



Full wwPDB X-ray Structure Validation Report ⓘ

Nov 19, 2023 – 04:28 PM JST

PDB ID : 6LOL
Title : The crystal structure of full length IpaH9.8
Authors : Ye, Y.; Huang, H.
Deposited on : 2020-01-06
Resolution : 2.75 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 : 4.02b-467
Xtriage (Phenix) : 1.13
EDS : 2.36
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.36

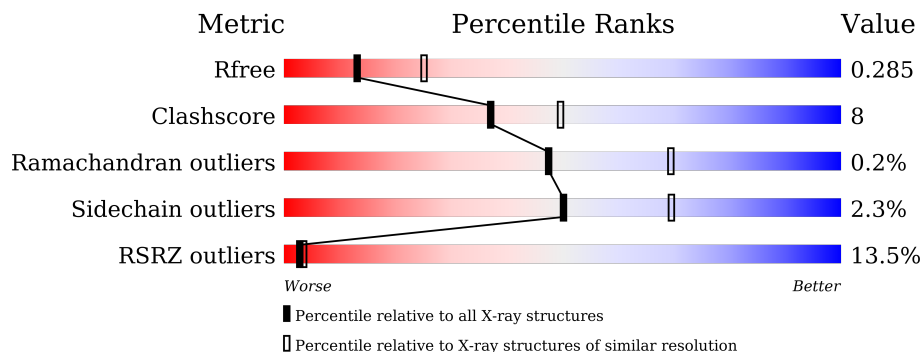
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.75 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}	130704	1235 (2.78-2.74)
Clashscore	141614	1277 (2.78-2.74)
Ramachandran outliers	138981	1257 (2.78-2.74)
Sidechain outliers	138945	1257 (2.78-2.74)
RSRZ outliers	127900	1207 (2.78-2.74)

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	526	

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 3460 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 ipaH9.8.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	466	3460	2185	598	670	7	0	0	0

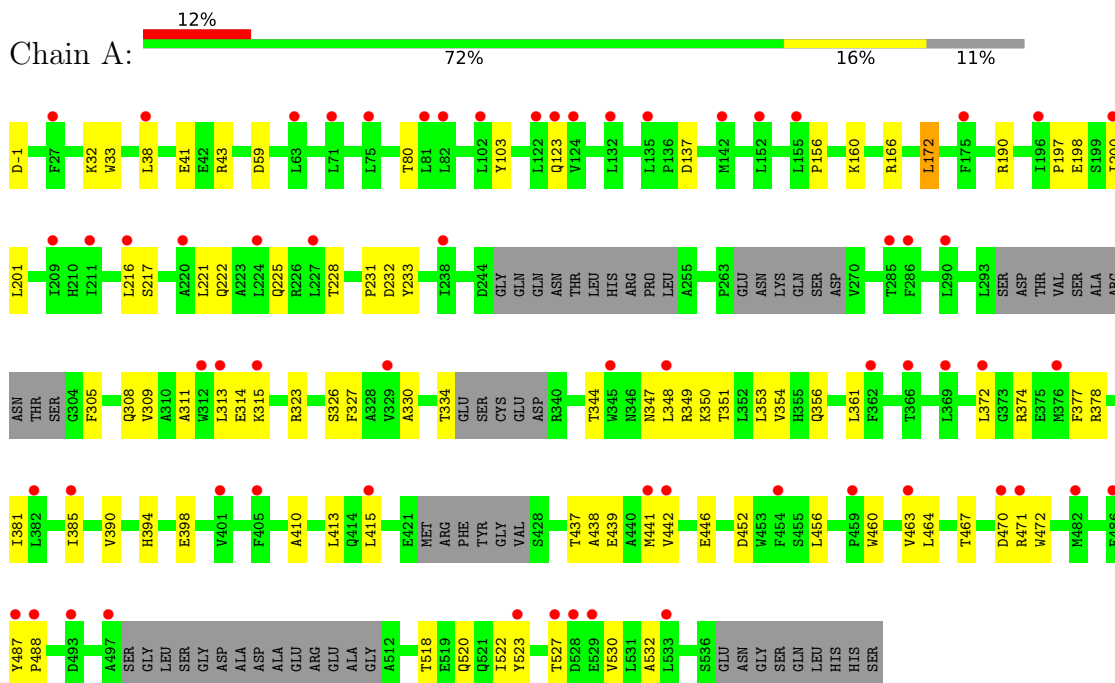
There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	ASP	-	expression tag	UNP D2AJU0
A	0	PRO	-	expression tag	UNP D2AJU0

