



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 7, 2026 – 02:11 AM UTC

PDB ID : 4TRM / pdb_00004trm
Title : Structure of the apo form of InhA from Mycobacterium tuberculosis
Authors : Chollet, A.; Julien, S.; Mourey, L.; Maveyraud, L.
Deposited on : 2014-06-17
Resolution : 1.80 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

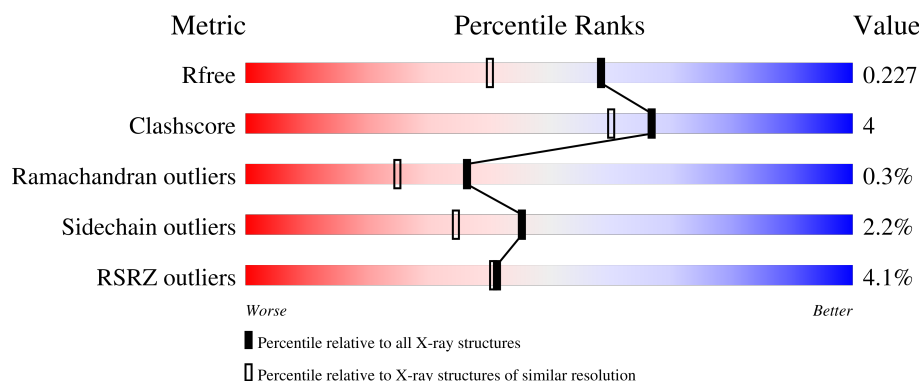
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.80 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	7662 (1.80-1.80)
Clashscore	190562	8479 (1.80-1.80)
Ramachandran outliers	187476	8391 (1.80-1.80)
Sidechain outliers	187428	8390 (1.80-1.80)
RSRZ outliers	180081	7663 (1.80-1.80)

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	269	3% 84% 9% 5%
1	B	269	2% 88% 8% .
1	C	269	4% 86% 9% 5%
1	D	269	3% 85% 7% . 7%
1	E	269	4% 87% 10% ..

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Mol	Chain	Length	Quality of chain
1	F	269	6% 84% 9% 7%

2 Entry composition [i](#)

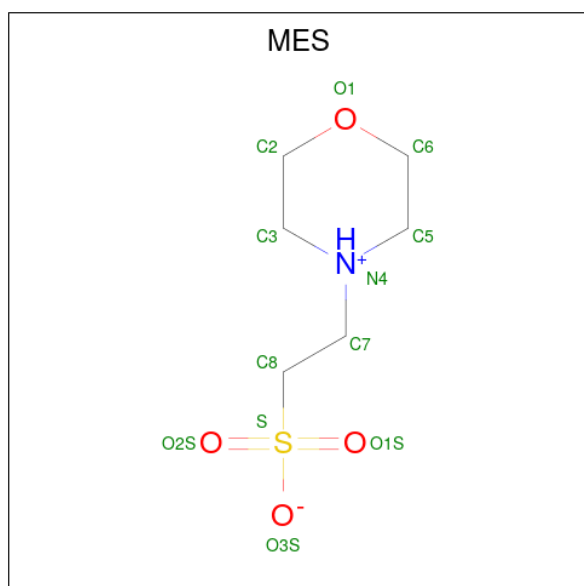
There are 3 unique types of molecules in this entry. The entry contains 12105 atoms, of which 0 are hydrogens and 0 are deuteriums.

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	256	Total	C	N	O	S	0	5	0
			1930	1227	334	360	9			
1	B	259	Total	C	N	O	S	0	7	0
			1926	1218	332	367	9			
1	C	256	Total	C	N	O	S	0	4	0
			1870	1185	325	351	9			
1	D	249	Total	C	N	O	S	0	0	0
			1790	1138	308	335	9			
1	E	266	Total	C	N	O	S	0	5	0
			1973	1253	342	369	9			
1	F	251	Total	C	N	O	S	0	2	0
			1810	1150	315	335	10			

- Molecule 2 is 2-(N-MORPHOLINO)-ETHANESULFONIC ACID (CCD ID: MES) (formula: $C_6H_{13}NO_4S$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	S	0	0
			12	6	1	4	1		

