



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 11:58 AM UTC

PDB ID : 3M1L / pdb_00003m1l
Title : Crystal structure of a C-terminal truncated mutant of a putative ketoacyl reductase (FabG4) from Mycobacterium tuberculosis H37Rv at 2.5 Angstrom resolution
Authors : Dutta, D.; Bhattacharyya, S.; Saha, B.; Das, A.K.
Deposited on : 2010-03-05
Resolution : 2.52 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
Mogul	:	2022.3.0, CSD as543be (2022)
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

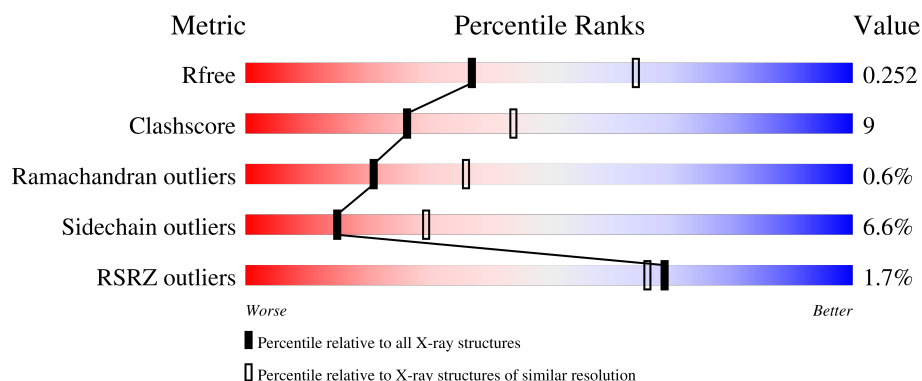
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.52 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	7383 (2.54-2.50)
Clashscore	190562	8079 (2.54-2.50)
Ramachandran outliers	187476	7944 (2.54-2.50)
Sidechain outliers	187428	7946 (2.54-2.50)
RSRZ outliers	180081	7387 (2.54-2.50)

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	432	
1	B	432	

2 Entry composition [i](#)

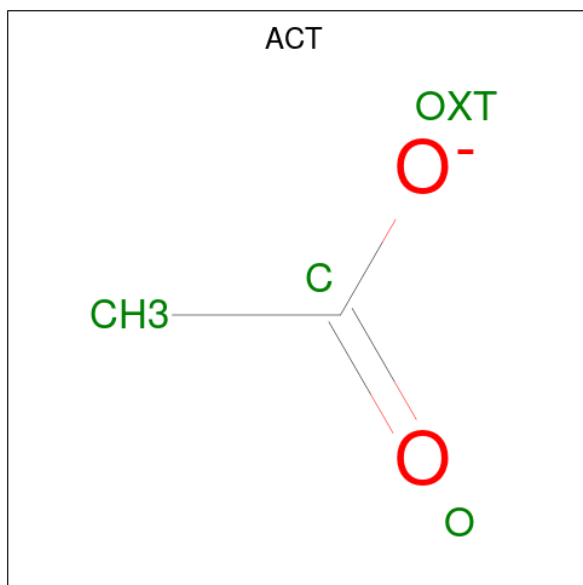
There are 3 unique types of molecules in this entry. The entry contains 5496 atoms, of which 0 are hydrogens and 0 are deuteriums.

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-oxoacyl-(Acyl-carrier-protein) reductase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	364	Total	C	N	O	S	0	2	0
			2652	1663	473	512	4			
1	B	362	Total	C	N	O	S	0	1	0
			2633	1654	467	507	5			

- Molecule 2 is ACETATE ION (CCD ID: ACT) (formula: C₂H₃O₂).



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

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	119	Total	O	0	0
			119	119		

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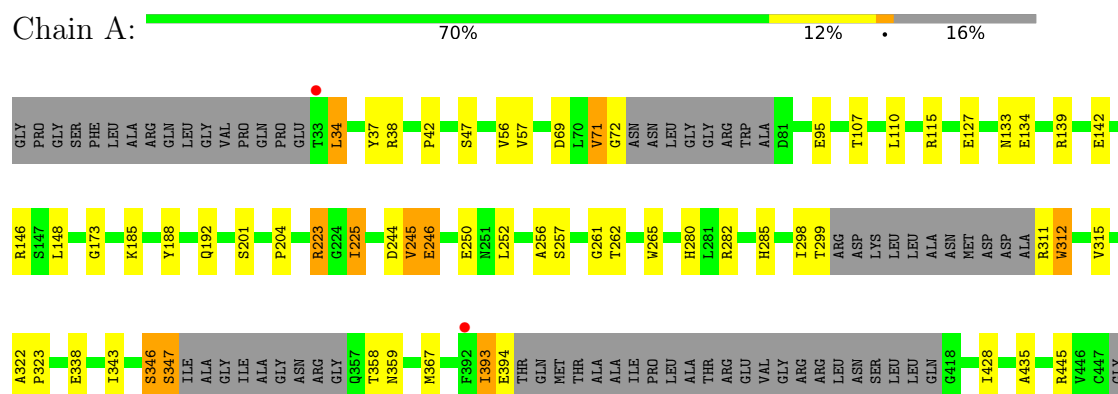
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	88	Total	O	0	0
			88	88		

