



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 12, 2026 – 07:53 PM UTC

PDB ID : 5ILB / pdb_00005ilb
Title : Crystal structure of protease domain of Deg2 linked with the PDZ domain of Deg9
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Deposited on : 2016-03-04
Resolution : 1.85 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

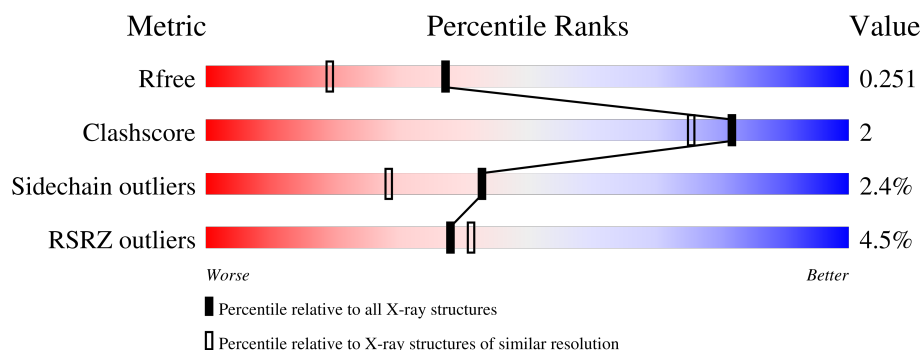
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.85 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	3428 (1.86-1.86)
Clashscore	190562	3579 (1.86-1.86)
Sidechain outliers	187428	3553 (1.86-1.86)
RSRZ outliers	180081	3429 (1.86-1.86)

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	474	4% 93% 5% •
1	B	474	5% 89% 8% ••

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 8263 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 *Protease Do-like 2, chloroplastic, Protease Do-like 9*.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	466	Total	C	N	O	S	0	9	0
			3621	2312	606	689	14			
1	B	463	Total	C	N	O	S	0	3	0
			3597	2290	612	681	14			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	109	GLY	-	expression tag	UNP O82261
B	109	GLY	-	expression tag	UNP O82261

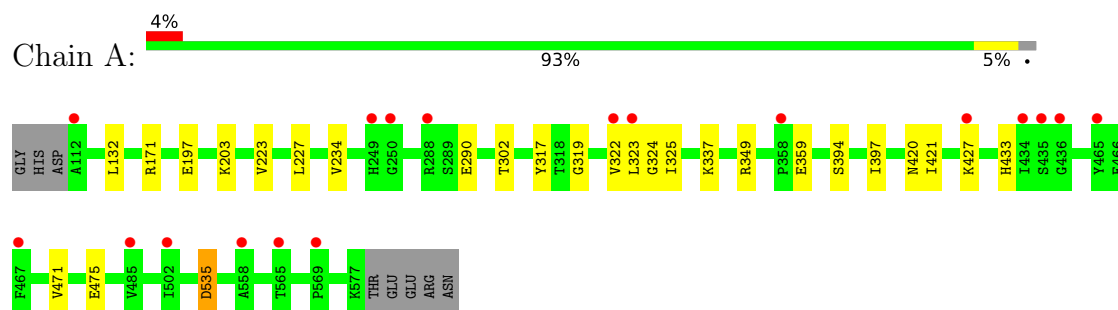
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	519	Total	O	0	0
			519	519		
2	B	526	Total	O	0	0
			526	526		

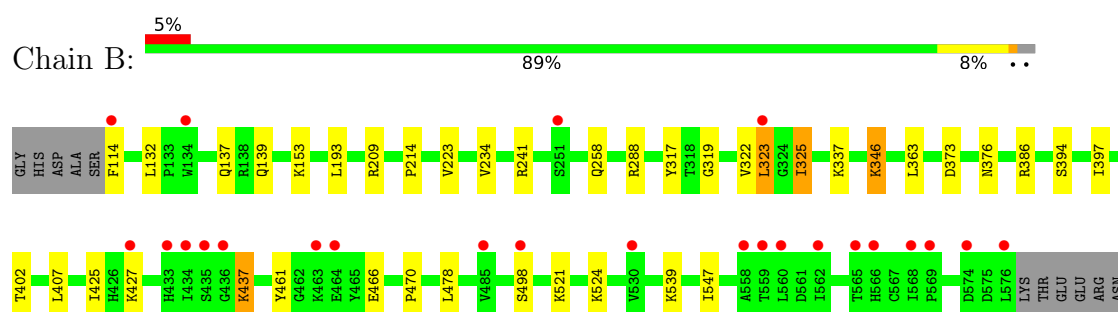
3 Residue-property plots

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: Protease Do-like 2, chloroplastic, Protease Do-like 9