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: E3 ubiquitin-protein ligase ipaH9.8



4 Data and refinement statistics

Property	Value	Source
Space group	I 2 2 2	Depositor
Cell constants a, b, c, α , β , γ	106.64Å 122.25Å 149.97Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	86.91 – 2.75 86.91 – 2.75	Depositor EDS
% Data completeness (in resolution range)	99.5 (86.91-2.75) 99.6 (86.91-2.75)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.33 (at 2.73Å)	Xtrriage
Refinement program	PHENIX 1.17.1_3660	Depositor
R, R_{free}	0.246 , 0.285 0.246 , 0.285	Depositor DCC
R_{free} test set	1319 reflections (5.13%)	wwPDB-VP
Wilson B-factor (Å ²)	80.1	Xtrriage
Anisotropy	0.591	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 80.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	3460	wwPDB-VP
Average B, all atoms (Å ²)	102.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.95% 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	0.26	0/3523	0.49	3/4818 (0.1%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	313	LEU	CA-CB-CG	6.16	129.47	115.30
1	A	172	LEU	CA-CB-CG	5.95	128.99	115.30
1	A	313	LEU	CB-CG-CD2	5.68	120.66	111.00

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	3460	0	3216	52	0
All	All	3460	0	3216	52	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 8.

All (52) 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:518:THR:O	1:A:522:ILE:HD12	1.86	0.75
1:A:456:LEU:HD11	1:A:520:GLN:HA	1.68	0.75
1:A:330:ALA:HB2	1:A:348:LEU:HD11	1.69	0.74
1:A:33:TRP:NE1	1:A:43:ARG:NH1	2.35	0.73
1:A:200:ILE:HD12	1:A:201:LEU:HG	1.77	0.66
1:A:200:ILE:HD12	1:A:201:LEU:N	2.11	0.65
1:A:200:ILE:CD1	1:A:201:LEU:HG	2.29	0.63
1:A:410:ALA:HA	1:A:415:LEU:HD12	1.81	0.62
1:A:530:VAL:HG23	1:A:532:ALA:H	1.66	0.60
1:A:216:LEU:HB2	1:A:221:LEU:HD11	1.84	0.59
1:A:308:GLN:HA	1:A:311:ALA:H	1.67	0.59
1:A:472:TRP:HZ3	1:A:523:TYR:HE1	1.50	0.58
1:A:413:LEU:HD21	1:A:439:GLU:HG3	1.84	0.58
1:A:217:SER:O	1:A:221:LEU:HD13	2.04	0.57
1:A:472:TRP:HZ3	1:A:523:TYR:CE1	2.22	0.57
1:A:231:PRO:O	1:A:232:ASP:HB3	2.04	0.56
1:A:353:LEU:HD21	1:A:372:LEU:HD21	1.89	0.55
1:A:166:ARG:CZ	1:A:190:ARG:HE	2.22	0.53
1:A:38:LEU:O	1:A:41:GLU:HG2	2.09	0.53
1:A:527:THR:HA	1:A:530:VAL:HG12	1.91	0.52
1:A:323:ARG:HA	1:A:326:SER:H	1.75	0.52
1:A:390:VAL:HG11	1:A:398:GLU:HA	1.91	0.52
1:A:530:VAL:HG23	1:A:532:ALA:N	2.25	0.51
1:A:347:ASN:O	1:A:351:THR:HG23	2.12	0.49
1:A:103:TYR:CD1	1:A:123:GLN:HG3	2.48	0.49
1:A:166:ARG:NH1	1:A:190:ARG:HE	2.11	0.49
1:A:305:PHE:O	1:A:309:VAL:HG13	2.13	0.49
1:A:311:ALA:HA	1:A:314:GLU:HB2	1.94	0.48
1:A:-1:ASP:OD1	1:A:-1:ASP:N	2.37	0.48
1:A:385:ILE:HD11	1:A:441:MET:HE2	1.96	0.48
1:A:323:ARG:HB2	1:A:327:PHE:CZ	2.48	0.48
1:A:197:PRO:O	1:A:200:ILE:HG13	2.16	0.46
1:A:200:ILE:HD12	1:A:201:LEU:H	1.81	0.45
1:A:166:ARG:HA	1:A:190:ARG:O	2.16	0.45
1:A:350:LYS:O	1:A:354:VAL:HG13	2.16	0.45
1:A:377:PHE:O	1:A:381:ILE:HG12	2.18	0.44
1:A:198:GLU:OE2	1:A:394:HIS:ND1	2.46	0.43
1:A:487:TYR:N	1:A:488:PRO:HD2	2.34	0.43
1:A:374:ARG:O	1:A:378:ARG:HG3	2.19	0.43
1:A:463:VAL:O	1:A:467:THR:N	2.39	0.42
1:A:356:GLN:HB3	1:A:361:LEU:HD23	2.02	0.42
1:A:460:TRP:O	1:A:464:LEU:HG	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:442:VAL:O	1:A:446:GLU:HB2	2.19	0.42
1:A:59:ASP:HB3	1:A:80:THR:OG1	2.21	0.41
1:A:172:LEU:HB3	1:A:197:PRO:HD3	2.02	0.41
1:A:228:THR:HG22	1:A:233:TYR:CZ	2.56	0.41
1:A:438:ALA:O	1:A:442:VAL:HG12	2.21	0.41
1:A:437:THR:O	1:A:441:MET:HG3	2.20	0.41
1:A:137:ASP:HA	1:A:156:PRO:HB3	2.02	0.41
1:A:32:LYS:O	1:A:32:LYS:HD2	2.22	0.40
1:A:103:TYR:HD1	1:A:123:GLN:HG3	1.84	0.40
1:A:344:THR:O	1:A:348:LEU:HD22	2.22	0.40

