



# Full wwPDB X-ray Structure Validation Report ⓘ

Mar 14, 2018 – 01:51 pm GMT

PDB ID : 3T5C  
Title : Crystal structure of N-terminal domain of FACL13 from Mycobacterium tuberculosis in different space group C2  
Authors : Goyal, A.; Sankaranarayanan, R.  
Deposited on : 2011-07-27  
Resolution : 2.09 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	:	trunk31020
Percentile statistics	:	20171227.v01 (using entries in the PDB archive December 27th 2017)
Refmac	:	5.8.0158
CCP4	:	7.0 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	trunk31020

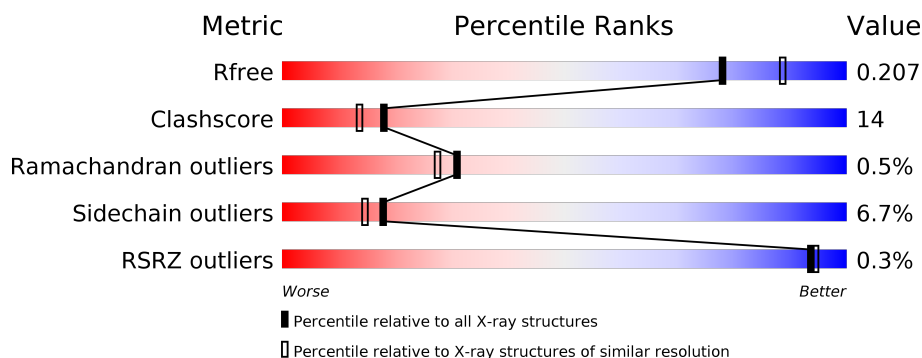
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.09 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))
$R_{free}$	111664	4608 (2.10-2.10)
Clashscore	122126	5109 (2.10-2.10)
Ramachandran outliers	120053	5059 (2.10-2.10)
Sidechain outliers	120020	5060 (2.10-2.10)
RSRZ outliers	108989	4497 (2.10-2.10)

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. 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	396	<div> <div></div> <div>70%26%..</div> </div>
1	B	396	<div> <div></div> <div>72%24%..</div> </div>

## 2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 6461 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 PROBABLE CHAIN-FATTY-ACID-CoA LIGASE FADD13.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	392	Total	C	N	O	S	0	0	0
			2985	1892	511	564	18			
1	B	392	Total	C	N	O	S	0	0	0
			2985	1892	511	564	18			

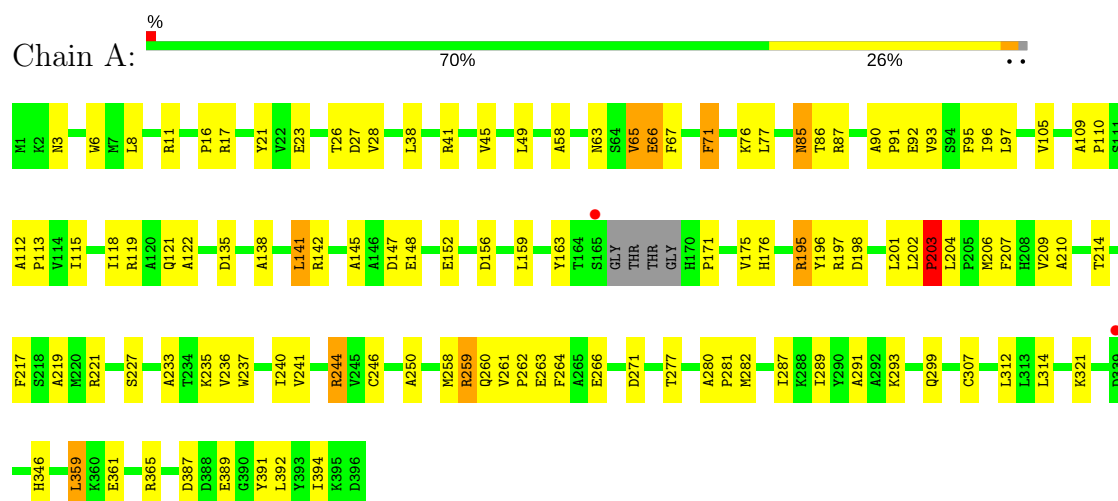
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	241	Total	O	0	0
			241	241		
2	B	250	Total	O	0	0
			250	250		

