU	CB-CG-CD1	-8.58	96.42	111.00
1	A	159	VAL	N-CA-C	8.55	134.07	111.00
1	A	162	VAL	CB-CA-C	-8.48	95.28	111.40
1	A	236	ASP	CB-CG-OD2	-8.27	110.86	118.30
1	A	84	VAL	CB-CA-C	-8.25	95.72	111.40
1	A	177	ASP	CB-CG-OD2	-8.03	111.08	118.30
1	A	219	ARG	CG-CD-NE	-7.87	95.27	111.80
1	A	62	LEU	CB-CG-CD1	7.32	123.44	111.00
1	A	43	ARG	NE-CZ-NH1	7.30	123.95	120.30
1	A	4	LEU	CB-CA-C	-7.26	96.41	110.20
1	A	236	ASP	CB-CA-C	-7.22	95.96	110.40
1	A	184	ARG	NE-CZ-NH1	6.55	123.58	120.30
1	A	29	ASP	CB-CG-OD2	6.45	124.10	118.30
1	A	52	LEU	CB-CG-CD2	6.43	121.94	111.00
1	A	90	VAL	CG1-CB-CG2	-6.40	100.66	110.90
1	A	160	MET	N-CA-CB	-6.40	99.08	110.60
1	A	35	LEU	CA-CB-CG	6.38	129.98	115.30
1	A	87	ILE	CG1-CB-CG2	-6.37	97.39	111.40
1	A	160	MET	CB-CA-C	6.34	123.08	110.40
1	A	83	ASP	CB-CG-OD2	6.28	123.95	118.30
1	A	84	VAL	CG1-CB-CG2	6.22	120.85	110.90
1	A	239	SER	N-CA-CB	-6.03	101.46	110.50
1	A	228	LYS	CD-CE-NZ	5.83	125.10	111.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	208	LEU	N-CA-C	-5.73	95.54	111.00
1	A	169	ALA	N-CA-C	-5.70	95.60	111.00
1	A	239	SER	C-N-CA	-5.60	110.53	122.30
1	A	225	GLU	OE1-CD-OE2	-5.58	116.61	123.30
1	A	247	HIS	N-CA-C	-5.56	95.99	111.00
1	A	43	ARG	NE-CZ-NH2	-5.54	117.53	120.30
1	A	1	MET	CA-CB-CG	5.50	122.65	113.30
1	A	78	GLU	N-CA-CB	5.42	120.35	110.60
1	A	70	ASP	C-N-CA	-5.35	108.31	121.70
1	A	119	ASP	CB-CG-OD2	-5.35	113.48	118.30
1	A	125	LEU	CB-CG-CD1	-5.34	101.93	111.00
1	A	176	LEU	CB-CG-CD2	5.28	119.98	111.00
1	A	128	VAL	CB-CA-C	-5.26	101.41	111.40
1	A	24	VAL	CA-CB-CG2	5.22	118.72	110.90
1	A	0	ALA	CB-CA-C	-5.17	102.34	110.10
1	A	130	HIS	CB-CA-C	5.16	120.73	110.40
1	A	0	ALA	N-CA-CB	-5.16	102.88	110.10
1	A	160	MET	N-CA-C	5.09	124.75	111.00
1	A	119	ASP	CB-CA-C	-5.07	100.25	110.40
1	A	184	ARG	CD-NE-CZ	5.06	130.69	123.60
1	A	80	ILE	CB-CA-C	-5.04	101.52	111.60

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
1	A	159	VAL	CA

All (10) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	-1	ALA	Peptide
1	A	120	ILE	Peptide
1	A	147	TYR	Peptide
1	A	158	ASN	Peptide
1	A	159	VAL	Peptide
1	A	161	GLY	Peptide
1	A	163	ALA	Peptide
1	A	208	LEU	Peptide
1	A	42	GLU	Peptide
1	A	43	ARG	Peptide

## 5.2 Too-close contacts ⓘ

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1704	0	1741	251	0
2	A	1	0	0	0	0
3	A	11	0	0	0	0
All	All	1716	0	1741	251	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 73.

