



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 12:23 PM UTC

PDB ID : 3VYO / pdb_00003vyo
Title : Crystal structure of Mycobacterium tuberculosis L,D-transpeptidase LdtMt2 N140 truncation mutant (residues 140-408)
Authors : Li, W.J.; Li, D.F.; Bi, L.J.; Wang, D.C.
Deposited on : 2012-09-30
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

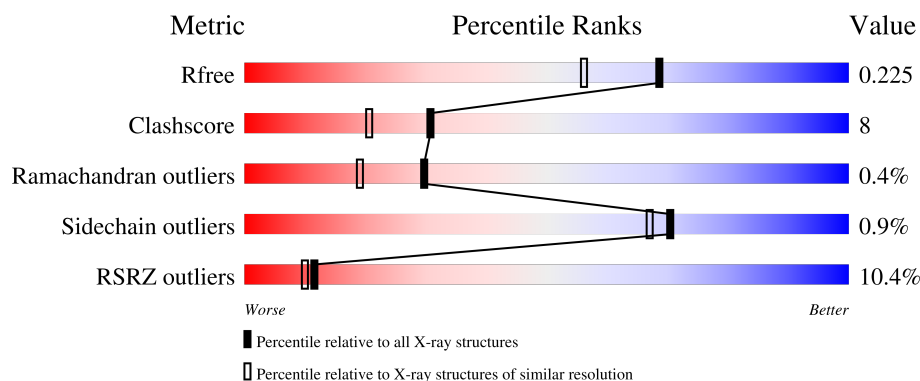
1 Overall quality at a glance

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	9% 82% 14% • •
1	B	269	10% 81% 15% • •

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 4519 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 conserved lipoprotein LPPS.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	261	Total	C	N	O	S	Se	0	0	0
			2003	1262	347	387	1	6			
1	B	262	Total	C	N	O	S	Se	0	0	0
			2004	1262	346	389	1	6			

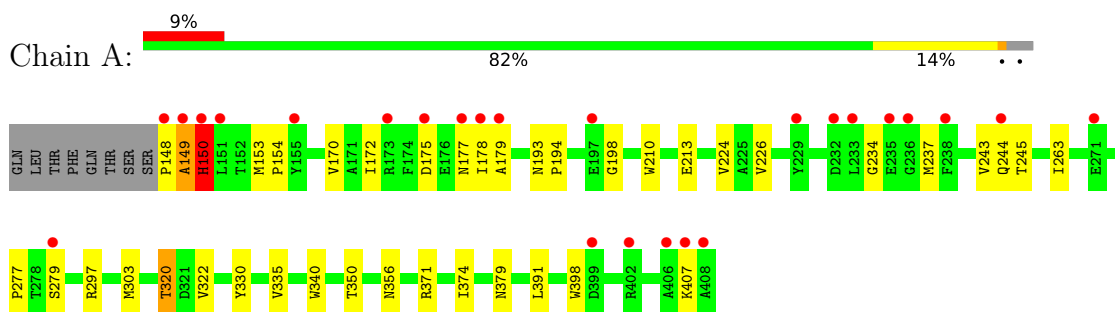
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	264	Total	O	0	0
			264	264		
2	B	248	Total	O	0	0
			248	248		

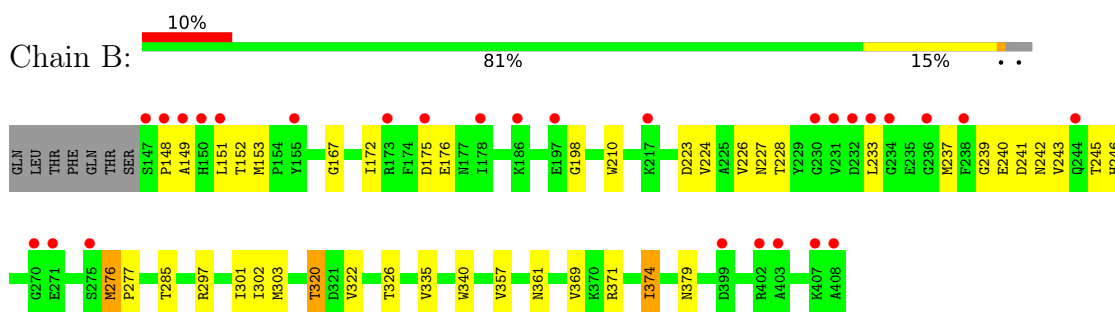
3 Residue-property plots [i](#)

