



Full wwPDB X-ray Structure Validation Report ⓘ

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PDB ID : 6VZF / pdb_00006vzf
Title : Crystal Structure of Atg11 Coiled-Coil 3
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Deposited on : 2020-02-28
Resolution : 2.03 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

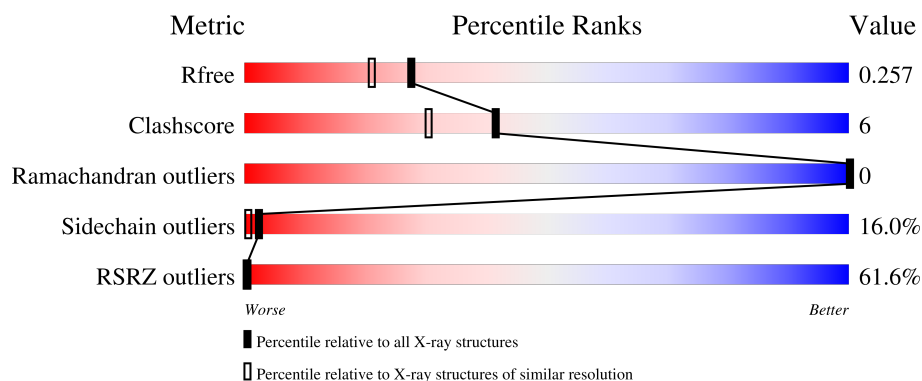
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.03 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	13299 (2.04-2.00)
Clashscore	190562	1022 (2.02-2.02)
Ramachandran outliers	187476	1014 (2.02-2.02)
Sidechain outliers	187428	1014 (2.02-2.02)
RSRZ outliers	180081	13314 (2.04-2.00)

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	105	50% 58%20%•18%

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 734 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 Autophagy-related protein 11.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	86	706	431	127	146	2	0	0	0

There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	696	GLY	-	expression tag	UNP Q12527
A	697	ALA	-	expression tag	UNP Q12527
A	698	MET	-	expression tag	UNP Q12527

- Molecule 2 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	O	S		
2	A	1	5	4	1	0	0

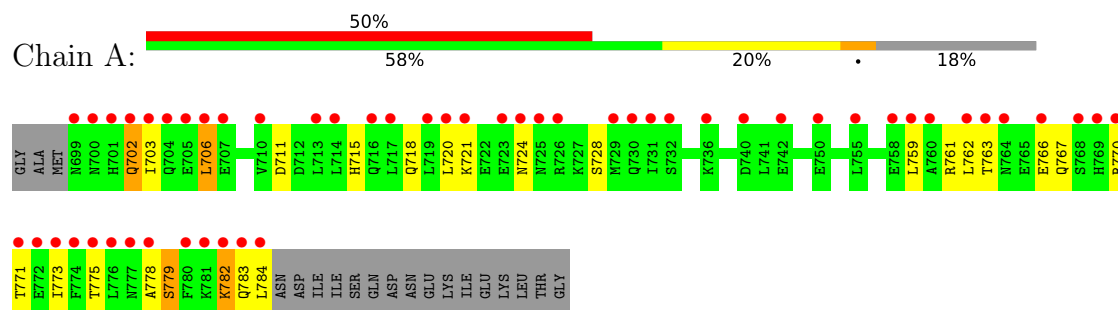
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	23	Total 23	O 23	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: Autophagy-related protein 11



4 Data and refinement statistics

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, α , β , γ	103.11Å 103.11Å 37.52Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	34.61 – 2.03 34.61 – 2.03	Depositor EDS
% Data completeness (in resolution range)	99.9 (34.61-2.03) 99.9 (34.61-2.03)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.90 (at 2.03Å)	Xtriage
Refinement program	REFMAC 5.8.0258	Depositor
R, R_{free}	0.226 , 0.246 0.233 , 0.257	Depositor DCC
R_{free} test set	781 reflections (5.19%)	wwPDB-VP
Wilson B-factor (Å ²)	38.9	Xtriage
Anisotropy	0.109	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 56.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.054 for -h,-k,l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	734	wwPDB-VP
Average B, all atoms (Å ²)	88.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.29% 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: SO4

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	1.12	0/711	1.75	6/954 (0.6%)

There are no bond length outliers.