All (251) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:148:LEU:CG	1:A:148:LEU:CD2	1.77	1.59
1:A:148:LEU:CD1	1:A:148:LEU:CG	1.75	1.59
1:A:159:VAL:CG1	1:A:160:MET:HB3	1.58	1.33
1:A:161:GLY:CA	1:A:162:VAL:O	1.83	1.26
1:A:160:MET:HG2	1:A:161:GLY:CA	1.66	1.25
1:A:157:TYR:C	1:A:158:ASN:OD1	1.79	1.19
1:A:160:MET:CG	1:A:161:GLY:HA2	1.72	1.19
1:A:158:ASN:H	1:A:159:VAL:CG2	1.58	1.16
1:A:159:VAL:HG12	1:A:160:MET:CB	1.76	1.16
1:A:160:MET:HG2	1:A:162:VAL:O	1.49	1.09
1:A:158:ASN:N	1:A:159:VAL:HG23	1.66	1.09
1:A:159:VAL:CB	1:A:160:MET:HB3	1.82	1.09
1:A:16:ASN:ND2	1:A:18:ARG:H	1.51	1.08
1:A:162:VAL:HG12	1:A:163:ALA:N	1.69	1.08
1:A:159:VAL:CG1	1:A:160:MET:CB	2.30	1.08
1:A:161:GLY:HA2	1:A:162:VAL:O	1.50	1.08
1:A:255:ILE:HG22	1:A:256:LYS:HE3	1.26	1.07
1:A:120:ILE:HG13	1:A:121:SER:HB2	1.38	1.03
1:A:158:ASN:CA	1:A:159:VAL:HG23	1.86	1.03
1:A:255:ILE:CG2	1:A:256:LYS:HE3	1.89	1.02
1:A:148:LEU:CD2	1:A:148:LEU:HG	1.91	1.00
1:A:160:MET:HG2	1:A:161:GLY:HA2	1.00	0.99
1:A:160:MET:CG	1:A:161:GLY:CA	2.33	0.98
1:A:33:ALA:O	1:A:35:LEU:HD13	1.65	0.96