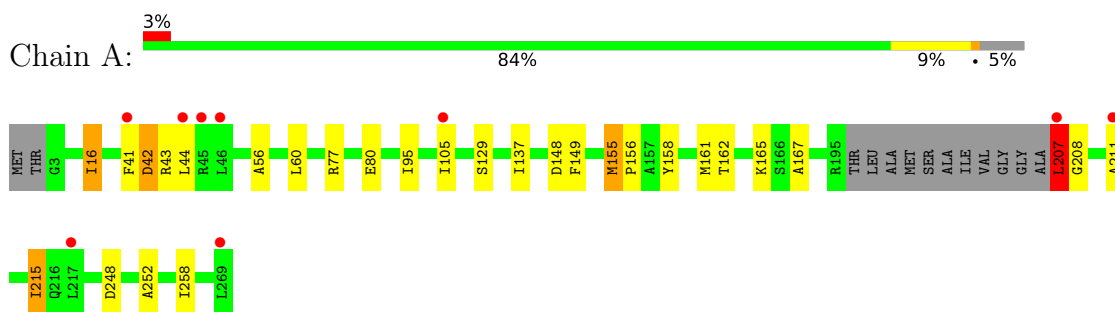
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	198	Total	O	0	0
			198	198		
3	B	177	Total	O	0	0
			177	177		
3	C	110	Total	O	0	0
			110	110		
3	D	99	Total	O	0	0
			99	99		
3	E	125	Total	O	0	0
			125	125		
3	F	85	Total	O	0	0
			85	85		

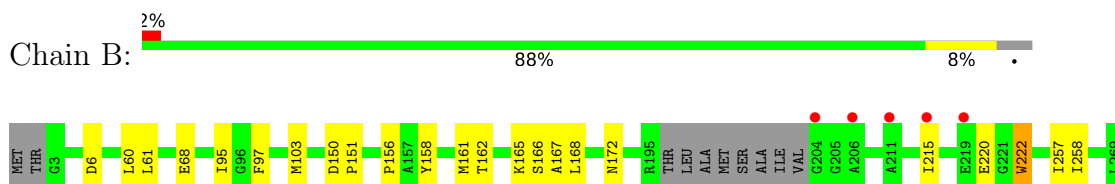
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

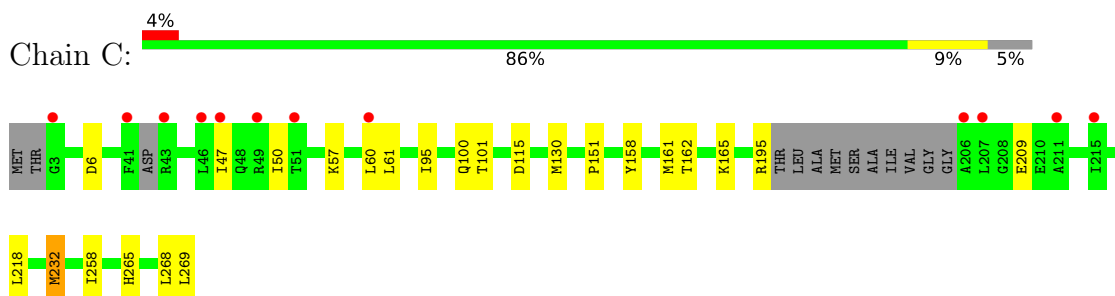
- Molecule 1: Enoyl-[acyl-carrier-protein] reductase [NADH]



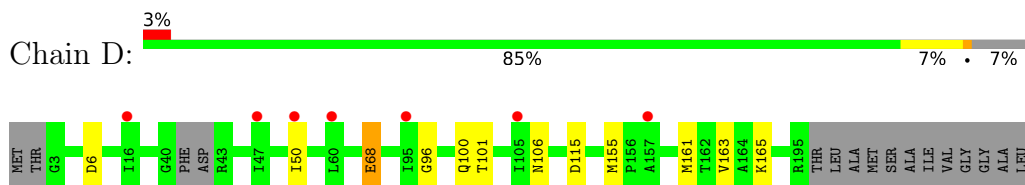
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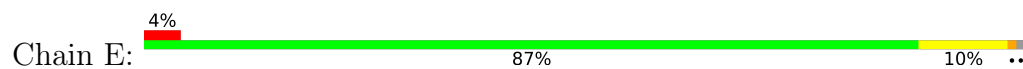