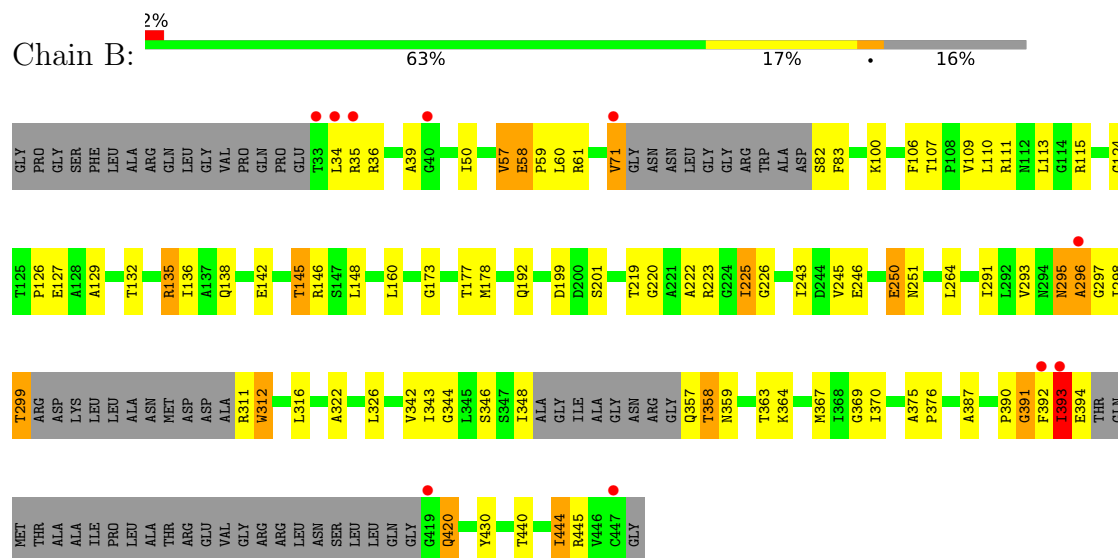
3 Residue-property plots

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: 3-oxoacyl-(Acyl-carrier-protein) reductase



- Molecule 1: 3-oxoacyl-(Acyl-carrier-protein) reductase



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	65.08Å 80.50Å 153.95Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.95 – 2.52 19.95 – 2.52	Depositor EDS
% Data completeness (in resolution range)	98.9 (19.95-2.52) 99.0 (19.95-2.52)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.24 (at 2.53Å)	Xtriage
Refinement program	REFMAC 5.5.0095	Depositor
R, R_{free}	0.181 , 0.241 0.195 , 0.252	Depositor DCC
R_{free} test set	1138 reflections (4.10%)	wwPDB-VP
Wilson B-factor (Å ²)	37.5	Xtriage
Anisotropy	0.552	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 37.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	5496	wwPDB-VP
Average B, all atoms (Å ²)	37.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.30% 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

5.1 Standard geometry

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

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	1.06	4/2690 (0.1%)	1.10	7/3653 (0.2%)
1	B	1.14	12/2671 (0.4%)	1.12	11/3628 (0.3%)
All	All	1.10	16/5361 (0.3%)	1.11	18/7281 (0.2%)

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

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	2
1	B	0	1
All	All	0	3

All (16) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	393	ILE	CA-CB	7.10	1.64	1.54
1	B	293	VAL	C-O	-6.07	1.17	1.24
1	B	297	GLY	CA-C	-6.00	1.48	1.51
1	B	58	GLU	N-CA	5.93	1.50	1.46
1	A	42	PRO	CA-C	5.80	1.56	1.52
1	B	295	ASN	C-O	-5.80	1.17	1.23
1	B	295	ASN	CA-C	-5.60	1.45	1.52
1	B	132	THR	C-O	-5.57	1.17	1.24
1	B	136	ILE	C-O	-5.55	1.17	1.24
1	B	129	ALA	CA-CB	-5.45	1.45	1.53
1	A	393	ILE	CA-CB	5.41	1.62	1.54
1	B	135	ARG	C-O	-5.26	1.18	1.24
1	B	296	ALA	CA-C	-5.25	1.46	1.52
1	B	291	ILE	C-O	-5.19	1.18	1.24