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:159:VAL:HG12	1:A:160:MET:N	1.74	0.93
1:A:148:LEU:CD1	1:A:148:LEU:CB	2.45	0.93
1:A:159:VAL:HB	1:A:160:MET:HB3	1.47	0.93
1:A:160:MET:HG3	1:A:163:ALA:CB	1.99	0.93
1:A:158:ASN:H	1:A:159:VAL:HG23	1.21	0.92
1:A:16:ASN:HD21	1:A:18:ARG:CG	1.83	0.91
1:A:44:SER:C	1:A:46:LYS:H	1.75	0.90
1:A:160:MET:SD	1:A:161:GLY:N	2.45	0.89
1:A:123:TYR:CZ	1:A:127:ILE:HD11	2.08	0.89
1:A:158:ASN:C	1:A:159:VAL:HG23	1.90	0.87
1:A:16:ASN:HD21	1:A:18:ARG:HG3	1.38	0.87
1:A:44:SER:O	1:A:48:LEU:HD23	1.75	0.87
1:A:160:MET:CG	1:A:163:ALA:HB3	2.06	0.86
1:A:43:ARG:HG2	1:A:44:SER:N	1.89	0.86
1:A:162:VAL:O	1:A:163:ALA:HB3	1.75	0.85
1:A:162:VAL:HG12	1:A:163:ALA:H	1.39	0.84
1:A:160:MET:CG	1:A:163:ALA:CB	2.57	0.83
1:A:161:GLY:HA3	1:A:162:VAL:O	1.78	0.82
1:A:75:ASN:O	1:A:79:GLN:HB2	1.79	0.82
1:A:123:TYR:CE1	1:A:127:ILE:HD11	2.16	0.81
1:A:205:ASN:N	1:A:207:ILE:HG13	1.95	0.81
1:A:157:TYR:O	1:A:158:ASN:OD1	1.98	0.79
1:A:158:ASN:N	1:A:159:VAL:CG2	2.29	0.79
1:A:44:SER:C	1:A:46:LYS:N	2.31	0.79
1:A:160:MET:SD	1:A:161:GLY:CA	2.70	0.79
1:A:18:ARG:HD3	1:A:195:THR:O	1.83	0.79
1:A:186:ASN:HD21	1:A:242:THR:HA	1.48	0.78
1:A:48:LEU:HA	1:A:51:LEU:HB2	1.66	0.77
1:A:160:MET:HG2	1:A:163:ALA:HB3	1.66	0.77
1:A:253:HIS:O	1:A:256:LYS:HG2	1.84	0.77
1:A:7:LYS:HD2	1:A:88:ASP:OD1	1.83	0.77
1:A:172:LYS:O	1:A:176:LEU:HD12	1.85	0.76
1:A:162:VAL:O	1:A:163:ALA:CB	2.35	0.74
1:A:43:ARG:HG2	1:A:44:SER:CA	2.16	0.74
1:A:158:ASN:CA	1:A:159:VAL:CG2	2.64	0.74
1:A:159:VAL:HG12	1:A:160:MET:HB2	1.69	0.74
1:A:164:LYS:O	1:A:165:ALA:C	2.23	0.73
1:A:16:ASN:HD21	1:A:18:ARG:H	1.36	0.73
1:A:3:ASN:HB3	1:A:7:LYS:HE2	1.71	0.72
1:A:160:MET:HG3	1:A:163:ALA:HB3	1.70	0.71
1:A:47:GLU:O	1:A:50:LYS:HB3	1.91	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:205:ASN:C	1:A:207:ILE:N	2.39	0.70
1:A:52:LEU:HD23	1:A:52:LEU:O	1.91	0.70
1:A:119:ASP:HB3	1:A:123:TYR:H	1.55	0.70
1:A:40:ARG:HA	1:A:65:ILE:O	1.92	0.70
1:A:160:MET:HG2	1:A:161:GLY:HA3	1.71	0.70
1:A:118:GLN:O	1:A:120:ILE:N	2.25	0.70
1:A:159:VAL:HG11	1:A:160:MET:HB3	1.66	0.70
1:A:159:VAL:CG1	1:A:160:MET:N	2.49	0.69
1:A:77:PHE:HA	1:A:80:ILE:HD12	1.75	0.69
1:A:16:ASN:ND2	1:A:18:ARG:N	2.34	0.69
1:A:205:ASN:O	1:A:208:LEU:HD23	1.92	0.69
1:A:161:GLY:C	1:A:162:VAL:O	2.28	0.69
1:A:158:ASN:C	1:A:159:VAL:CG2	2.62	0.68
1:A:184:ARG:NH1	1:A:238:SER:OG	2.21	0.68
1:A:119:ASP:C	1:A:121:SER:H	1.96	0.68