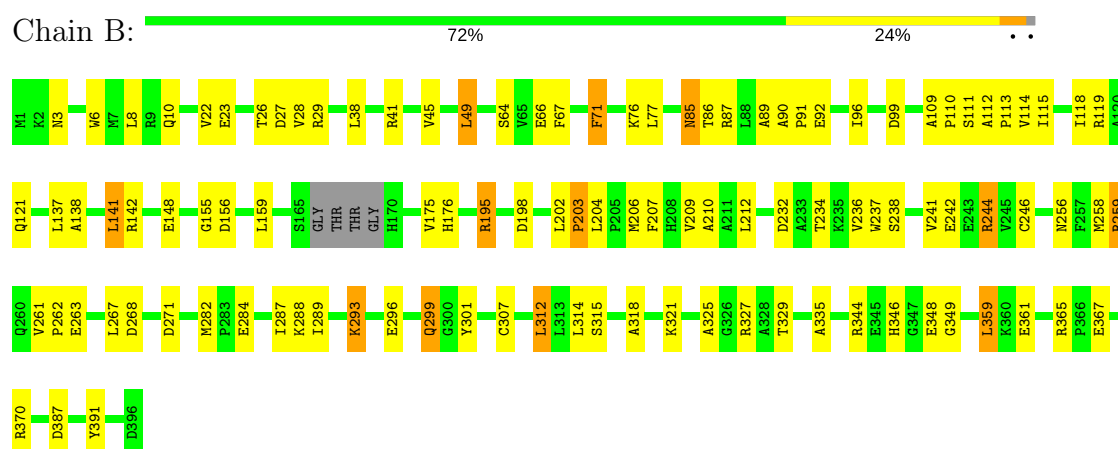
### 3 Residue-property plots

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: PROBABLE CHAIN-FATTY-ACID-CoA LIGASE FADD13



#### • Molecule 1: PROBABLE CHAIN-FATTY-ACID-CoA LIGASE FADD13



## 4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	129.23Å 91.98Å 84.41Å 90.00° 130.79° 90.00°	Depositor
Resolution (Å)	25.00 – 2.09 24.85 – 2.09	Depositor EDS
% Data completeness (in resolution range)	(Not available) (25.00-2.09) 97.4 (24.85-2.09)	Depositor EDS
$R_{merge}$	0.04	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.64 (at 2.08Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, $R_{free}$	0.208 , 0.263 0.205 , 0.207	Depositor DCC
$R_{free}$ test set	2190 reflections (5.04%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	34.3	Xtriage
Anisotropy	0.264	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.30 , 33.6	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.487 for -h-2*k,l	Xtriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	6461	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	36.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.98% 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	0.33	0/3047	0.59	0/4142
1	B	0.33	0/3047	0.59	0/4142
All	All	0.33	0/6094	0.59	0/8284

There are no bond length outliers.

There are no bond angle outliers.

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	2985	0	2963	86	1
1	B	2985	0	2963	87	0
2	A	241	0	0	1	0
2	B	250	0	0	3	0
All	All	6461	0	5926	172	1

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 14.

All (172) 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:282:MET:H	1:A:299:GLN:HE22	1.13	0.93

