



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 04:41 PM UTC

PDB ID : 7Y7U / pdb_00007y7u
Title : Dimeric structure of a Quorum-Quenching metallo-hydrolase, LrsL
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Deposited on : 2022-06-22
Resolution : 1.89 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
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

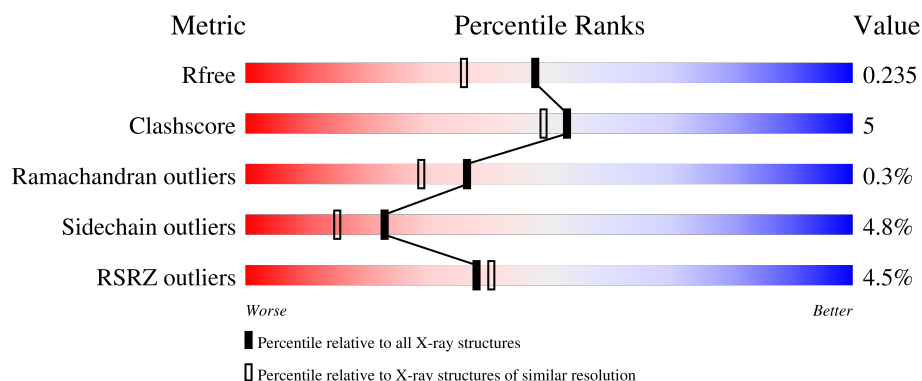
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.89 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	7789 (1.90-1.90)
Clashscore	190562	8410 (1.90-1.90)
Ramachandran outliers	187476	8333 (1.90-1.90)
Sidechain outliers	187428	8333 (1.90-1.90)
RSRZ outliers	180081	7790 (1.90-1.90)

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	295	4% 84% 12% ..
1	B	295	5% 83% 13% ..

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 4713 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 MBL fold metallo-hydrolase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	288	Total	C	N	O	S	0	1	0
			2204	1405	374	419	6			
1	B	288	Total	C	N	O	S	0	2	0
			2212	1409	375	422	6			

There are 22 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	45	MET	-	initiating methionine	UNP A0A222F232
A	46	THR	-	expression tag	UNP A0A222F232
A	47	GLY	-	expression tag	UNP A0A222F232
A	332	THR	-	expression tag	UNP A0A222F232
A	333	GLY	-	expression tag	UNP A0A222F232
A	334	LEU	-	expression tag	UNP A0A222F232
A	335	GLU	-	expression tag	UNP A0A222F232
A	336	VAL	-	expression tag	UNP A0A222F232
A	337	LEU	-	expression tag	UNP A0A222F232
A	338	PHE	-	expression tag	UNP A0A222F232
A	339	GLN	-	expression tag	UNP A0A222F232
B	45	MET	-	initiating methionine	UNP A0A222F232
B	46	THR	-	expression tag	UNP A0A222F232
B	47	GLY	-	expression tag	UNP A0A222F232
B	332	THR	-	expression tag	UNP A0A222F232
B	333	GLY	-	expression tag	UNP A0A222F232
B	334	LEU	-	expression tag	UNP A0A222F232
B	335	GLU	-	expression tag	UNP A0A222F232
B	336	VAL	-	expression tag	UNP A0A222F232
B	337	LEU	-	expression tag	UNP A0A222F232
B	338	PHE	-	expression tag	UNP A0A222F232
B	339	GLN	-	expression tag	UNP A0A222F232

- Molecule 2 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	2	Total 2	Zn 2	0	0
2	B	2	Total 2	Zn 2	0	0

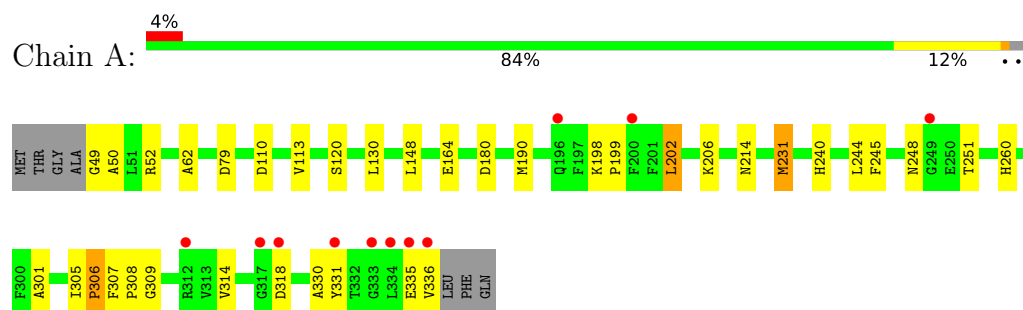
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	157	Total 157	O 157	0	0
3	B	136	Total 136	O 136	0	0