- Molecule 1: Protease Do-like 2, chloroplastic, Protease Do-like 9



4 Data and refinement statistics

Property	Value	Source
Space group	H 3	Depositor
Cell constants a, b, c, α , β , γ	106.09Å 106.09Å 255.96Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	26.52 – 1.85 26.52 – 1.85	Depositor EDS
% Data completeness (in resolution range)	96.0 (26.52-1.85) 95.9 (26.52-1.85)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.46 (at 1.85Å)	Xtriage
Refinement program	PHENIX 1.8.1_1168	Depositor
R, R_{free}	0.190 , 0.225 (Not available) , 0.251	Depositor DCC
R_{free} test set	4396 reflections (4.81%)	wwPDB-VP
Wilson B-factor (Å ²)	18.3	Xtriage
Anisotropy	0.078	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 50.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.35$	Xtriage
Estimated twinning fraction	0.000 for -h,1/3*h-1/3*k-1/3*l,-4/3*h-8/3*k +1/3*l 0.002 for -1/3*h+1/3*k+1/3*l,-k,8/3*h+4/ 3*k+1/3*l 0.004 for -2/3*h-1/3*k-1/3*l,-1/3*h-2/3*k+ 1/3*l,-4/3*h+4/3*k+1/3*l 0.000 for 1/3*h+2/3*k-1/3*l,-k,-8/3*h-4/3* k-1/3*l 0.003 for -1/3*h-2/3*k+1/3*l,-2/3*h-1/3*k- 1/3*l,4/3*h-4/3*k-1/3*l 0.000 for -h,2/3*h+1/3*k+1/3*l,4/3*h+8/3 *k-1/3*l 0.020 for h,-h-k,-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	8263	wwPDB-VP
Average B, all atoms (Å ²)	22.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.03% 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	0.33	0/3720	0.71	2/5050 (0.0%)
1	B	0.33	0/3679	0.72	2/4993 (0.0%)
All	All	0.33	0/7399	0.72	4/10043 (0.0%)

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	132	LEU	CA-C-N	6.07	125.77	119.82
1	B	132	LEU	C-N-CA	6.07	125.77	119.82
1	A	132	LEU	CA-C-N	5.78	125.48	119.82
1	A	132	LEU	C-N-CA	5.78	125.48	119.82

There are no chirality outliers.

There are no planarity outliers.

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	3621	0	3623	14	0
1	B	3597	0	3596	23	0
2	A	519	0	0	5	1
2	B	526	0	0	9	1
All	All	8263	0	7219	36	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 2.