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:282:MET:H	1:B:299:GLN:HE22	1.02	0.92
1:B:26:THR:HG23	1:B:28:VAL:HG23	1.59	0.84
1:A:244:ARG:HG3	1:A:271:ASP:OD2	1.81	0.81
1:A:282:MET:N	1:A:299:GLN:HE22	1.83	0.77
1:B:85:ASN:ND2	1:B:87:ARG:H	1.83	0.76
1:A:201:LEU:HD22	1:A:240:ILE:HD11	1.67	0.76
1:A:26:THR:HG23	1:A:28:VAL:HG23	1.67	0.75
1:A:85:ASN:ND2	1:A:87:ARG:H	1.85	0.75
1:A:112:ALA:HB3	1:A:113:PRO:HD3	1.72	0.70
1:A:23:GLU:HB3	1:A:26:THR:HG22	1.73	0.70
1:B:244:ARG:HG3	1:B:271:ASP:OD2	1.91	0.69
1:B:41:ARG:O	1:B:45:VAL:HG23	1.93	0.69
1:B:118:ILE:O	1:B:121:GLN:HG2	1.93	0.68
1:B:282:MET:N	1:B:299:GLN:HE22	1.84	0.67
1:A:260:GLN:HE21	1:A:260:GLN:HA	1.60	0.67
1:B:175:VAL:HG22	1:B:361:GLU:O	1.95	0.67
1:B:175:VAL:HG21	1:B:361:GLU:HG2	1.77	0.66
1:B:112:ALA:HB3	1:B:113:PRO:HD3	1.79	0.65
1:A:109:ALA:HB3	1:A:110:PRO:HD3	1.79	0.64
1:A:11:ARG:HD2	1:A:219:ALA:O	1.98	0.64
1:B:282:MET:H	1:B:299:GLN:NE2	1.86	0.63
1:A:261:VAL:HG23	1:A:262:PRO:HD2	1.80	0.63
1:B:176:HIS:CE1	1:B:307:CYS:HB2	2.34	0.62
1:A:118:ILE:O	1:A:121:GLN:HG2	2.00	0.62
1:A:237:TRP:O	1:A:241:VAL:HG23	1.99	0.61
1:B:206:MET:O	1:B:212:LEU:HD13	2.00	0.61
1:A:259:ARG:HH12	1:A:289:ILE:HD12	1.64	0.61
1:B:296:GLU:HB3	1:B:315:SER:HB2	1.83	0.61
1:B:259:ARG:HD2	1:B:259:ARG:C	2.21	0.60
1:B:232:ASP:O	1:B:236:VAL:HG13	2.02	0.60
1:B:85:ASN:HD22	1:B:85:ASN:C	2.05	0.60
1:B:23:GLU:HB3	1:B:26:THR:HG22	1.84	0.59
1:B:90:ALA:HB3	1:B:91:PRO:HD3	1.84	0.59
1:A:11:ARG:CD	1:A:219:ALA:O	2.51	0.59
1:A:142:ARG:HG2	1:A:142:ARG:HH11	1.68	0.59
1:B:367:GLU:HA	1:B:370:ARG:HH21	1.69	0.58
1:A:359:LEU:N	1:A:359:LEU:HD12	2.19	0.58
1:B:92:GLU:O	1:B:96:ILE:HG12	2.04	0.57
1:A:314:LEU:HD22	1:A:314:LEU:N	2.20	0.56
1:B:258:MET:O	1:B:261:VAL:HG12	2.05	0.56
1:A:175:VAL:HG22	1:A:361:GLU:O	2.06	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:235:LYS:HB3	2:A:534:HOH:O	2.04	0.56
1:B:6:TRP:O	1:B:10:GLN:HG3	2.06	0.56
1:B:85:ASN:HD22	1:B:86:THR:N	2.04	0.56
1:A:138:ALA:O	1:A:142:ARG:HG3	2.06	0.56
1:B:261:VAL:HG23	1:B:262:PRO:HD2	1.87	0.56
1:B:109:ALA:HB3	1:B:110:PRO:HD3	1.87	0.55
1:B:238:SER:O	1:B:242:GLU:HG3	2.06	0.55
1:A:95:PHE:CE2	1:A:171:PRO:HD2	2.42	0.55
1:A:17:ARG:NH2	1:B:327:ARG:HH12	2.05	0.55
1:A:259:ARG:NH1	1:A:289:ILE:HD12	2.22	0.54
1:A:90:ALA:HB3	1:A:91:PRO:HD3	1.89	0.54
1:A:63:ASN:ND2	1:A:206:MET:HG3	2.23	0.53
1:A:233:ALA:HA	1:A:236:VAL:HG22	1.90	0.53
1:B:195:ARG:HG2	1:B:198:ASP:OD2	2.08	0.53
1:B:76:LYS:NZ	1:B:148:GLU:OE2	2.37	0.53
1:B:26:THR:HG23	1:B:28:VAL:CG2	2.37	0.53