1:A:43:ARG:HG2	1:A:44:SER:HA	1.76	0.68
1:A:164:LYS:O	1:A:166:SER:N	2.27	0.67
1:A:205:ASN:C	1:A:207:ILE:H	1.98	0.67
1:A:43:ARG:CG	1:A:44:SER:N	2.58	0.67
1:A:162:VAL:CG1	1:A:163:ALA:N	2.45	0.66
1:A:160:MET:CB	1:A:161:GLY:HA2	2.23	0.66
1:A:16:ASN:ND2	1:A:18:ARG:CG	2.57	0.66
1:A:70:ASP:OD2	1:A:70:ASP:N	2.28	0.66
1:A:69:SER:O	1:A:72:GLU:HB2	1.95	0.66
1:A:79:GLN:HG3	1:A:83:ASP:OD2	1.95	0.66
1:A:172:LYS:O	1:A:175:ALA:HB3	1.97	0.65
1:A:5:GLU:O	1:A:6:ASN:HB2	1.96	0.65
1:A:24:VAL:HG12	1:A:28:LEU:HD12	1.79	0.64
1:A:3:ASN:ND2	1:A:236:ASP:CG	2.51	0.64
1:A:39:TYR:O	1:A:65:ILE:N	2.30	0.64
1:A:158:ASN:N	1:A:158:ASN:OD1	2.30	0.64
1:A:47:GLU:O	1:A:50:LYS:CB	2.45	0.63
1:A:16:ASN:ND2	1:A:18:ARG:HG2	2.14	0.62
1:A:39:TYR:CE2	1:A:64:GLN:HB2	2.34	0.62
1:A:124:SER:O	1:A:128:VAL:HG23	2.00	0.62
1:A:211:ILE:HG23	1:A:215:ALA:HB2	1.82	0.62
1:A:164:LYS:O	1:A:167:LEU:N	2.33	0.62
1:A:177:ASP:HB3	1:A:178:LEU:HD23	1.81	0.62
1:A:159:VAL:HG12	1:A:160:MET:CA	2.29	0.61
1:A:22:PHE:O	1:A:26:LYS:HG3	2.00	0.61
1:A:14:ILE:CG1	1:A:14:ILE:O	2.47	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:29:ASP:OD2	1:A:57:GLN:NE2	2.27	0.61
1:A:43:ARG:O	1:A:45:ARG:N	2.33	0.61
1:A:39:TYR:O	1:A:64:GLN:HA	2.00	0.61
1:A:211:ILE:O	1:A:212:GLU:C	2.38	0.61
1:A:160:MET:HG3	1:A:163:ALA:HB2	1.83	0.60
1:A:157:TYR:N	1:A:158:ASN:OD1	2.33	0.60
1:A:67:VAL:HB	1:A:124:SER:OG	2.01	0.59
1:A:159:VAL:HG12	1:A:160:MET:H	1.65	0.59
1:A:119:ASP:C	1:A:121:SER:N	2.56	0.59
1:A:160:MET:CG	1:A:163:ALA:HB2	2.33	0.59
1:A:67:VAL:CG1	1:A:124:SER:OG	2.51	0.59
1:A:50:LYS:O	1:A:53:GLU:HB2	2.02	0.59
1:A:3:ASN:ND2	1:A:236:ASP:OD2	2.35	0.58
1:A:119:ASP:N	1:A:119:ASP:OD1	2.35	0.58
1:A:7:LYS:HB3	1:A:88:ASP:CG	2.23	0.58
1:A:118:GLN:HA	1:A:118:GLN:OE1	2.03	0.58
1:A:44:SER:O	1:A:46:LYS:N	2.36	0.58
1:A:178:LEU:N	1:A:178:LEU:HD23	2.19	0.58
1:A:16:ASN:HD22	1:A:17:LYS:H	1.52	0.58
1:A:217:LEU:O	1:A:219:ARG:HG2	2.04	0.57
1:A:44:SER:O	1:A:47:GLU:N	2.37	0.57
1:A:40:ARG:NE	1:A:66:ASP:OD1	2.33	0.57
1:A:24:VAL:CG1	1:A:28:LEU:HD12	2.35	0.57
1:A:48:LEU:C	1:A:50:LYS:N	2.53	0.56
1:A:123:TYR:O	1:A:127:ILE:HG12	2.05	0.56
1:A:211:ILE:O	1:A:213:GLU:N	2.37	0.56
1:A:119:ASP:CB	1:A:122:SER:HB3	2.36	0.56
1:A:20:ILE:HG21	1:A:226:VAL:HG11	1.87	0.56
1:A:44:SER:O	1:A:48:LEU:CD2	2.51	0.56
1:A:186:ASN:ND2	1:A:242:THR:HA	2.19	0.56
1:A:253:HIS:O	1:A:256:LYS:CG	2.53	0.56
1:A:147:TYR:CE1	1:A:193:ILE:HD12	2.41	0.56
1:A:3:ASN:O	1:A:4:LEU:HB2	2.06	0.55