- Molecule 1: Enoyl-[acyl-carrier-protein] reductase [NADH]

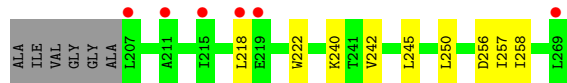
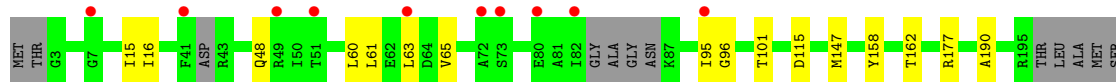
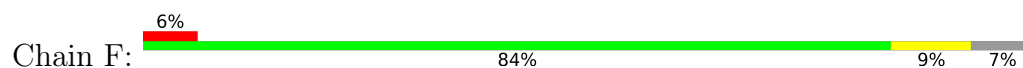




- Molecule 1: Enoyl-[acyl-carrier-protein] reductase [NADH]



- Molecule 1: Enoyl-[acyl-carrier-protein] reductase [NADH]



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	100.10Å 81.79Å 187.24Å 90.00° 95.62° 90.00°	Depositor
Resolution (Å)	35.00 – 1.80 35.00 – 1.80	Depositor EDS
% Data completeness (in resolution range)	97.9 (35.00-1.80) 98.0 (35.00-1.80)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.06	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.01 (at 1.79Å)	Xtriage
Refinement program	REFMAC 5.8.0049	Depositor
R, R_{free}	0.189 , 0.219 0.196 , 0.227	Depositor DCC
R_{free} test set	6889 reflections (4.94%)	wwPDB-VP
Wilson B-factor (Å ²)	29.0	Xtriage
Anisotropy	0.111	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 61.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.27$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	12105	wwPDB-VP
Average B, all atoms (Å ²)	41.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.96% 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: MES

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	1.42	8/1967 (0.4%)	1.13	4/2671 (0.1%)
1	B	1.36	8/1962 (0.4%)	1.08	3/2666 (0.1%)
1	C	1.07	0/1904	0.96	2/2589 (0.1%)
1	D	1.13	1/1825 (0.1%)	0.99	3/2486 (0.1%)
1	E	1.17	4/2011 (0.2%)	1.05	8/2736 (0.3%)
1	F	1.22	3/1844 (0.2%)	1.04	3/2509 (0.1%)
All	All	1.24	24/11513 (0.2%)	1.04	23/15657 (0.1%)

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	E	0	1

All (24) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	E	49	ARG	CZ-NH2	-8.32	1.22	1.33
1	F	222	TRP	N-CA	-8.19	1.36	1.46
1	B	222	TRP	N-CA	-7.83	1.37	1.46
1	B	168	LEU	N-CA	7.10	1.54	1.46
1	E	166	SER	N-CA	6.68	1.54	1.46
1	A	149	PHE	N-CA	-6.46	1.38	1.46
1	A	137	ILE	N-CA	6.22	1.54	1.46
1	A	148	ASP	N-CA	6.12	1.54	1.45
1	B	156	PRO	C-O	5.96	1.30	1.23
1	A	167	ALA	N-CA	-5.90	1.39	1.46
1	F	15	ILE	N-CA	5.80	1.52	1.46
1	E	49	ARG	CD-NE	5.67	1.54	1.46

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	166	SER	CA-CB	5.45	1.61	1.53
1	D	163	VAL	C-O	5.42	1.30	1.24
1	A	129	SER	CB-OG	-5.41	1.31	1.42
1	B	172	ASN	N-CA	5.34	1.52	1.46
1	B	103	MET	N-CA	-5.29	1.40	1.46
1	F	190	ALA	CA-C	5.27	1.59	1.52
1	A	56	ALA	N-CA	5.25	1.52	1.45
1	A	155	MET	C-O	-5.21	1.19	1.24
1	E	16	ILE	C-O	-5.18	1.17	1.23
1	B	97	PHE	CA-C	5.13	1.59	1.52
1	A	252	ALA	C-O	5.01	1.30	1.24
1	B	167	ALA	N-CA	-5.01	1.40	1.46