1:B:64:SER:OG	1:B:66:GLU:HG2	2.08	0.53
1:A:41:ARG:O	1:A:45:VAL:HG23	2.09	0.53
1:A:92:GLU:O	1:A:96:ILE:HG12	2.09	0.52
1:A:26:THR:HG23	1:A:28:VAL:CG2	2.36	0.52
1:A:85:ASN:C	1:A:85:ASN:HD22	2.11	0.52
1:B:142:ARG:HH11	1:B:142:ARG:HG2	1.75	0.52
1:A:214:THR:HA	1:A:217:PHE:CE2	2.45	0.52
1:A:359:LEU:HD12	1:A:359:LEU:H	1.75	0.52
1:A:85:ASN:HD21	1:A:87:ARG:HB2	1.75	0.51
1:A:96:ILE:HD12	1:A:163:TYR:CG	2.45	0.51
1:A:176:HIS:CE1	1:A:307:CYS:HB2	2.45	0.51
1:A:85:ASN:HD22	1:A:86:THR:N	2.09	0.51
1:B:209:VAL:HG13	1:B:210:ALA:N	2.25	0.51
1:B:22:VAL:HG13	1:B:29:ARG:HG2	1.93	0.50
1:A:21:TYR:CE2	1:A:65:VAL:HG22	2.46	0.50
1:A:244:ARG:HG3	1:A:271:ASP:CG	2.32	0.50
1:A:258:MET:O	1:A:261:VAL:HG12	2.12	0.49
1:A:76:LYS:NZ	1:A:148:GLU:OE2	2.44	0.49
1:B:261:VAL:CG2	1:B:262:PRO:HD2	2.41	0.49
1:B:359:LEU:HD12	1:B:359:LEU:H	1.76	0.49
1:B:244:ARG:HG3	1:B:271:ASP:CG	2.32	0.49
1:A:392:LEU:HD21	1:A:394:ILE:HD11	1.95	0.49
1:B:359:LEU:HD12	1:B:359:LEU:N	2.28	0.49
1:A:291:ALA:C	1:A:293:LYS:H	2.16	0.49
1:B:293:LYS:HD2	1:B:293:LYS:N	2.28	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:45:VAL:HG21	1:A:141:LEU:HD12	1.95	0.48
1:B:256:ASN:O	1:B:259:ARG:HG3	2.12	0.48
1:A:16:PRO:HG2	1:A:17:ARG:HE	1.78	0.48
1:A:201:LEU:HD22	1:A:240:ILE:CD1	2.41	0.48
1:A:21:TYR:CD2	1:A:65:VAL:HG22	2.48	0.48
1:B:321:LYS:HE2	1:B:391:TYR:CG	2.48	0.48
1:A:261:VAL:HG22	1:A:263:GLU:H	1.78	0.48
1:A:250:ALA:O	1:A:277:THR:HA	2.13	0.48
1:A:221:ARG:O	1:A:221:ARG:HG3	2.13	0.47
1:B:175:VAL:O	1:B:175:VAL:HG23	2.14	0.47
1:B:259:ARG:CD	1:B:259:ARG:C	2.83	0.47
1:A:261:VAL:CG2	1:A:262:PRO:HD2	2.43	0.47
1:B:89:ALA:HB3	1:B:92:GLU:OE1	2.14	0.47
1:A:314:LEU:H	1:A:314:LEU:HD22	1.80	0.47
1:A:346:HIS:HD2	1:A:387:ASP:O	1.97	0.47
1:B:85:ASN:HA	1:B:207:PHE:CE2	2.50	0.46
1:B:175:VAL:CG2	1:B:361:GLU:HG2	2.44	0.46
1:A:260:GLN:NE2	1:A:260:GLN:HA	2.28	0.46
1:A:45:VAL:O	1:A:49:LEU:HD23	2.15	0.46
1:B:259:ARG:NH1	1:B:289:ILE:HD12	2.30	0.46
1:B:284:GLU:HG2	1:B:288:LYS:HE3	1.96	0.46
1:B:321:LYS:HE2	1:B:391:TYR:CD1	2.51	0.46
1:B:315:SER:HB2	2:B:569:HOH:O	2.14	0.45
1:B:236:VAL:HG23	1:B:237:TRP:N	2.31	0.45
1:B:348:GLU:HG2	1:B:349:GLY:H	1.81	0.45
1:B:85:ASN:C	1:B:85:ASN:ND2	2.69	0.45
1:B:111:SER:O	1:B:114:VAL:HG12	2.16	0.45
1:B:246:CYS:HA	1:B:271:ASP:O	2.17	0.45
1:B:301:TYR:HB2	1:B:325:ALA:HB1	1.99	0.45
1:A:66:GLU:HG2	1:A:66:GLU:H	1.48	0.45
1:A:195:ARG:HG2	1:A:198:ASP:OD2	2.17	0.45
1:A:236:VAL:O	1:A:240:ILE:HG12	2.17	0.44
1:B:259:ARG:HH12	1:B:289:ILE:HD12	1.82	0.44