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	133	ASN	C-O	-5.16	1.18	1.24
1	A	134	GLU	C-O	-5.02	1.18	1.24

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	57	VAL	N-CA-C	7.63	118.37	110.36
1	B	297	GLY	N-CA-C	7.03	116.61	110.21
1	A	312	TRP	N-CA-C	6.38	118.23	111.28
1	B	296	ALA	N-CA-C	-6.19	98.19	108.34
1	B	109	VAL	N-CA-C	5.87	118.83	112.96
1	B	322	ALA	CA-C-N	-5.74	113.71	119.56
1	B	322	ALA	C-N-CA	-5.74	113.71	119.56
1	A	34	LEU	N-CA-C	5.68	118.51	110.50
1	B	106	PHE	N-CA-C	5.65	119.35	112.23
1	B	420	GLN	N-CA-C	5.55	117.22	108.23
1	B	107	THR	CA-C-N	-5.45	114.06	119.56
1	B	107	THR	C-N-CA	-5.45	114.06	119.56
1	B	295	ASN	N-CA-C	-5.39	99.15	108.69
1	B	57	VAL	N-CA-C	5.38	116.01	110.36
1	A	225	ILE	CB-CA-C	-5.31	105.08	112.04
1	A	107	THR	CA-C-N	-5.29	114.37	119.87
1	A	107	THR	C-N-CA	-5.29	114.37	119.87
1	A	285	HIS	N-CA-C	5.07	116.95	110.61

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	346	SER	Peptide
1	A	38	ARG	Peptide
1	B	296	ALA	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	2652	0	2672	40	0
1	B	2633	0	2659	59	0
2	A	4	0	3	0	0
3	A	119	0	0	5	0
3	B	88	0	0	6	0
All	All	5496	0	5334	99	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 9.