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	A	728	SER	CA-C-N	5.40	127.46	120.44
1	A	728	SER	C-N-CA	5.40	127.46	120.44
1	A	779	SER	CA-C-N	5.22	127.73	120.63
1	A	779	SER	C-N-CA	5.22	127.73	120.63
1	A	721	LYS	CA-C-N	5.19	127.50	120.65
1	A	721	LYS	C-N-CA	5.19	127.50	120.65

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	706	0	694	9	0
2	A	5	0	0	0	0
3	A	23	0	0	0	0
All	All	734	0	694	9	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 6.

All (9) 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:759:LEU:O	1:A:763:THR:HG23	1.94	0.67
1:A:702:GLN:HE21	1:A:702:GLN:HA	1.76	0.50
1:A:779:SER:O	1:A:783:GLN:NE2	2.47	0.47
1:A:771:THR:O	1:A:775:THR:HG23	2.16	0.46
1:A:715:HIS:O	1:A:718:GLN:HB3	2.16	0.45
1:A:778:ALA:O	1:A:782:LYS:HG2	2.16	0.44
1:A:702:GLN:HA	1:A:702:GLN:NE2	2.33	0.43
1:A:706:LEU:N	1:A:706:LEU:CD2	2.82	0.43
1:A:767:GLN:O	1:A:771:THR:HG23	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	84/105 (80%)	79 (94%)	5 (6%)	0	100	100

There are no Ramachandran outliers to report.

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	81/98 (83%)	68 (84%)	13 (16%)	2 0

All (13) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	702	GLN
1	A	703	ILE
1	A	706	LEU
1	A	711	ASP
1	A	720	LEU
1	A	724	ASN
1	A	761	ARG
1	A	762	LEU
1	A	766	GLU
1	A	770	ARG
1	A	773	ILE
1	A	782	LYS
1	A	784	LEU

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	699	ASN
1	A	702	GLN
1	A	757	GLN

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 [i](#)

1 ligand is modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. 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 $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	SO4	A	901	-	4,4,4	0.52	0	6,6,6	0.36	0

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	86/105 (81%)	2.74	53 (61%) 0 0	26, 93, 144, 150	0

All (53) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	703	ILE	10.3
1	A	713	LEU	10.2
1	A	784	LEU	7.4
1	A	776	LEU	7.3
1	A	717	LEU	6.9
1	A	774	PHE	6.6
1	A	780	PHE	6.6
1	A	706	LEU	5.3
1	A	710	VAL	5.2
1	A	773	ILE	5.0
1	A	720	LEU	5.0
1	A	716	GLN	4.9
1	A	714	LEU	4.7
1	A	783	GLN	4.6
1	A	781	LYS	4.5
1	A	775	THR	4.3
1	A	771	THR	4.2
1	A	778	ALA	4.0
1	A	762	LEU	3.6
1	A	766	GLU	3.6
1	A	719	LEU	3.5
1	A	758	GLU	3.3
1	A	701	HIS	3.2
1	A	729	MET	3.2
1	A	770	ARG	3.1
1	A	750	GLU	3.1
1	A	764	ASN	3.1

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Mol	Chain	Res	Type	RSRZ
1	A	699	ASN	2.9
1	A	777	ASN	2.9
1	A	763	THR	2.8
1	A	702	GLN	2.8
1	A	731	ILE	2.8
1	A	725	ASN	2.7
1	A	769	HIS	2.7
1	A	707	GLU	2.7
1	A	759	LEU	2.7
1	A	772	GLU	2.6
1	A	760	ALA	2.6
1	A	755	LEU	2.6
1	A	768	SER	2.6
1	A	730	GLN	2.5
1	A	732	SER	2.4
1	A	705	GLU	2.4
1	A	723	GLU	2.4
1	A	704	GLN	2.4
1	A	742	GLU	2.3
1	A	724	ASN	2.3
1	A	726	ARG	2.3
1	A	721	LYS	2.2
1	A	740	ASP	2.2
1	A	782	LYS	2.1
1	A	700	ASN	2.1
1	A	736	LYS	2.1

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	SO4	A	901	5/5	0.97	0.08	35,35,39,39	5

6.5 Other polymers [i](#)

There are no such residues in this entry.