All (23) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	95	ILE	N-CA-C	9.21	121.64	111.88
1	E	95[A]	ILE	N-CA-C	-7.77	97.23	108.11
1	E	95[B]	ILE	N-CA-C	-7.77	97.23	108.11
1	B	95[A]	ILE	N-CA-C	-6.96	97.60	107.75
1	B	95[B]	ILE	N-CA-C	-6.96	97.60	107.75
1	A	16	ILE	CB-CA-C	6.82	117.59	110.91
1	F	95	ILE	N-CA-C	-6.70	98.73	108.11
1	D	96	GLY	N-CA-C	-6.65	97.42	113.18
1	E	104	GLY	N-CA-C	6.08	123.15	114.64
1	A	95[A]	ILE	N-CA-C	-5.99	99.72	108.11
1	A	95[B]	ILE	N-CA-C	-5.99	99.72	108.11
1	B	150	ASP	O-C-N	5.63	125.86	121.47
1	C	209	GLU	N-CA-C	5.41	116.85	111.07
1	E	16	ILE	CB-CA-C	-5.37	106.34	111.44
1	D	220	GLU	N-CA-C	5.34	117.86	111.71
1	E	209	GLU	N-CA-C	5.34	117.10	111.28
1	F	96	GLY	N-CA-C	-5.34	100.53	113.18
1	F	16	ILE	CB-CA-C	-5.17	106.53	111.44
1	A	207	LEU	CB-CA-C	5.07	119.73	110.10
1	E	155	MET	CB-CA-C	5.06	120.14	110.17
1	D	155	MET	CB-CG-SD	5.05	127.85	112.70
1	E	158[A]	TYR	N-CA-CB	5.04	119.00	110.49
1	E	158[B]	TYR	N-CA-CB	5.04	119.00	110.49

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	E	49	ARG	Sidechain

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	1930	0	1927	17	0
1	B	1926	0	1871	8	0
1	C	1870	0	1809	17	0
1	D	1790	0	1737	11	0
1	E	1973	0	1940	20	0
1	F	1810	0	1744	20	0
2	A	12	0	13	0	0
3	A	198	0	0	5	0
3	B	177	0	0	1	0
3	C	110	0	0	1	0
3	D	99	0	0	0	0
3	E	125	0	0	1	0
3	F	85	0	0	1	0
All	All	12105	0	11041	83	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 4.

