



Full wwPDB X-ray Structure Validation Report ⓘ

Jun 8, 2026 – 06:55 PM JST

PDB ID : 26LM / pdb_000026lm
Title : Crystal structure of cblb
Authors : Zhiyan, D.; Danyan, C.; Lingyu, S.; Bing, X.
Deposited on : 2026-05-06
Resolution : 1.95 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

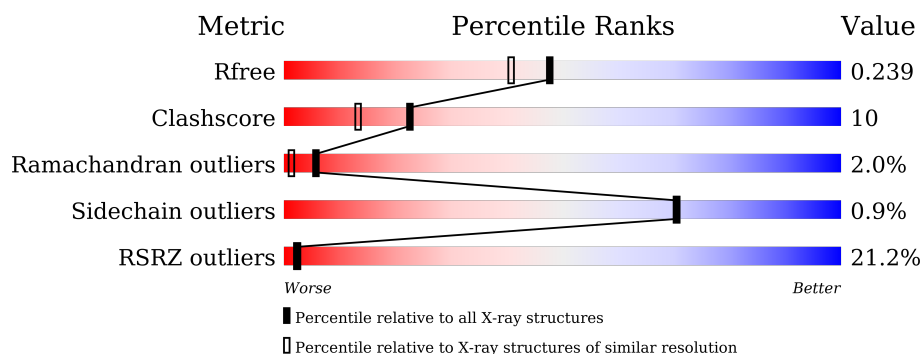
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.95 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	3494 (1.96-1.96)
Clashscore	190562	3612 (1.96-1.96)
Ramachandran outliers	187476	3587 (1.96-1.96)
Sidechain outliers	187428	3587 (1.96-1.96)
RSRZ outliers	180081	3495 (1.96-1.96)

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	399	<div> <div>21%</div> <div>81%</div> <div>16%</div> <div>..</div> </div>

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 3345 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 E3 ubiquitin-protein ligase CBL-B.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	396	Total	C	N	O	S	0	2	0
			3165	2040	526	577	22			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	35	MET	-	initiating methionine	UNP Q13191
A	363	PHE	TYR	conflict	UNP Q13191
A	428	GLU	-	expression tag	UNP Q13191
A	429	ASN	-	expression tag	UNP Q13191
A	430	LEU	-	expression tag	UNP Q13191
A	431	TYR	-	expression tag	UNP Q13191
A	432	PHE	-	expression tag	UNP Q13191
A	433	GLN	-	expression tag	UNP Q13191

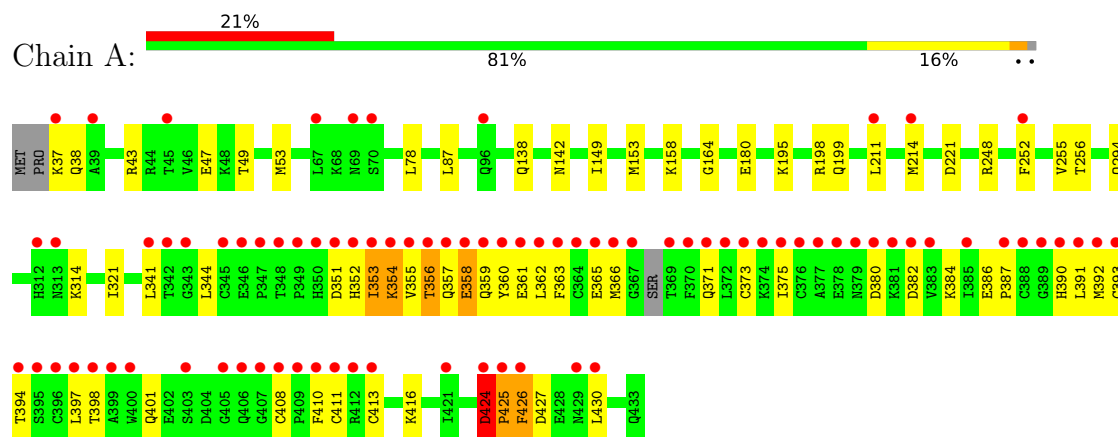
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	180	Total	O	0	0
			180	180		

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: E3 ubiquitin-protein ligase CBL-B



4 Data and refinement statistics

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, α , β , γ	71.84Å 71.84Å 212.31Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.42 – 1.95 50.42 – 1.95	Depositor EDS
% Data completeness (in resolution range)	100.0 (50.42-1.95) 100.0 (50.42-1.95)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.71 (at 1.95Å)	Xtriage
Refinement program	PHENIX (1.19.2_4158: ???)	Depositor
R, R_{free}	0.212 , 0.240 0.213 , 0.239	Depositor DCC
R_{free} test set	2000 reflections (4.81%)	wwPDB-VP
Wilson B-factor (Å ²)	36.1	Xtriage
Anisotropy	0.368	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 43.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	3345	wwPDB-VP
Average B, all atoms (Å ²)	46.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.91% 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 [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	0.37	0/3244	0.63	1/4396 (0.0%)

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

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	A	424	ASP	C-N-CD	-6.02	100.33	125.00

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	354	LYS	Peptide
1	A	424	ASP	Peptide

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	3165	0	3041	65	0
2	A	180	0	0	4	1
All	All	3345	0	3041	65	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 10.