All (99) 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:223:ARG:HG3	1:A:223:ARG:HH11	1.07	1.17
1:A:282[B]:ARG:HH11	1:A:282[B]:ARG:HG2	0.98	1.13
1:B:393:ILE:HG23	1:B:394:GLU:H	1.18	1.06
1:A:367:MET:HG3	3:A:551:HOH:O	1.63	0.98
1:A:282[B]:ARG:HG2	1:A:282[B]:ARG:NH1	1.81	0.91
1:A:282[B]:ARG:HH11	1:A:282[B]:ARG:CG	1.80	0.90
1:B:124:GLY:H	1:B:138:GLN:NE2	1.71	0.89
1:B:219:THR:OG1	1:B:295:ASN:O	1.92	0.87
1:B:57:VAL:O	1:B:61:ARG:HG3	1.74	0.86
3:A:554:HOH:O	1:B:358:THR:HG21	1.79	0.83
1:B:223:ARG:NH1	1:B:251:ASN:HD22	1.77	0.81
1:B:393:ILE:CG2	1:B:394:GLU:H	1.91	0.81
1:A:223:ARG:HG3	1:A:223:ARG:NH1	1.88	0.81
1:B:178:MET:HA	1:B:178:MET:HE2	1.66	0.74
1:B:393:ILE:HG23	1:B:394:GLU:N	1.98	0.73
1:B:348:ILE:HG23	1:B:391:GLY:HA3	1.69	0.72
1:B:344:GLY:HA3	1:B:367[B]:MET:HE3	1.70	0.72
1:B:364:LYS:HE2	3:B:453:HOH:O	1.92	0.69
1:A:223:ARG:HH11	1:A:223:ARG:CG	1.96	0.69
1:B:124:GLY:H	1:B:138:GLN:HE21	1.39	0.69
1:A:282[B]:ARG:NH1	1:A:282[B]:ARG:CG	2.49	0.66
1:B:245:VAL:HG13	3:B:457:HOH:O	1.95	0.65
1:B:348:ILE:HD13	1:B:393:ILE:HG22	1.82	0.62
1:B:346:SER:OG	1:B:367[A]:MET:CE	2.49	0.61
1:A:346:SER:HB3	1:A:367:MET:CE	2.30	0.61
1:B:346:SER:OG	1:B:367[A]:MET:HE1	2.00	0.61
1:A:115:ARG:HD2	3:A:485:HOH:O	2.01	0.60
1:A:315:VAL:HG11	1:A:359:ASN:HB3	1.84	0.59
1:B:245:VAL:CG1	3:B:462:HOH:O	2.50	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:346:SER:O	1:A:347:SER:HB2	2.04	0.58
1:A:282[A]:ARG:HG3	3:A:553:HOH:O	2.05	0.57
1:A:110:LEU:HD11	1:A:148:LEU:HD11	1.87	0.55
1:B:124:GLY:N	1:B:138:GLN:NE2	2.50	0.55
1:B:192:GLN:NE2	1:B:445:ARG:H	2.04	0.55
1:B:124:GLY:N	1:B:138:GLN:HE21	2.03	0.54
1:B:145:THR:HG21	1:B:160:LEU:HB2	1.90	0.54
1:B:220:GLY:O	1:B:226:GLY:HA3	2.08	0.53
1:B:393:ILE:CG2	1:B:394:GLU:N	2.65	0.53
1:B:173:GLY:HA2	1:B:201:SER:OG	2.09	0.52
1:A:71:VAL:CG1	1:A:72:GLY:N	2.71	0.52
1:A:322:ALA:N	1:A:323:PRO:HD2	2.24	0.52
1:B:223:ARG:HH11	1:B:251:ASN:HD22	1.55	0.52
1:B:178:MET:HE2	1:B:178:MET:CA	2.36	0.52
1:B:192:GLN:HB3	1:B:444:ILE:HD12	1.91	0.52
1:A:244:ASP:OD1	1:A:299:THR:HG21	2.10	0.51
1:B:348:ILE:HD13	1:B:393:ILE:CG2	2.41	0.51
1:A:262:THR:OG1	1:A:280:HIS:HE1	1.94	0.51
1:B:127:GLU:H	1:B:127:GLU:CD	2.20	0.50
1:B:60:LEU:HD21	1:B:178:MET:HG3	1.92	0.50
1:B:50:ILE:O	1:B:61:ARG:NH1	2.44	0.50
1:B:225:ILE:HG12	3:B:460:HOH:O	2.11	0.49
1:B:36:ARG:HH21	1:B:420:GLN:CG	2.25	0.49
1:A:244:ASP:CG	1:A:299:THR:HG21	2.36	0.49
1:B:367[A]:MET:HE3	1:B:387:ALA:HB1	1.94	0.49
1:A:127:GLU:H	1:A:127:GLU:CD	2.20	0.49
1:B:83:PHE:HB2	1:B:113:LEU:HA	1.95	0.48
1:B:222:ALA:HB3	1:B:299:THR:HG21	1.96	0.48
1:B:375:ALA:HB3	1:B:376:PRO:HD3	1.94	0.48
1:B:142:GLU:OE1	1:B:146:ARG:NH1	2.47	0.48
1:B:110:LEU:HD11	1:B:148:LEU:HD11	1.94	0.48
1:B:393:ILE:O	1:B:394:GLU:HG2	2.13	0.47
1:B:342:VAL:C	1:B:343:ILE:HG13	2.38	0.47
1:A:47:SER:HB3	1:A:69:ASP:OD2	2.15	0.47
1:A:246:GLU:HG3	1:A:265:TRP:HB3	1.97	0.47
1:A:34:LEU:HD13	1:A:188:TYR:OH	2.14	0.47
1:B:126:PRO:O	1:B:135:ARG:HG3	2.15	0.46
1:B:344:GLY:HA3	1:B:367[B]:MET:CE	2.44	0.46
1:B:444:ILE:HD13	1:B:444:ILE:HA	1.68	0.45
1:B:61:ARG:HH22	1:B:71:VAL:HG12	1.81	0.45
1:A:139:ARG:HD2	1:A:139:ARG:O	2.17	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:58:GLU:HB2	1:B:59:PRO:CD	2.47	0.45
1:A:315:VAL:HG21	1:A:359:ASN:HD22	1.83	0.44
1:A:173:GLY:HA2	1:A:201:SER:OG	2.18	0.44
1:A:256:ALA:O	1:A:261:GLY:N	2.49	0.44
1:B:357:GLN:HA	3:B:521:HOH:O	2.18	0.44
1:A:146:ARG:HH21	1:A:445:ARG:HH22	1.65	0.43
1:B:243:ILE:HG12	1:B:264:LEU:HB3	2.01	0.43
1:B:250:GLU:H	1:B:250:GLU:HG3	1.50	0.43
1:A:311:ARG:HD3	1:A:312:TRP:N	2.33	0.43
1:B:312:TRP:CZ3	1:B:316:LEU:HD11	2.53	0.43
1:A:204:PRO:HD3	1:A:435:ALA:HB2	2.01	0.43
1:A:311:ARG:HD3	1:A:312:TRP:H	1.84	0.43
1:B:312:TRP:HZ3	1:B:316:LEU:HD11	1.84	0.43
1:B:359:ASN:O	1:B:363:THR:HG23	2.19	0.43
1:A:37:TYR:CD1	1:A:185:LYS:HE3	2.55	0.42
1:A:142:GLU:OE1	1:A:146:ARG:NH1	2.53	0.42
1:B:111:ARG:NH1	1:B:111:ARG:HB2	2.34	0.42
1:B:177:THR:HA	1:B:430:TYR:CE1	2.55	0.41
1:B:58:GLU:HB2	1:B:59:PRO:HD3	2.02	0.41
1:A:244:ASP:HB2	1:A:252:LEU:HD13	2.02	0.41
1:B:369:GLY:O	1:B:370:ILE:C	2.64	0.41
1:A:71:VAL:HG13	1:A:72:GLY:H	1.86	0.40
1:A:244:ASP:OD2	1:A:299:THR:HG21	2.21	0.40
1:B:199:ASP:HB2	3:B:454:HOH:O	2.21	0.40
1:A:245:VAL:HG13	3:A:529:HOH:O	2.21	0.40
1:A:343:ILE:HG21	1:A:428:ILE:HG23	2.04	0.40
1:B:316:LEU:HD23	1:B:363:THR:HG22	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	356/432 (82%)	343 (96%)	13 (4%)	0	100	100
1	B	353/432 (82%)	338 (96%)	11 (3%)	4 (1%)	11	21
All	All	709/864 (82%)	681 (96%)	24 (3%)	4 (1%)	21	36