1:A:159:VAL:CG1	1:A:160:MET:HB2	2.32	0.55
1:A:231:ALA:O	1:A:235:SER:HB3	2.06	0.55
1:A:162:VAL:CG1	1:A:163:ALA:H	2.16	0.55
1:A:160:MET:SD	1:A:160:MET:C	2.86	0.54
1:A:46:LYS:O	1:A:50:LYS:HB2	2.08	0.54
1:A:159:VAL:CB	1:A:160:MET:CB	2.72	0.54
1:A:37:PHE:HB2	1:A:62:LEU:HD12	1.89	0.54
1:A:90:VAL:O	1:A:90:VAL:HG22	2.05	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:248:VAL:O	1:A:248:VAL:HG12	2.08	0.54
1:A:249:ASP:OD1	1:A:249:ASP:N	2.39	0.54
1:A:48:LEU:C	1:A:50:LYS:H	2.02	0.54
1:A:160:MET:SD	1:A:161:GLY:HA3	2.47	0.54
1:A:157:TYR:CA	1:A:158:ASN:OD1	2.54	0.53
1:A:46:LYS:HD3	1:A:49:GLU:OE2	2.08	0.53
1:A:159:VAL:HB	1:A:160:MET:CB	2.30	0.53
1:A:26:LYS:O	1:A:30:GLN:HB2	2.08	0.53
1:A:158:ASN:H	1:A:159:VAL:HG21	1.66	0.53
1:A:46:LYS:C	1:A:48:LEU:N	2.52	0.53
1:A:148:LEU:CD1	1:A:148:LEU:HB2	2.34	0.53
1:A:27:VAL:O	1:A:31:LEU:HD22	2.07	0.53
1:A:16:ASN:ND2	1:A:17:LYS:N	2.56	0.53
1:A:46:LYS:O	1:A:47:GLU:C	2.46	0.53
1:A:164:LYS:HA	1:A:167:LEU:HB3	1.90	0.52
1:A:45:ARG:HG3	1:A:45:ARG:O	2.08	0.52
1:A:160:MET:HG2	1:A:163:ALA:CB	2.32	0.52
1:A:10:VAL:HB	1:A:90:VAL:HB	1.92	0.52
1:A:17:LYS:HB2	1:A:51:LEU:HD21	1.91	0.52
1:A:166:SER:O	1:A:169:ALA:HB3	2.09	0.52
1:A:68:GLN:HA	1:A:68:GLN:OE1	2.09	0.52
1:A:79:GLN:O	1:A:82:LYS:HB2	2.09	0.52
1:A:76:GLY:O	1:A:79:GLN:HB3	2.10	0.51
1:A:40:ARG:HG2	1:A:40:ARG:HH11	1.76	0.51
1:A:75:ASN:O	1:A:79:GLN:CB	2.56	0.51
1:A:195:THR:O	1:A:195:THR:OG1	2.29	0.51
1:A:34:LYS:H	1:A:34:LYS:HD3	1.75	0.51
1:A:143:VAL:CG2	1:A:233:LEU:HB3	2.42	0.50
1:A:36:VAL:HG12	1:A:37:PHE:N	2.26	0.50
1:A:223:GLN:O	1:A:227:GLY:N	2.32	0.50
1:A:14:ILE:HG13	1:A:14:ILE:O	2.11	0.50
1:A:194:ARG:NH1	1:A:212:GLU:OE2	2.40	0.50
1:A:16:ASN:ND2	1:A:17:LYS:H	2.10	0.50
1:A:44:SER:C	1:A:48:LEU:HD23	2.32	0.50
1:A:80:ILE:O	1:A:80:ILE:HG22	2.12	0.49
1:A:2:LEU:HG	1:A:237:LEU:CD2	2.43	0.49
1:A:119:ASP:HB3	1:A:122:SER:HB3	1.93	0.49
1:A:33:ALA:O	1:A:35:LEU:CD1	2.50	0.49
1:A:12:MET:O	1:A:93:SER:HB2	2.13	0.49
1:A:205:ASN:C	1:A:205:ASN:OD1	2.51	0.49
1:A:-1:ALA:O	1:A:2:LEU:HD13	2.13	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:24:VAL:HA	1:A:227:GLY:HA3	1.95	0.48
1:A:41:LYS:HG3	1:A:42:GLU:HG3	1.93	0.48
1:A:90:VAL:HG13	1:A:142:ILE:HG12	1.95	0.48
1:A:148:LEU:HD23	1:A:148:LEU:N	2.29	0.48
1:A:230:ALA:O	1:A:231:ALA:C	2.51	0.48
1:A:252:PHE:O	1:A:255:ILE:HB	2.14	0.48
1:A:24:VAL:HG21	1:A:91:TYR:CZ	2.49	0.48
1:A:6:ASN:CG	1:A:34:LYS:HZ1	2.18	0.47