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: Probable conserved lipoprotein LPPS



- Molecule 1: Probable conserved lipoprotein LPPS



4 Data and refinement statistics

Property	Value	Source
Space group	I 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	116.37Å 121.19Å 122.99Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	42.27 – 1.80 42.27 – 1.80	Depositor EDS
% Data completeness (in resolution range)	99.9 (42.27-1.80) 99.9 (42.27-1.80)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	0.11	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.28 (at 1.79Å)	Xtriage
Refinement program	CNS 1.2	Depositor
R, R_{free}	0.209 , 0.225 0.209 , 0.225	Depositor DCC
R_{free} test set	4046 reflections (5.03%)	wwPDB-VP
Wilson B-factor (Å ²)	19.8	Xtriage
Anisotropy	0.684	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 34.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.064 for -h,-l,-k	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	4519	wwPDB-VP
Average B, all atoms (Å ²)	27.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.30% 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

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	0.35	0/2054	0.89	6/2799 (0.2%)
1	B	0.35	0/2054	0.89	6/2800 (0.2%)
All	All	0.35	0/4108	0.89	12/5599 (0.2%)

There are no bond length outliers.

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	335	VAL	N-CA-C	-7.86	96.40	107.80
1	A	335	VAL	N-CA-C	-7.46	96.99	107.80
1	A	320	THR	N-CA-C	6.54	119.57	108.90
1	A	374	ILE	N-CA-C	6.34	117.61	108.48
1	B	320	THR	N-CA-C	6.30	119.21	109.07
1	A	379	ASN	N-CA-C	6.10	120.67	113.23
1	B	379	ASN	N-CA-C	5.97	120.52	113.23
1	B	374	ILE	N-CA-C	5.73	116.73	108.48
1	A	170	VAL	N-CA-C	-5.48	101.65	108.89
1	B	285	THR	CA-C-N	5.20	125.18	120.03
1	B	285	THR	C-N-CA	5.20	125.18	120.03
1	A	150	HIS	N-CA-C	-5.16	106.66	113.17

There are no chirality outliers.

There are no planarity outliers.

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	2003	0	1898	29	0
1	B	2004	0	1897	38	0
2	A	264	0	0	4	3
2	B	248	0	0	7	0
All	All	4519	0	3795	65	3

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 8.