1:B:175:VAL:HG21	1:B:361:GLU:CG	2.44	0.44
1:B:261:VAL:HG22	1:B:263:GLU:H	1.83	0.44
1:B:45:VAL:O	1:B:49:LEU:HD22	2.17	0.44
1:A:196:TYR:O	1:A:197:ARG:HB2	2.16	0.44
1:A:209:VAL:HG13	1:A:210:ALA:N	2.33	0.43
1:A:201:LEU:CD2	1:A:240:ILE:HD11	2.44	0.43
1:A:264:PHE:O	1:A:293:LYS:NZ	2.44	0.43
1:A:93:VAL:O	1:A:97:LEU:HG	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:246:CYS:HA	1:A:271:ASP:O	2.18	0.43
1:B:314:LEU:N	1:B:314:LEU:HD22	2.33	0.43
1:B:138:ALA:O	1:B:142:ARG:HG3	2.18	0.43
1:B:259:ARG:NH1	1:B:289:ILE:CD1	2.81	0.43
1:A:135:ASP:N	1:A:135:ASP:OD2	2.52	0.43
1:A:389:GLU:HB2	1:A:391:TYR:HD1	1.84	0.43
1:A:3:ASN:HB3	1:A:6:TRP:CG	2.53	0.43
1:A:85:ASN:HA	1:A:207:PHE:CE2	2.54	0.43
1:B:23:GLU:OE1	1:B:26:THR:HG21	2.18	0.43
1:A:85:ASN:C	1:A:85:ASN:ND2	2.72	0.43
1:A:71:PHE:C	1:A:71:PHE:CD1	2.92	0.42
1:A:23:GLU:HG3	1:A:227:SER:HB3	2.01	0.42
1:A:115:ILE:O	1:A:119:ARG:HG3	2.20	0.42
1:A:28:VAL:HG11	1:A:65:VAL:CG2	2.49	0.42
1:A:41:ARG:HA	1:A:145:ALA:O	2.19	0.42
1:A:45:VAL:O	1:A:49:LEU:CD2	2.67	0.42
1:B:142:ARG:NH2	2:B:450:HOH:O	2.47	0.42
1:B:237:TRP:O	1:B:241:VAL:HG23	2.20	0.42
1:B:335:ALA:HB1	1:B:344:ARG:O	2.20	0.42
1:B:346:HIS:HD2	1:B:387:ASP:O	2.02	0.42
1:B:296:GLU:HB3	1:B:315:SER:CB	2.49	0.42
1:B:85:ASN:HD21	1:B:87:ARG:HB2	1.84	0.42
1:A:282:MET:HG2	1:A:299:GLN:NE2	2.34	0.42
1:B:287:ILE:HD13	1:B:318:ALA:HB1	2.02	0.42
1:A:204:LEU:HD21	1:A:250:ALA:HA	2.02	0.41
1:A:23:GLU:HB3	1:A:26:THR:CG2	2.46	0.41
1:B:85:ASN:HA	1:B:207:PHE:CD2	2.55	0.41
1:B:314:LEU:HD22	1:B:314:LEU:H	1.84	0.41
1:B:92:GLU:H	1:B:92:GLU:CD	2.23	0.41
1:A:321:LYS:HE2	1:A:391:TYR:CG	2.55	0.41
1:B:85:ASN:HD22	1:B:87:ARG:H	1.65	0.41
1:B:99:ASP:OD2	1:B:365:ARG:NH1	2.49	0.41
1:A:175:VAL:O	1:A:175:VAL:HG23	2.20	0.41
1:A:175:VAL:HG21	1:A:361:GLU:HG2	2.02	0.41
1:A:260:GLN:CA	1:A:260:GLN:HE21	2.27	0.41
1:A:282:MET:HB2	1:A:287:ILE:HD11	2.02	0.41
1:B:137:LEU:HG	1:B:141:LEU:HD22	2.02	0.41
1:B:3:ASN:HB3	1:B:6:TRP:CG	2.56	0.41
1:B:203:PRO:HD2	1:B:204:LEU:HD22	2.02	0.41
1:B:267:LEU:HD13	1:B:268:ASP:N	2.36	0.41
1:B:312:LEU:HB2	1:B:329:THR:HG22	2.04	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:155:GLY:HA2	2:B:492:HOH:O	2.22	0.40
1:B:175:VAL:CG2	1:B:175:VAL:O	2.70	0.40
1:A:58:ALA:HB3	1:A:105:VAL:HG22	2.03	0.40
1:B:115:ILE:O	1:B:119:ARG:HG3	2.22	0.40
1:B:232:ASP:OD2	1:B:234:THR:HB	2.21	0.40
1:A:203:PRO:HG2	1:A:204:LEU:HD22	2.02	0.40
1:B:71:PHE:CD1	1:B:71:PHE:C	2.94	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:122:ALA:O	1:A:122:ALA:O[2_554]	2.11	0.09