1:A:160:MET:CG	1:A:161:GLY:HA3	2.37	0.47
1:A:41:LYS:HB2	1:A:41:LYS:HE2	1.63	0.47
1:A:193:ILE:HG12	1:A:222:ASP:HA	1.97	0.46
1:A:36:VAL:HG12	1:A:37:PHE:H	1.79	0.46
1:A:189:SER:OG	1:A:245:ASN:ND2	2.39	0.46
1:A:147:TYR:HE1	1:A:193:ILE:HD12	1.80	0.45
1:A:16:ASN:HD21	1:A:18:ARG:N	2.06	0.45
1:A:80:ILE:CG2	1:A:80:ILE:O	2.62	0.45
1:A:231:ALA:O	1:A:232:TYR:C	2.55	0.44
1:A:2:LEU:HG	1:A:237:LEU:HD22	1.99	0.44
1:A:79:GLN:NE2	1:A:79:GLN:HA	2.32	0.44
1:A:20:ILE:O	1:A:20:ILE:HG22	2.17	0.44
1:A:24:VAL:HG21	1:A:91:TYR:CE1	2.52	0.44
1:A:211:ILE:C	1:A:213:GLU:N	2.65	0.44
1:A:232:TYR:CZ	1:A:238:SER:HB3	2.53	0.44
1:A:133:LYS:C	1:A:135:LEU:N	2.70	0.44
1:A:87:ILE:HD13	1:A:87:ILE:HG21	1.71	0.44
1:A:16:ASN:ND2	1:A:18:ARG:HG3	2.18	0.44
1:A:47:GLU:O	1:A:51:LEU:HG	2.18	0.44
1:A:38:THR:HG21	1:A:65:ILE:HD12	2.00	0.44
1:A:46:LYS:O	1:A:48:LEU:N	2.51	0.43
1:A:144:ALA:O	1:A:187:ALA:HA	2.18	0.43
1:A:210:GLU:O	1:A:214:ARG:HG3	2.18	0.43
1:A:205:ASN:O	1:A:205:ASN:OD1	2.36	0.43
1:A:3:ASN:ND2	1:A:236:ASP:OD1	2.52	0.43
1:A:28:LEU:O	1:A:31:LEU:N	2.49	0.43
1:A:119:ASP:HA	1:A:122:SER:CB	2.48	0.43
1:A:26:LYS:O	1:A:27:VAL:C	2.57	0.43
1:A:88:ASP:OD2	1:A:88:ASP:N	2.52	0.42
1:A:9:TYR:N	1:A:9:TYR:CD1	2.87	0.42
1:A:159:VAL:HG11	1:A:160:MET:CB	2.34	0.42
1:A:222:ASP:HB2	1:A:223:GLN:OE1	2.20	0.42
1:A:77:PHE:CZ	1:A:128:VAL:HG13	2.55	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:171:VAL:HG21	1:A:187:ALA:HB2	2.01	0.42
1:A:19:SER:O	1:A:22:PHE:HB3	2.19	0.42
1:A:78:GLU:O	1:A:82:LYS:HG3	2.19	0.42
1:A:133:LYS:HA	1:A:136:MET:CG	2.50	0.41
1:A:9:TYR:CD2	1:A:234:LEU:HD13	2.55	0.41
1:A:183:ILE:HG22	1:A:183:ILE:O	2.19	0.41
1:A:20:ILE:O	1:A:20:ILE:CG2	2.67	0.41
1:A:40:ARG:HH11	1:A:40:ARG:CG	2.33	0.41
1:A:18:ARG:HD2	1:A:196:LEU:HA	2.03	0.41
1:A:40:ARG:NH1	1:A:40:ARG:CG	2.84	0.41
1:A:148:LEU:CD2	1:A:148:LEU:N	2.83	0.41
1:A:208:LEU:C	1:A:210:GLU:N	2.72	0.41
1:A:119:ASP:HB3	1:A:123:TYR:N	2.31	0.41
1:A:119:ASP:CA	1:A:122:SER:HB3	2.50	0.41
1:A:46:LYS:HD3	1:A:46:LYS:HA	1.89	0.41
1:A:18:ARG:HG2	1:A:18:ARG:NH1	2.37	0.41
1:A:20:ILE:O	1:A:23:GLY:N	2.54	0.40
1:A:36:VAL:CG1	1:A:37:PHE:N	2.84	0.40
1:A:219:ARG:NH2	1:A:225:GLU:OE1	2.54	0.40
1:A:188:ILE:CD1	1:A:229:THR:HG22	2.51	0.40
1:A:176:LEU:O	1:A:177:ASP:C	2.58	0.40
1:A:55:LEU:HG	1:A:55:LEU:H	1.61	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	213/260 (82%)	169 (79%)	28 (13%)	16 (8%)	<b>1</b> <b>1</b>