All (65) 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:151:LEU:HB2	2:B:687:HOH:O	1.69	0.93
1:B:176:GLU:HG3	1:B:237:MSE:HE2	1.72	0.72
1:A:279:SER:HB2	2:A:523:HOH:O	1.89	0.70
1:A:224:VAL:HB	1:A:245:THR:HG22	1.77	0.67
1:A:213:GLU:HG3	1:A:398:TRP:CG	2.34	0.63
1:A:350:THR:HG23	2:A:523:HOH:O	1.98	0.62
1:B:152:THR:HB	1:B:237:MSE:HE3	1.81	0.60
1:A:150:HIS:H	1:A:150:HIS:CD2	2.18	0.60
1:A:148:PRO:O	1:A:149:ALA:HB3	2.01	0.59
1:B:176:GLU:HG3	1:B:237:MSE:CE	2.33	0.58
1:B:228:THR:O	1:B:239:GLY:HA3	2.03	0.58
1:B:277:PRO:HD2	2:B:502:HOH:O	2.03	0.58
1:A:226:VAL:HB	1:A:243:VAL:CG2	2.36	0.56
1:B:303:MSE:HE2	1:B:322:VAL:HG11	1.86	0.55
1:A:320:THR:HG23	1:B:340:TRP:HB2	1.88	0.55
1:B:226:VAL:HB	1:B:243:VAL:HG22	1.88	0.54
1:A:279:SER:OG	1:A:356:ASN:HB2	2.08	0.54
1:A:153:MSE:N	1:A:175:ASP:OD1	2.39	0.54
1:B:153:MSE:HE2	1:B:240:GLU:OE1	2.09	0.53
1:B:227:ASN:HA	1:B:242:ASN:OD1	2.09	0.53
1:A:179:ALA:HB3	1:A:234:GLY:HA3	1.92	0.52
1:A:198:GLY:HA3	1:A:210:TRP:CZ2	2.45	0.51
1:A:297:ARG:HH11	1:A:297:ARG:HG2	1.74	0.51
1:A:148:PRO:O	1:A:149:ALA:CB	2.59	0.51
1:B:198:GLY:HA3	1:B:210:TRP:CZ2	2.47	0.50
1:A:154:PRO:HG3	2:A:582:HOH:O	2.12	0.49
1:B:301:ILE:HG12	1:B:302:ILE:N	2.29	0.48
1:A:226:VAL:HB	1:A:243:VAL:HG23	1.95	0.48
1:B:276:MSE:HE2	2:B:690:HOH:O	2.13	0.48
1:B:151:LEU:O	1:B:237:MSE:HE1	2.15	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:263:ILE:HD13	1:A:277:PRO:HA	1.97	0.47
1:B:167:GLY:HA3	1:B:374:ILE:HD11	1.96	0.47
1:A:340:TRP:HB2	1:B:320:THR:HG23	1.96	0.46
1:B:226:VAL:HB	1:B:243:VAL:CG2	2.45	0.46
1:B:223:ASP:OD2	1:B:246:HIS:HD2	1.98	0.46
1:B:224:VAL:HB	1:B:245:THR:HG22	1.97	0.46
1:B:371:ARG:HG3	1:B:371:ARG:HH11	1.80	0.46
1:A:193:ASN:HA	1:A:194:PRO:C	2.40	0.45
1:B:152:THR:CG2	1:B:233:LEU:HD12	2.47	0.45
1:B:148:PRO:N	2:B:687:HOH:O	2.50	0.45
1:B:176:GLU:CG	1:B:237:MSE:HE2	2.44	0.44
1:A:177:ASN:O	1:A:237:MSE:HE1	2.18	0.43
1:A:178:ILE:HD13	1:A:237:MSE:CE	2.49	0.43
1:A:178:ILE:HD13	1:A:237:MSE:HE1	2.00	0.43
1:B:175:ASP:OD1	1:B:176:GLU:HG2	2.18	0.43
1:B:369:VAL:HB	2:B:588:HOH:O	2.18	0.43
1:A:330:TYR:CE1	1:A:391:LEU:HG	2.54	0.42
1:A:371:ARG:HG3	1:A:371:ARG:HH11	1.84	0.42
1:B:198:GLY:HA3	1:B:210:TRP:CH2	2.54	0.42
1:B:151:LEU:HB3	1:B:240:GLU:HB2	2.01	0.42
1:B:241:ASP:CG	1:B:242:ASN:N	2.78	0.42
1:B:357:VAL:HA	2:B:748:HOH:O	2.20	0.42
1:A:303:MSE:HE2	1:A:322:VAL:HG11	2.00	0.42
1:B:361:ASN:ND2	2:B:502:HOH:O	2.49	0.42
1:B:152:THR:HB	1:B:237:MSE:CE	2.49	0.42
1:B:240:GLU:O	1:B:241:ASP:HB2	2.20	0.42
1:B:276:MSE:HA	1:B:277:PRO:HD3	1.93	0.42
1:A:226:VAL:HB	1:A:243:VAL:HG22	2.01	0.41
1:A:237:MSE:HE2	2:A:578:HOH:O	2.20	0.41
1:B:152:THR:HG21	1:B:233:LEU:HD12	2.02	0.41
1:B:241:ASP:CG	1:B:242:ASN:H	2.29	0.41
1:B:172:ILE:HD12	1:B:172:ILE:N	2.36	0.40
1:B:297:ARG:HG2	1:B:326:THR:HG23	2.03	0.40
1:A:172:ILE:HD12	1:A:172:ILE:N	2.37	0.40
1:A:150:HIS:CD2	1:A:150:HIS:N	2.87	0.40