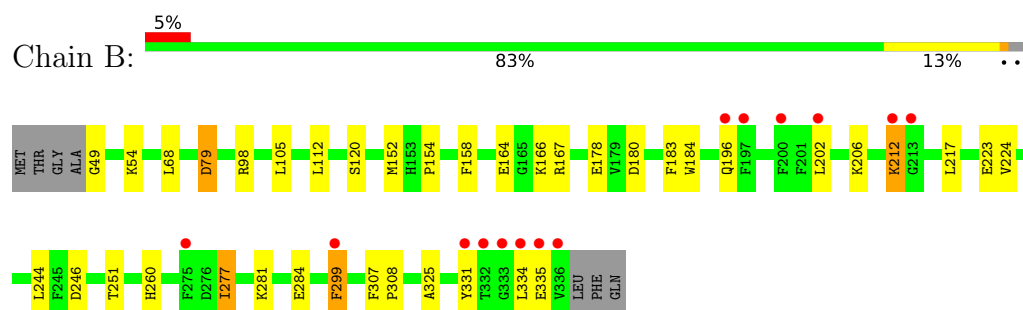
3 Residue-property plots [i](#)

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: MBL fold metallo-hydrolase



- Molecule 1: MBL fold metallo-hydrolase



4 Data and refinement statistics

Property	Value	Source
Space group	P 41 3 2	Depositor
Cell constants a, b, c, α , β , γ	166.51Å 166.51Å 166.51Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	46.18 – 1.89 46.18 – 1.89	Depositor EDS
% Data completeness (in resolution range)	99.9 (46.18-1.89) 100.0 (46.18-1.89)	Depositor EDS
R_{merge}	0.15	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.05 (at 1.88Å)	Xtriage
Refinement program	REFMAC 5.8.0349	Depositor
R, R_{free}	0.191 , 0.235 0.191 , 0.235	Depositor DCC
R_{free} test set	2000 reflections (3.14%)	wwPDB-VP
Wilson B-factor (Å ²)	44.7	Xtriage
Anisotropy	0.000	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 42.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	4713	wwPDB-VP
Average B, all atoms (Å ²)	47.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.70% 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: ZN

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	0.72	2/2258 (0.1%)	1.14	2/3070 (0.1%)
1	B	0.70	2/2266 (0.1%)	1.18	4/3081 (0.1%)
All	All	0.71	4/4524 (0.1%)	1.16	6/6151 (0.1%)

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	236	HIS	CE1-NE2	6.46	1.39	1.32
1	A	260	HIS	CE1-NE2	6.17	1.38	1.32
1	A	240	HIS	CE1-NE2	6.02	1.38	1.32
1	B	260	HIS	CE1-NE2	5.70	1.38	1.32

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	248	ASN	CB-CA-C	-8.48	100.50	111.86
1	B	299	PHE	CA-CB-CG	5.86	119.66	113.80
1	B	224	VAL	N-CA-C	-5.81	106.15	111.67
1	A	318	ASP	CB-CA-C	5.81	120.38	110.56
1	B	167	ARG	CB-CG-CD	-5.26	99.20	111.30
1	B	284	GLU	CB-CA-C	5.06	119.45	110.85

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	2204	0	2127	24	0
1	B	2212	0	2130	20	0
2	A	2	0	0	0	0
2	B	2	0	0	0	0
3	A	157	0	0	2	0
3	B	136	0	0	6	0
All	All	4713	0	4257	43	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 5.