All (83) 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:158[A]:TYR:HD2	1:B:162:THR:HG1	0.97	0.96
1:E:158[B]:TYR:HD2	1:E:162:THR:HG1	1.09	0.90
1:E:43:ARG:O	1:E:46:LEU:HB3	1.76	0.84
1:C:101:THR:HG21	1:C:115:ASP:OD2	1.85	0.77
1:D:101:THR:HG21	1:D:115:ASP:OD2	1.86	0.76
1:E:156:PRO:O	1:E:157:ALA:O	2.03	0.76
1:F:101:THR:HG21	1:F:115:ASP:OD2	1.86	0.75
1:A:158:TYR:HD2	1:A:162:THR:HG1	1.33	0.74
1:F:158:TYR:HD2	1:F:162:THR:HG1	1.34	0.73
1:A:207:LEU:HD23	1:A:208:GLY:H	1.52	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:211:ALA:O	1:A:215:ILE:HD12	1.93	0.68
1:F:245:LEU:HD12	1:F:250:LEU:CD1	2.24	0.68
1:F:245:LEU:HD12	1:F:250:LEU:HD12	1.77	0.66
1:C:158:TYR:HD2	1:C:162:THR:HG1	1.45	0.63
1:D:68:GLU:H	1:D:68:GLU:CD	2.04	0.63
1:A:158:TYR:HD2	1:A:162:THR:OG1	1.82	0.62
1:E:156:PRO:HG3	1:F:177:ARG:NH2	2.16	0.61
1:C:195:ARG:CB	1:C:232:MET:HE3	2.31	0.59
1:B:158[A]:TYR:HD2	1:B:162:THR:OG1	1.78	0.58
1:C:158:TYR:HD2	1:C:162:THR:OG1	1.85	0.58
1:E:106:ASN:HB2	1:E:107:PRO:HD2	1.86	0.58
1:A:77:ARG:HA	1:A:80:GLU:HG2	1.86	0.57
1:F:158:TYR:HD2	1:F:162:THR:OG1	1.88	0.57
1:A:207:LEU:CD2	1:A:208:GLY:H	2.20	0.54
1:E:46:LEU:C	1:E:46:LEU:HD23	2.33	0.54
1:D:269:LEU:HD11	1:F:218:LEU:HB2	1.90	0.54
1:C:57:LYS:HA	3:C:371:HOH:O	2.06	0.53
1:A:158:TYR:HE2	3:A:565:HOH:O	1.89	0.53
1:D:269:LEU:HD21	1:F:218:LEU:HD12	1.91	0.52
1:A:80:GLU:HG3	3:A:559:HOH:O	2.08	0.52
1:B:258:ILE:HD12	1:B:258:ILE:N	2.27	0.49
1:C:268[B]:LEU:O	1:C:269[B]:LEU:HB2	2.13	0.49
1:D:258:ILE:HD12	1:D:258:ILE:N	2.26	0.49
1:C:218:LEU:HA	1:E:269:LEU:CD1	2.43	0.48
1:C:268[B]:LEU:C	1:C:269[B]:LEU:O	2.56	0.48
1:F:147[A]:MET:HE1	1:F:242:VAL:CG2	2.44	0.48
1:D:100:GLN:O	1:D:106:ASN:ND2	2.40	0.48
1:C:269[B]:LEU:HD12	1:C:269[B]:LEU:N	2.29	0.47
1:E:156:PRO:C	1:E:157:ALA:O	2.57	0.47
1:D:265:HIS:HE1	1:E:256:ASP:OD2	1.97	0.47
1:C:265:HIS:HE1	1:F:256:ASP:OD2	1.97	0.47
1:A:258[A]:ILE:N	1:A:258[A]:ILE:HD12	2.30	0.46
1:C:258:ILE:N	1:C:258:ILE:HD12	2.30	0.46
1:E:100:GLN:HB3	1:E:105:ILE:CG1	2.46	0.46
1:A:41:PHE:O	1:A:42:ASP:CB	2.64	0.46
1:E:161:MET:HE2	1:E:165:LYS:HE2	1.98	0.46
1:F:147[A]:MET:HE1	1:F:242:VAL:HG22	1.97	0.46
1:C:218:LEU:HA	1:E:269:LEU:HD12	1.99	0.45
1:D:256:ASP:OD2	1:E:265:HIS:HE1	1.99	0.45
1:A:16:ILE:CG1	3:A:507:HOH:O	2.65	0.45
1:B:60:LEU:C	1:B:61:LEU:HD12	2.42	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:257:ILE:C	1:D:258:ILE:HD12	2.42	0.44
1:E:202:ILE:HG21	1:E:215:ILE:HD12	1.99	0.44
1:E:158[A]:TYR:HE2	1:E:218:LEU:HD22	1.82	0.44
1:B:161:MET:HE2	1:B:165:LYS:HE2	1.98	0.44
1:C:60:LEU:C	1:C:61:LEU:HD12	2.42	0.44
1:F:60:LEU:C	1:F:61:LEU:HD12	2.42	0.44
1:A:161:MET:HE2	1:A:165:LYS:HE2	1.99	0.43
1:C:268[B]:LEU:O	1:C:269[B]:LEU:O	2.35	0.43
1:F:240:LYS:HE2	3:F:385:HOH:O	2.18	0.43
1:C:47:ILE:HA	1:C:50:ILE:HD12	2.00	0.43
1:F:258:ILE:HD12	1:F:258:ILE:N	2.32	0.43
1:F:63:LEU:HD21	1:F:65:VAL:HG12	1.99	0.43
1:C:130:MET:HE3	1:C:130:MET:HB2	1.86	0.42
1:F:63:LEU:HD23	1:F:65:VAL:HG13	2.00	0.42
1:A:248:ASP:OD1	3:A:522:HOH:O	2.22	0.42
1:A:207:LEU:CG	1:A:208:GLY:H	2.33	0.42
1:A:155:MET:HB2	1:A:156:PRO:HD2	2.02	0.42
1:D:269:LEU:HD21	1:F:218:LEU:HA	2.01	0.42
1:F:63:LEU:HD21	1:F:65:VAL:CG1	2.50	0.42
1:B:257:ILE:C	1:B:258:ILE:HD12	2.44	0.42
1:F:257:ILE:C	1:F:258:ILE:HD12	2.45	0.42
1:E:156:PRO:HD3	1:F:177:ARG:NH1	2.35	0.41
1:A:158:TYR:CE2	3:A:565:HOH:O	2.56	0.41
1:B:158[A]:TYR:CD2	1:B:162:THR:OG1	2.59	0.41
1:B:222:TRP:HZ2	3:B:459:HOH:O	2.02	0.41
1:E:155:MET:HG3	1:E:158[A]:TYR:HD2	1.85	0.41
1:A:42:ASP:O	1:A:43:ARG:C	2.63	0.41
1:E:142:GLY:HA2	3:E:373:HOH:O	2.20	0.41
1:E:158[A]:TYR:O	1:E:159:ASN:C	2.64	0.40
1:D:161:MET:HE2	1:D:165:LYS:HE2	2.02	0.40
1:E:43:ARG:CB	1:E:46:LEU:HD13	2.51	0.40
1:C:161:MET:HE2	1:C:165:LYS:HE2	2.04	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	257/269 (96%)	249 (97%)	7 (3%)	1 (0%)	30	19
1	B	261/269 (97%)	254 (97%)	6 (2%)	1 (0%)	30	19
1	C	253/269 (94%)	244 (96%)	8 (3%)	1 (0%)	30	19
1	D	243/269 (90%)	236 (97%)	7 (3%)	0	100	100
1	E	267/269 (99%)	252 (94%)	12 (4%)	3 (1%)	11	3
1	F	245/269 (91%)	237 (97%)	8 (3%)	0	100	100
All	All	1526/1614 (94%)	1472 (96%)	48 (3%)	6 (0%)	36	19