## 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	388/396 (98%)	370 (95%)	15 (4%)	3 (1%)	21	16
1	B	388/396 (98%)	373 (96%)	14 (4%)	1 (0%)	43	43
All	All	776/792 (98%)	743 (96%)	29 (4%)	4 (0%)	31	27

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	280	ALA
1	A	203	PRO
1	B	203	PRO
1	A	281	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	312/314 (99%)	289 (93%)	23 (7%)	15	11
1	B	312/314 (99%)	293 (94%)	19 (6%)	20	17
All	All	624/628 (99%)	582 (93%)	42 (7%)	18	15

All (42) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	8	LEU
1	A	27	ASP
1	A	38	LEU
1	A	65	VAL
1	A	66	GLU
1	A	67	PHE
1	A	71	PHE
1	A	77	LEU
1	A	85	ASN
1	A	141	LEU
1	A	147	ASP
1	A	152	GLU
1	A	156	ASP
1	A	159	LEU
1	A	195	ARG
1	A	202	LEU
1	A	203	PRO
1	A	244	ARG
1	A	259	ARG
1	A	266	GLU
1	A	312	LEU
1	A	359	LEU
1	A	365	ARG
1	B	8	LEU
1	B	27	ASP
1	B	38	LEU
1	B	49	LEU

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Mol	Chain	Res	Type
1	B	67	PHE
1	B	71	PHE
1	B	77	LEU
1	B	85	ASN
1	B	141	LEU
1	B	156	ASP
1	B	159	LEU
1	B	195	ARG
1	B	202	LEU
1	B	244	ARG
1	B	259	ARG
1	B	293	LYS
1	B	299	GLN
1	B	312	LEU
1	B	359	LEU

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	85	ASN
1	A	230	GLN
1	A	260	GLN
1	A	299	GLN
1	A	346	HIS
1	B	85	ASN
1	B	299	GLN
1	B	346	HIS

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

## 5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

## 5.6 Ligand geometry

There are no ligands in this entry.

## 5.7 Other polymers

There are no such residues in this entry.

## 5.8 Polymer linkage issues

There are no chain breaks in this entry.

## 6 Fit of model and data [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	392/396 (98%)	-0.41	2 (0%) 90 92	21, 35, 56, 65	0
1	B	392/396 (98%)	-0.41	0 100 100	21, 34, 54, 64	0
All	All	784/792 (98%)	-0.41	2 (0%) 93 94	21, 35, 55, 65	0

All (2) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	165	SER	2.4
1	A	339	ASP	2.2

### 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](#)

There are no ligands in this entry.

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

There are no such residues in this entry.