All (16) Ramachandran outliers are listed below:



Mol	Chain	Res	Type
1	A	43	ARG
1	A	119	ASP
1	A	121	SER
1	A	159	VAL
1	A	160	MET
1	A	162	VAL
1	A	250	SER
1	A	0	ALA
1	A	45	ARG
1	A	67	VAL
1	A	165	ALA
1	A	239	SER
1	A	255	ILE
1	A	6	ASN
1	A	208	LEU
1	A	44	SER

### 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	184/211 (87%)	133 (72%)	51 (28%)	<b>0</b> <b>1</b>

All (51) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	-3	LEU
1	A	14	ILE
1	A	18	ARG
1	A	30	GLN
1	A	31	LEU
1	A	34	LYS
1	A	35	LEU
1	A	38	THR
1	A	40	ARG
1	A	41	LYS
1	A	43	ARG

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Mol	Chain	Res	Type
1	A	44	SER
1	A	45	ARG
1	A	46	LYS
1	A	49	GLU
1	A	50	LYS
1	A	59	GLU
1	A	62	LEU
1	A	67	VAL
1	A	70	ASP
1	A	73	VAL
1	A	78	GLU
1	A	84	VAL
1	A	88	ASP
1	A	90	VAL
1	A	93	SER
1	A	118	GLN
1	A	119	ASP
1	A	121	SER
1	A	136	MET
1	A	146	THR
1	A	148	LEU
1	A	157	TYR
1	A	158	ASN
1	A	159	VAL
1	A	167	LEU
1	A	176	LEU
1	A	178	LEU
1	A	180	PRO
1	A	189	SER
1	A	192	PRO
1	A	194	ARG
1	A	195	THR
1	A	219	ARG
1	A	220	ASN
1	A	228	LYS
1	A	233	LEU
1	A	239	SER
1	A	245	ASN
1	A	255	ILE
1	A	256	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	3	ASN
1	A	16	ASN
1	A	79	GLN
1	A	86	ASN
1	A	130	HIS
1	A	170	ASN
1	A	186	ASN
1	A	245	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](#)

Of 1 ligands modelled in this entry, 1 is 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 [i](#)

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 [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, 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	221/260 (85%)	-0.35	2 (0%) 84 85	13, 38, 69, 90	0

All (2) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	67	VAL	2.1
1	A	159	VAL	2.0

### 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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
2	NA	A	257	1/1	0.92	0.16	59,59,59,59	0

### 6.5 Other polymers [i](#)

There are no such residues in this entry.