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	391	GLY
1	B	393	ILE
1	B	39	ALA
1	B	390	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	266/313 (85%)	251 (94%)	15 (6%)	19	37
1	B	265/313 (85%)	245 (92%)	20 (8%)	12	24
All	All	531/626 (85%)	496 (93%)	35 (7%)	15	30

All (35) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	56	VAL
1	A	71	VAL
1	A	95	GLU
1	A	223	ARG
1	A	225	ILE
1	A	245	VAL
1	A	246	GLU
1	A	250	GLU
1	A	257	SER
1	A	298	ILE
1	A	338	GLU
1	A	347	SER

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Mol	Chain	Res	Type
1	A	358	THR
1	A	393	ILE
1	A	394	GLU
1	B	34	LEU
1	B	35	ARG
1	B	71	VAL
1	B	82	SER
1	B	100	LYS
1	B	115	ARG
1	B	145	THR
1	B	225	ILE
1	B	246	GLU
1	B	250	GLU
1	B	298	ILE
1	B	299	THR
1	B	311	ARG
1	B	312	TRP
1	B	326	LEU
1	B	358	THR
1	B	392	PHE
1	B	393	ILE
1	B	440	THR
1	B	444	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	192	GLN
1	A	251	ASN
1	A	280	HIS
1	A	319	ASN
1	A	372	GLN
1	B	138	GLN
1	B	192	GLN
1	B	239	HIS
1	B	251	ASN
1	B	280	HIS

5.3.3 RNA ⓘ

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](#)

1 ligand is modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	ACT	A	1	-	3,3,3	0.85	0	3,3,3	1.24	0

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 [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	364/432 (84%)	-0.43	2 (0%) 87 85	13, 31, 59, 83	2 (0%)
1	B	362/432 (83%)	-0.31	10 (2%) 55 51	13, 37, 64, 111	1 (0%)
All	All	726/864 (84%)	-0.37	12 (1%) 69 66	13, 34, 62, 111	3 (0%)

All (12) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	392	PHE	3.5
1	B	33	THR	3.4
1	B	34	LEU	2.8
1	B	296	ALA	2.8
1	B	393	ILE	2.7
1	B	419	GLY	2.5
1	A	392	PHE	2.5
1	B	71	VAL	2.3
1	B	40	GLY	2.2
1	B	35	ARG	2.2
1	A	33	THR	2.2
1	B	447	CYS	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 oligosaccharides in this entry.

6.4 Ligands

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	ACT	A	1	4/4	0.88	0.14	38,39,39,39	0

6.5 Other polymers

There are no such residues in this entry.