There are no symmetry-related clashes.

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	452/526 (86%)	414 (92%)	37 (8%)	1 (0%)	47 69

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	470	ASP

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	344/462 (74%)	336 (98%)	8 (2%)	50 69

All (8) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	160	LYS
1	A	222	GLN
1	A	225	GLN
1	A	315	LYS
1	A	334	THR
1	A	349	ARG
1	A	452	ASP
1	A	471	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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 monosaccharides 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

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	466/526 (88%)	0.78	63 (13%) 3 3	46, 91, 180, 203	0

All (63) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	312	TRP	6.4
1	A	488	PRO	6.0
1	A	175	PHE	5.6
1	A	362	PHE	5.3
1	A	471	ARG	4.9
1	A	211	ILE	4.2
1	A	482	MET	3.9
1	A	369	LEU	3.9
1	A	224	LEU	3.8
1	A	523	TYR	3.8
1	A	290	LEU	3.7
1	A	329	VAL	3.7
1	A	366	THR	3.7
1	A	132	LEU	3.7
1	A	102	LEU	3.5
1	A	441	MET	3.5
1	A	82	LEU	3.4
1	A	286	PHE	3.3
1	A	470	ASP	3.2
1	A	285	THR	3.2
1	A	71	LEU	3.2
1	A	122	LEU	3.0
1	A	442	VAL	2.9
1	A	405	PHE	2.9
1	A	463	VAL	2.8
1	A	415	LEU	2.8
1	A	216	LEU	2.8

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Mol	Chain	Res	Type	RSRZ
1	A	155	LEU	2.8
1	A	209	ILE	2.7
1	A	376	MET	2.6
1	A	345	TRP	2.6
1	A	313	LEU	2.6
1	A	529	GLU	2.6
1	A	124	VAL	2.6
1	A	142	MET	2.5
1	A	196	ILE	2.5
1	A	227	LEU	2.5
1	A	454	PHE	2.4
1	A	401	VAL	2.3
1	A	459	PRO	2.3
1	A	487	TYR	2.3
1	A	220	ALA	2.3
1	A	528	ASP	2.3
1	A	348	LEU	2.3
1	A	200	ILE	2.3
1	A	38	LEU	2.3
1	A	75	LEU	2.3
1	A	27	PHE	2.2
1	A	493	ASP	2.2
1	A	486	GLU	2.2
1	A	123	GLN	2.2
1	A	497	ALA	2.2
1	A	385	ILE	2.1
1	A	238	ILE	2.1
1	A	372	LEU	2.1
1	A	527	THR	2.1
1	A	81	LEU	2.1
1	A	63	LEU	2.0
1	A	315	LYS	2.0
1	A	135	LEU	2.0
1	A	152	LEU	2.0
1	A	382	LEU	2.0
1	A	533	LEU	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 [i](#)

There are no monosaccharides 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.