All (36) 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:214:PRO:O	1:B:386:ARG:NH2	2.13	0.81
1:B:346:LYS:HE3	1:B:376:ASN:HB3	1.74	0.68
1:A:203:LYS:NZ	2:A:615:HOH:O	2.34	0.61
1:A:223:VAL:HG22	1:A:234[B]:VAL:HG12	1.82	0.61
1:B:139:GLN:NE2	2:B:601:HOH:O	2.21	0.60
1:B:223:VAL:HG22	1:B:234[B]:VAL:HG12	1.85	0.59
1:A:317:TYR:CZ	1:A:319:GLY:HA2	2.41	0.56
1:B:524:LYS:NZ	2:B:616:HOH:O	2.32	0.51
1:A:535:ASP:OD2	1:A:535:ASP:N	2.42	0.51
1:B:288:ARG:NH2	2:B:632:HOH:O	2.43	0.50
1:B:317:TYR:CZ	1:B:319:GLY:HA2	2.47	0.50
1:A:324:GLY:N	1:A:359:GLU:OE2	2.45	0.49
1:A:359:GLU:HB3	1:A:421:ILE:HG21	1.93	0.49
1:B:539:LYS:NZ	2:B:644:HOH:O	2.46	0.48
1:A:349:ARG:NH2	2:A:632:HOH:O	2.46	0.48
1:A:197:GLU:OE2	2:A:602:HOH:O	2.20	0.47
1:B:137:GLN:NE2	2:B:646:HOH:O	2.48	0.47
1:A:394:SER:HA	1:A:397:ILE:HD12	1.97	0.47
1:B:437:LYS:O	2:B:602:HOH:O	2.21	0.46
1:A:337:LYS:HE3	1:A:337:LYS:HB2	1.71	0.45
1:B:323:LEU:HD13	1:B:325:ILE:HG12	1.98	0.45
1:B:394:SER:HA	1:B:397:ILE:HD12	1.99	0.44
1:B:153:LYS:HD3	1:B:193:LEU:HD13	2.00	0.43
1:A:471[B]:VAL:HG21	1:B:478:LEU:HD12	2.00	0.43
1:B:241:ARG:HD2	1:B:258:GLN:OE1	2.18	0.42
1:B:402:THR:HG21	1:B:425:ILE:HG23	2.01	0.42
1:B:461:TYR:CZ	1:B:470:PRO:HD3	2.55	0.42
1:A:171:ARG:HD3	2:A:625:HOH:O	2.19	0.42
1:B:539:LYS:HE3	1:B:547:ILE:HD13	2.02	0.42
1:B:209:ARG:NH1	2:B:608:HOH:O	2.26	0.42
1:A:302:THR:HG23	2:A:601:HOH:O	2.19	0.41
1:B:373:ASP:OD1	2:B:603:HOH:O	2.22	0.41
1:B:337:LYS:HB2	1:B:337:LYS:HE2	1.82	0.41
1:A:427:LYS:HD2	1:A:427:LYS:HA	1.84	0.40
1:B:427:LYS:HD2	1:B:427:LYS:HA	1.89	0.40
1:B:498:SER:OG	2:B:604:HOH:O	2.22	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:983:HOH:O	2:B:1045:HOH:O[9_544]	2.10	0.10

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

There are no protein backbone outliers to report in this entry.

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	400/410 (98%)	391 (98%)	9 (2%)	44	30
1	B	397/410 (97%)	387 (98%)	10 (2%)	42	27
All	All	797/820 (97%)	778 (98%)	19 (2%)	43	28

All (19) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	227	LEU
1	A	290	GLU
1	A	322	VAL
1	A	323	LEU
1	A	325	ILE
1	A	420	ASN
1	A	433	HIS
1	A	475	GLU
1	A	535	ASP
1	B	114	PHE
1	B	322	VAL
1	B	323	LEU
1	B	325	ILE
1	B	346	LYS
1	B	363	LEU
1	B	407	LEU

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Mol	Chain	Res	Type
1	B	437	LYS
1	B	466	GLU
1	B	521	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	163	HIS
1	A	426	HIS
1	B	157	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

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

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	466/474 (98%)	0.11	18 (3%) 43 47	2, 19, 42, 56	9 (1%)
1	B	463/474 (97%)	0.11	24 (5%) 33 35	2, 20, 42, 58	3 (0%)
All	All	929/948 (97%)	0.11	42 (4%) 38 41	2, 19, 42, 58	12 (1%)

All (42) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	114	PHE	5.2
1	A	436	GLY	4.6
1	A	434	ILE	4.0
1	B	435	SER	3.6
1	A	435	SER	3.5
1	A	465	TYR	3.5
1	B	434	ILE	3.4
1	B	565	THR	3.3
1	B	560	LEU	3.2
1	B	562	ILE	3.1
1	A	250	GLY	3.0
1	B	436	GLY	2.9
1	A	502	ILE	2.9
1	B	576	LEU	2.9
1	A	467	PHE	2.8
1	B	574	ASP	2.7
1	B	427	LYS	2.7
1	A	322	VAL	2.7
1	A	485	VAL	2.7
1	B	485	VAL	2.6
1	A	323	LEU	2.6
1	A	249	HIS	2.5
1	A	558	ALA	2.5
1	A	569	PRO	2.5

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Mol	Chain	Res	Type	RSRZ
1	B	134	TRP	2.5
1	B	569	PRO	2.4
1	B	251	SER	2.4
1	B	464	GLU	2.4
1	A	288	ARG	2.3
1	B	323	LEU	2.3
1	B	566	HIS	2.3
1	A	565	THR	2.3
1	B	433	HIS	2.3
1	A	112	ALA	2.2
1	A	358	PRO	2.2
1	B	568	ILE	2.2
1	A	427	LYS	2.2
1	B	530	VAL	2.2
1	B	559	THR	2.1
1	B	463	LYS	2.0
1	B	558	ALA	2.0
1	B	498	SER	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.