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	E	157	ALA
1	E	158[A]	TYR
1	E	158[B]	TYR
1	A	42	ASP
1	B	151	PRO
1	C	151	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	195/205 (95%)	190 (97%)	5 (3%)	40	28
1	B	188/205 (92%)	184 (98%)	4 (2%)	47	36
1	C	180/205 (88%)	177 (98%)	3 (2%)	53	45
1	D	173/205 (84%)	167 (96%)	6 (4%)	32	19
1	E	194/205 (95%)	188 (97%)	6 (3%)	35	23
1	F	172/205 (84%)	171 (99%)	1 (1%)	78	77

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
All	All	1102/1230 (90%)	1077 (98%)	25 (2%)	45 33

All (25) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	44	LEU
1	A	60	LEU
1	A	105	ILE
1	A	207	LEU
1	A	215	ILE
1	B	6	ASP
1	B	68	GLU
1	B	215	ILE
1	B	220	GLU
1	C	6	ASP
1	C	100	GLN
1	C	232	MET
1	D	6	ASP
1	D	50	ILE
1	D	68	GLU
1	D	217	LEU
1	D	232	MET
1	D	269	LEU
1	E	6	ASP
1	E	44	LEU
1	E	95[A]	ILE
1	E	95[B]	ILE
1	E	214	GLN
1	E	217	LEU
1	F	48	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (28) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	35	GLN
1	A	66	GLN
1	A	139	ASN
1	A	265	HIS
1	B	48	GLN
1	B	66	GLN
1	B	139	ASN

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Mol	Chain	Res	Type
1	B	214	GLN
1	B	265	HIS
1	C	24	HIS
1	C	48	GLN
1	C	66	GLN
1	C	121	HIS
1	C	214	GLN
1	C	265	HIS
1	D	24	HIS
1	D	66	GLN
1	D	121	HIS
1	D	265	HIS
1	E	48	GLN
1	E	86	ASN
1	E	214	GLN
1	E	265	HIS
1	F	24	HIS
1	F	48	GLN
1	F	139	ASN
1	F	216	GLN
1	F	265	HIS

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

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	MES	A	301	-	12,12,12	2.29	1 (8%)	15,16,16	2.30	5 (33%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	MES	A	301	-	-	5/6/14/14	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	301	MES	C8-S	-7.49	1.67	1.77

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	301	MES	C6-C5-N4	4.61	117.12	110.12
2	A	301	MES	C2-C3-N4	4.30	116.66	110.12
2	A	301	MES	O2S-S-C8	3.63	112.22	106.73
2	A	301	MES	O1-C6-C5	-3.25	104.77	111.77
2	A	301	MES	C7-N4-C5	2.72	118.48	111.24

There are no chirality outliers.