All (3) 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)
2:A:739:HOH:O	2:A:739:HOH:O[6_665]	1.84	0.36

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:756:HOH:O	2:A:756:HOH:O[8_566]	2.12	0.08
2:A:623:HOH:O	2:A:623:HOH:O[8_566]	2.13	0.07

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	259/269 (96%)	247 (95%)	11 (4%)	1 (0%)	30	19
1	B	260/269 (97%)	248 (95%)	11 (4%)	1 (0%)	30	19
All	All	519/538 (96%)	495 (95%)	22 (4%)	2 (0%)	30	19

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	149	ALA
1	A	149	ALA

5.3.2 Protein sidechains [i](#)

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	213/215 (99%)	210 (99%)	3 (1%)	59	52
1	B	213/215 (99%)	212 (100%)	1 (0%)	81	80
All	All	426/430 (99%)	422 (99%)	4 (1%)	70	67

All (4) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	150	HIS
1	A	244	GLN
1	A	407	LYS
1	B	276	MSE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	150	HIS
1	A	193	ASN
1	A	244	GLN
1	A	347	HIS
1	B	177	ASN
1	B	193	ASN
1	B	227	ASN
1	B	244	GLN
1	B	246	HIS
1	B	379	ASN
1	B	400	GLN

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](#)

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, 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	255/269 (94%)	0.49	25 (9%) 13 11	14, 21, 44, 66	0
1	B	256/269 (95%)	0.54	28 (10%) 10 9	14, 21, 46, 69	0
All	All	511/538 (94%)	0.52	53 (10%) 11 10	14, 21, 45, 69	0

All (53) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	148	PRO	7.1
1	A	408	ALA	7.1
1	A	148	PRO	6.8
1	B	147	SER	5.3
1	A	155	TYR	5.2
1	B	149	ALA	4.8
1	B	408	ALA	4.5
1	B	155	TYR	4.3
1	B	150	HIS	4.3
1	B	236	GLY	4.1
1	A	149	ALA	4.0
1	B	232	ASP	3.4
1	B	238	PHE	3.2
1	A	151	LEU	3.1
1	A	407	LYS	3.1
1	B	151	LEU	3.1
1	A	406	ALA	3.0
1	B	275	SER	3.0
1	B	230	GLY	2.9
1	A	197	GLU	2.8
1	B	399	ASP	2.8
1	A	399	ASP	2.8
1	A	233	LEU	2.7
1	B	402	ARG	2.7

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Mol	Chain	Res	Type	RSRZ
1	A	232	ASP	2.7
1	B	271	GLU	2.7
1	B	231	VAL	2.7
1	B	234	GLY	2.6
1	A	175	ASP	2.6
1	A	150	HIS	2.6
1	B	217	LYS	2.6
1	A	271	GLU	2.6
1	A	402	ARG	2.5
1	B	407	LYS	2.5
1	B	403	ALA	2.4
1	A	238	PHE	2.4
1	B	178	ILE	2.4
1	B	173	ARG	2.3
1	B	233	LEU	2.3
1	B	197	GLU	2.3
1	A	173	ARG	2.3
1	A	236	GLY	2.3
1	A	177	ASN	2.3
1	B	186	LYS	2.3
1	A	178	ILE	2.2
1	B	244	GLN	2.2
1	A	279	SER	2.2
1	A	179	ALA	2.2
1	A	244	GLN	2.1
1	B	270	GLY	2.1
1	A	235	GLU	2.1
1	A	229	TYR	2.0
1	B	175	ASP	2.0

6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

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

6.3 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

6.4 Ligands ⓘ

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.