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:A:359:GLN:HA	1:A:362:LEU:HB2	1.64	0.80
1:A:384:LYS:HE2	1:A:386:GLU:OE1	1.86	0.76
1:A:256:THR:HG23	1:A:355:VAL:HG22	1.70	0.73
1:A:158:LYS:NZ	2:A:502:HOH:O	2.20	0.73
1:A:256:THR:HG22	1:A:359:GLN:HE21	1.55	0.70
1:A:49:THR:O	1:A:53:MET:HG3	1.94	0.67
1:A:180:GLU:OE2	2:A:501:HOH:O	2.14	0.64
1:A:411:CYS:SG	1:A:413:CYS:HB2	2.37	0.64
1:A:195:LYS:O	1:A:199:GLN:HG3	1.99	0.63
1:A:357:GLN:HB2	1:A:359:GLN:CD	2.24	0.63
1:A:53:MET:HE3	1:A:78:LEU:HD22	1.81	0.62
1:A:355:VAL:HB	1:A:360:TYR:CZ	2.35	0.61
1:A:387:PRO:HD3	1:A:416:LYS:HB2	1.82	0.60
1:A:256:THR:CG2	1:A:355:VAL:HG22	2.31	0.60
1:A:371:GLN:HG2	1:A:380:ASP:OD1	2.02	0.60
1:A:294:GLN:NE2	2:A:505:HOH:O	2.26	0.58
1:A:198:ARG:HG2	1:A:198:ARG:HH11	1.68	0.57
1:A:355:VAL:HA	1:A:357:GLN:HE22	1.70	0.56
1:A:252:PHE:CD1	1:A:344:LEU:HD11	2.40	0.56
1:A:373:CYS:HB2	1:A:393:CYS:HB3	1.88	0.55
1:A:252:PHE:HZ	1:A:341:LEU:HG	1.73	0.54
1:A:255:VAL:CG1	1:A:355:VAL:HG21	2.39	0.53
1:A:366:MET:SD	1:A:366:MET:N	2.82	0.53
1:A:359:GLN:O	1:A:363:PHE:HB2	2.10	0.52
1:A:355:VAL:O	1:A:356:THR:HG23	2.10	0.51
1:A:357:GLN:OE1	1:A:359:GLN:NE2	2.44	0.51
1:A:375:ILE:HG12	1:A:410:PHE:CE2	2.46	0.51
1:A:149:ILE:O	1:A:153:MET:HG3	2.12	0.49
1:A:365:GLU:N	1:A:365:GLU:OE1	2.45	0.49
1:A:390:HIS:CE1	1:A:411:CYS:SG	3.05	0.49
1:A:425:PRO:C	1:A:427:ASP:H	2.20	0.49
1:A:392:MET:HE2	1:A:397:LEU:HB2	1.95	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:252:PHE:CE1	1:A:341:LEU:HD23	2.48	0.48
1:A:351:ASP:O	1:A:353:ILE:N	2.46	0.48
1:A:357:GLN:C	1:A:359:GLN:H	2.21	0.48
1:A:358:GLU:OE1	1:A:362:LEU:HD13	2.14	0.47
1:A:252:PHE:CD1	1:A:252:PHE:C	2.93	0.47
1:A:321:ILE:CG1	1:A:341:LEU:HB2	2.45	0.47
1:A:390:HIS:HE1	1:A:411:CYS:SG	2.38	0.47
1:A:321:ILE:HG12	1:A:341:LEU:HB2	1.99	0.45
1:A:221:ASP:OD1	2:A:503:HOH:O	2.21	0.45
1:A:391:LEU:O	1:A:392:MET:HB3	2.17	0.44
1:A:248:ARG:HD3	1:A:351:ASP:OD1	2.17	0.44
1:A:357:GLN:O	1:A:358:GLU:HG3	2.18	0.44
1:A:87:LEU:HD13	1:A:164:GLY:HA3	1.98	0.44
1:A:255:VAL:HG11	1:A:355:VAL:HG11	1.99	0.43
1:A:43:ARG:O	1:A:47:GLU:HG3	2.19	0.43
1:A:392:MET:CE	1:A:397:LEU:HB2	2.48	0.43
1:A:252:PHE:CZ	1:A:341:LEU:HD23	2.53	0.42
1:A:138:GLN:HE21	1:A:142:ASN:CG	2.27	0.42
1:A:353:ILE:HG22	1:A:356:THR:OG1	2.20	0.42
1:A:387:PRO:HD3	1:A:416:LYS:HG3	2.01	0.42
1:A:398:THR:HA	1:A:401:GLN:HG2	2.02	0.42
1:A:357:GLN:C	1:A:359:GLN:N	2.78	0.41
1:A:214:MET:HE3	1:A:214:MET:HB3	1.94	0.41
1:A:195:LYS:HG3	1:A:198:ARG:NH2	2.36	0.41
1:A:424:ASP:C	1:A:425:PRO:O	2.63	0.41
1:A:353:ILE:O	1:A:357:GLN:NE2	2.54	0.41
1:A:382:ASP:OD1	1:A:394:THR:HG23	2.20	0.41
1:A:255:VAL:HG11	1:A:355:VAL:HG21	2.01	0.41
1:A:314:LYS:NZ	1:A:314:LYS:HB3	2.36	0.41
1:A:408:CYS:SG	1:A:410:PHE:HB2	2.61	0.41
1:A:252:PHE:HZ	1:A:341:LEU:CG	2.34	0.41
1:A:211:LEU:HD21	1:A:361:GLU:HG3	2.01	0.40
1:A:387:PRO:HD3	1:A:416:LYS:CB	2.49	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)
2:A:597:HOH:O	2:A:656:HOH:O[6_454]	2.04	0.16