All (5) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	301	MES	C8-C7-N4-C5
2	A	301	MES	N4-C7-C8-S
2	A	301	MES	C7-C8-S-O1S
2	A	301	MES	C7-C8-S-O3S
2	A	301	MES	C7-C8-S-O2S

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

6.1 Protein, DNA and RNA chains

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2			OWAB(Å ²)	Q<0.9
1	A	256/269 (95%)	-0.10	9 (3%)	47	47	10, 28, 59, 83	5 (1%)
1	B	259/269 (96%)	-0.07	5 (1%)	66	66	10, 30, 55, 67	7 (2%)
1	C	256/269 (95%)	0.37	12 (4%)	36	35	13, 46, 77, 95	4 (1%)
1	D	249/269 (92%)	0.47	9 (3%)	46	46	22, 45, 76, 90	0
1	E	266/269 (98%)	0.31	12 (4%)	38	37	13, 45, 76, 95	5 (1%)
1	F	251/269 (93%)	0.49	16 (6%)	25	24	12, 45, 80, 102	2 (0%)
All	All	1537/1614 (95%)	0.24	63 (4%)	41	41	10, 39, 75, 102	23 (1%)

All (63) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	157	ALA	4.4
1	F	211	ALA	4.2
1	A	41	PHE	4.1
1	F	207	LEU	4.0
1	A	207	LEU	3.9
1	E	46	LEU	3.8
1	F	41	PHE	3.6
1	E	157	ALA	3.5
1	E	44	LEU	3.5
1	F	82	ILE	3.5
1	B	204	GLY	3.4
1	E	47	ILE	3.3
1	A	44	LEU	3.3
1	E	269	LEU	3.3
1	C	41	PHE	3.2
1	E	105	ILE	3.1
1	A	105	ILE	3.1
1	D	50	ILE	2.9
1	F	219	GLU	2.8

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Mol	Chain	Res	Type	RSRZ
1	D	105	ILE	2.8
1	F	215	ILE	2.8
1	F	269	LEU	2.8
1	D	47	ILE	2.7
1	F	49	ARG	2.7
1	E	158[A]	TYR	2.6
1	C	207	LEU	2.6
1	B	206	ALA	2.6
1	C	211	ALA	2.6
1	A	269	LEU	2.6
1	D	60	LEU	2.6
1	D	95	ILE	2.6
1	F	51	THR	2.5
1	E	207	LEU	2.5
1	E	42	ASP	2.5
1	A	211	ALA	2.4
1	A	217	LEU	2.4
1	B	215	ILE	2.4
1	C	47	ILE	2.4
1	C	60	LEU	2.4
1	A	45	ARG	2.4
1	B	211	ALA	2.4
1	F	7	GLY	2.4
1	D	269	LEU	2.3
1	F	73	SER	2.3
1	E	103	MET	2.3
1	B	219	GLU	2.3
1	F	218	LEU	2.3
1	D	218	LEU	2.3
1	F	80	GLU	2.2
1	F	72	ALA	2.2
1	E	41	PHE	2.2
1	E	97	PHE	2.2
1	C	46	LEU	2.2
1	D	16	ILE	2.1
1	C	51	THR	2.1
1	C	49	ARG	2.1
1	F	63	LEU	2.1
1	F	95	ILE	2.1
1	C	215	ILE	2.1
1	A	46	LEU	2.1
1	C	206	ALA	2.1

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Mol	Chain	Res	Type	RSRZ
1	C	3	GLY	2.0
1	C	43	ARG	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 oligosaccharides 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, 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($\text{\AA}^2$)	Q<0.9
2	MES	A	301	12/12	0.73	0.14	46,85,148,149	0

6.5 Other polymers [i](#)

There are no such residues in this entry.