All (43) 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:52:ARG:HG2	1:A:331:TYR:HE2	1.09	1.06
1:B:79:ASP:HB3	3:B:501:HOH:O	1.58	1.03
1:A:52:ARG:HG2	1:A:331:TYR:CE2	1.94	1.02
1:A:231:MET:HE2	1:A:244:LEU:HD21	1.67	0.77
1:B:229:ARG:HD3	3:B:613:HOH:O	1.85	0.76
1:A:52:ARG:CG	1:A:331:TYR:HE2	1.94	0.72
1:A:202:LEU:O	1:A:206:LYS:HG3	1.94	0.66
1:B:49:GLY:N	3:B:503:HOH:O	2.30	0.64
1:A:231:MET:CE	1:A:244:LEU:HD21	2.29	0.62
1:A:330:ALA:HB2	1:B:98:ARG:NH1	2.16	0.61
1:B:202:LEU:O	1:B:206:LYS:HG3	2.01	0.59
1:A:113:VAL:CG1	1:A:148:LEU:HD13	2.32	0.59
1:A:335:GLU:HA	1:A:335:GLU:OE2	2.06	0.55
1:B:79:ASP:CB	3:B:501:HOH:O	2.33	0.54
1:B:54:LYS:HD3	1:B:331:TYR:HE2	1.73	0.53
1:B:105:LEU:HD11	1:B:112:LEU:HG	1.90	0.52
1:B:54:LYS:HD3	1:B:331:TYR:CE2	2.46	0.51
1:A:110:ASP:OD1	1:A:110:ASP:N	2.41	0.51
1:B:152:MET:HB2	1:B:184:TRP:CD1	2.46	0.51
1:B:79:ASP:CG	3:B:501:HOH:O	2.52	0.51
1:A:190:MET:SD	1:A:202:LEU:HD13	2.52	0.50
1:A:307:PHE:CD1	1:A:308:PRO:HA	2.47	0.49
1:B:178:GLU:HB3	1:B:217:LEU:HB3	1.95	0.49
1:A:113:VAL:HG13	1:A:148:LEU:HD13	1.95	0.48
1:B:244:LEU:HD11	1:B:251:THR:CG2	2.44	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:49:GLY:N	3:A:507:HOH:O	2.46	0.47
1:A:305:ILE:O	1:A:306:PRO:C	2.57	0.47
1:B:154:PRO:HA	1:B:158:PHE:CD1	2.51	0.46
1:B:299:PHE:HE1	1:B:325:ALA:HB2	1.82	0.45
1:A:287:LYS:O	1:A:291:ASP:HB2	2.17	0.45
1:B:212:LYS:HD3	1:B:212:LYS:HA	1.81	0.45
1:B:234:PRO:HG2	1:B:281:LYS:HE3	1.98	0.44
1:A:335:GLU:O	1:A:336:VAL:HB	2.16	0.44
1:A:190:MET:SD	1:A:202:LEU:CD1	3.07	0.43
1:A:214:ASN:ND2	3:A:502:HOH:O	2.29	0.43
1:A:245:PHE:O	1:A:251:THR:HA	2.19	0.43
1:A:301:ALA:HA	1:A:309:GLY:O	2.19	0.42
1:A:307:PHE:CG	1:A:308:PRO:HA	2.54	0.42
1:A:198:LYS:N	1:A:199:PRO:CD	2.84	0.41
1:B:79:ASP:OD2	3:B:501:HOH:O	2.22	0.41
1:A:50:ALA:HA	1:A:62:ALA:O	2.21	0.41
1:B:307:PHE:CD1	1:B:308:PRO:HA	2.56	0.41
1:B:183:PHE:CZ	1:B:277:ILE:HD12	2.56	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	287/295 (97%)	277 (96%)	10 (4%)	0	100	100
1	B	288/295 (98%)	279 (97%)	7 (2%)	2 (1%)	18	10
All	All	575/590 (98%)	556 (97%)	17 (3%)	2 (0%)	36	29

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	334	LEU
1	B	335	GLU

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	221/225 (98%)	211 (96%)	10 (4%)	24	17
1	B	222/225 (99%)	211 (95%)	11 (5%)	22	14
All	All	443/450 (98%)	422 (95%)	21 (5%)	23	15

All (21) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	79	ASP
1	A	120	SER
1	A	130	LEU
1	A	164	GLU
1	A	180	ASP
1	A	202	LEU
1	A	231	MET
1	A	291	ASP
1	A	306	PRO
1	A	314	VAL
1	B	68	LEU
1	B	79	ASP
1	B	120	SER
1	B	164	GLU
1	B	166	LYS
1	B	180	ASP
1	B	196	GLN
1	B	212	LYS
1	B	223	GLU
1	B	246	ASP
1	B	277	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	260	HIS
1	A	292	GLN
1	B	196	GLN
1	B	292	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](#)

Of 4 ligands modelled in this entry, 4 are 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 [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	288/295 (97%)	0.17	12 (4%) 40 43	17, 41, 71, 113	1 (0%)
1	B	288/295 (97%)	0.41	14 (4%) 35 37	17, 45, 73, 121	2 (0%)
All	All	576/590 (97%)	0.29	26 (4%) 38 40	17, 43, 73, 121	3 (0%)

All (26) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	336	VAL	7.3
1	B	336	VAL	7.0
1	A	334	LEU	4.7
1	B	334	LEU	4.6
1	B	333	GLY	3.6
1	B	299	PHE	3.5
1	A	299	PHE	3.1
1	A	333	GLY	3.1
1	B	331	TYR	2.9
1	A	200	PHE	2.9
1	B	332	THR	2.8
1	B	212	LYS	2.8
1	A	331	TYR	2.7
1	A	317	GLY	2.5
1	A	318	ASP	2.5
1	B	200	PHE	2.5
1	B	213	GLY	2.4
1	B	275	PHE	2.4
1	B	335	GLU	2.4
1	B	197	PHE	2.3
1	A	196	GLN	2.3
1	A	249	GLY	2.2
1	B	196	GLN	2.2
1	A	312[A]	ARG	2.2

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Mol	Chain	Res	Type	RSRZ
1	B	202	LEU	2.1
1	A	335	GLU	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	ZN	A	401	1/1	0.99	0.02	35,35,35,35	0
2	ZN	B	402	1/1	0.99	0.03	43,43,43,43	0
2	ZN	B	401	1/1	1.00	0.02	42,42,42,42	0
2	ZN	A	402	1/1	1.00	0.02	36,36,36,36	0

6.5 Other polymers [i](#)

There are no such residues in this entry.