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	394/399 (99%)	366 (93%)	20 (5%)	8 (2%)	6 1

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	425	PRO
1	A	430	LEU
1	A	352	HIS
1	A	356	THR
1	A	38	GLN
1	A	354	LYS
1	A	426	PHE
1	A	358	GLU

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	331/359 (92%)	328 (99%)	3 (1%)	70 70

All (3) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	37	LYS
1	A	353	ILE
1	A	426	PHE

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	138	GLN
1	A	357	GLN
1	A	359	GLN
1	A	379	ASN

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 [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	396/399 (99%)	1.15	84 (21%) 2 2	20, 39, 93, 133	2 (0%)

All (84) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	355	VAL	9.9
1	A	410	PHE	7.6
1	A	396	CYS	7.3
1	A	370	PHE	7.2
1	A	341	LEU	6.9
1	A	377	ALA	6.7
1	A	354	LYS	6.5
1	A	356	THR	6.3
1	A	372	LEU	6.0
1	A	353	ILE	5.7
1	A	369	THR	5.6
1	A	378	GLU	5.4
1	A	397	LEU	5.4
1	A	373	CYS	5.3
1	A	347	PRO	5.2
1	A	313	ASN	5.1
1	A	375	ILE	5.0
1	A	364	CYS	4.9
1	A	69	ASN	4.9
1	A	390	HIS	4.8
1	A	411	CYS	4.8
1	A	425	PRO	4.7
1	A	357	GLN	4.7
1	A	348	THR	4.7
1	A	367	GLY	4.6
1	A	365	GLU	4.5
1	A	374	LYS	4.5

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Mol	Chain	Res	Type	RSRZ
1	A	400	TRP	4.4
1	A	376	CYS	4.4
1	A	360	TYR	4.4
1	A	395[A]	SER	4.4
1	A	252	PHE	4.3
1	A	391	LEU	4.3
1	A	393	CYS	4.3
1	A	349	PRO	4.3
1	A	399	ALA	4.2
1	A	406	GLN	4.0
1	A	359	GLN	4.0
1	A	345	CYS	4.0
1	A	342	THR	3.9
1	A	398	THR	3.8
1	A	312	HIS	3.7
1	A	363	PHE	3.7
1	A	362	LEU	3.6
1	A	366	MET	3.6
1	A	413	CYS	3.6
1	A	403	SER	3.5
1	A	394	THR	3.5
1	A	379	ASN	3.4
1	A	408	CYS	3.4
1	A	358	GLU	3.3
1	A	352	HIS	3.3
1	A	214	MET	3.3
1	A	351	ASP	3.3
1	A	343	GLY	3.2
1	A	430	LEU	3.2
1	A	350	HIS	3.1
1	A	45[A]	THR	2.9
1	A	429	ASN	2.9
1	A	421	ILE	2.9
1	A	389	GLY	2.9
1	A	387	PRO	2.8
1	A	371	GLN	2.8
1	A	412	ARG	2.8
1	A	39	ALA	2.8
1	A	67	LEU	2.8
1	A	388	CYS	2.8
1	A	392	MET	2.7
1	A	37	LYS	2.7

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Mol	Chain	Res	Type	RSRZ
1	A	96	GLN	2.7
1	A	211	LEU	2.7
1	A	383	VAL	2.7
1	A	346	GLU	2.5
1	A	380	ASP	2.5
1	A	424	ASP	2.5
1	A	409	PRO	2.4
1	A	405	GLY	2.4
1	A	407	GLY	2.4
1	A	70	SER	2.4
1	A	361	GLU	2.3
1	A	426	PHE	2.2
1	A	382	ASP	2.1
1	A	381	LYS	2.1
1	A	385	ILE	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](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.