



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 07:17 AM GMT

PDB ID : 3A7O
Title : The crystal structure of the coiled-coil domain of *Saccharomyces cerevisiae* Atg16
Authors : Fujioka, Y.; Noda, N.N.; Inagaki, F.
Deposited on : 2009-10-01
Resolution : 2.50 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
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

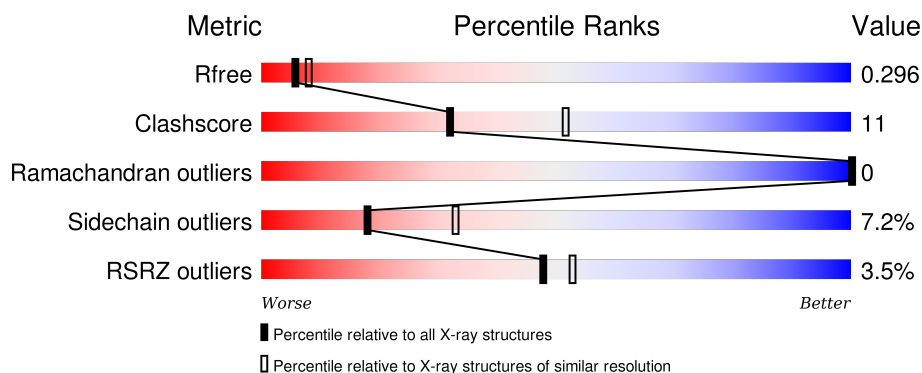
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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.



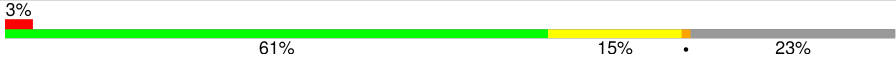
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	91344	3553 (2.50-2.50)
Clashscore	102246	4242 (2.50-2.50)
Ramachandran outliers	100387	4156 (2.50-2.50)
Sidechain outliers	100360	4158 (2.50-2.50)
RSRZ outliers	91569	3562 (2.50-2.50)

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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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	75	<div> <div> <div></div> <div>59%</div> <div>19%</div> <div>•</div> <div>21%</div> </div> </div>
1	B	75	<div> <div> <div>4%</div> <div>61%</div> <div>9%</div> <div>•</div> <div>28%</div> </div> </div>
1	C	75	<div> <div> <div>3%</div> <div>44%</div> <div>23%</div> <div>•</div> <div>29%</div> </div> </div>
1	D	75	<div> <div> <div>3%</div> <div>44%</div> <div>31%</div> <div>•</div> <div>23%</div> </div> </div>
1	E	75	<div> <div> <div>3%</div> <div>64%</div> <div>15%</div> <div>•</div> <div>20%</div> </div> </div>

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Mol	Chain	Length	Quality of chain
1	F	75	 A horizontal bar chart showing the quality of chain F. The bar is divided into four segments: a small red segment at the beginning labeled '3%', followed by a large green segment labeled '61%', a yellow segment labeled '15%', and a final grey segment labeled '23%'. A small black dot is located on the boundary between the yellow and grey segments.

2 Entry composition [i](#)

There are 2 unique types of molecules in this entry. The entry contains 2744 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 protein 16.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
1	A	59	Total	C	N	O	0	0	0
			455	282	81	92			
1	B	54	Total	C	N	O	0	0	0
			427	263	76	88			
1	C	53	Total	C	N	O	0	0	0
			414	256	73	85			
1	D	58	Total	C	N	O	0	0	0
			451	278	81	92			
1	E	60	Total	C	N	O	0	0	0
			452	279	79	94			
1	F	58	Total	C	N	O	0	0	0
			463	286	83	94			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	49	GLY	-	EXPRESSION TAG	UNP Q03818
B	49	GLY	-	EXPRESSION TAG	UNP Q03818
C	49	GLY	-	EXPRESSION TAG	UNP Q03818
D	49	GLY	-	EXPRESSION TAG	UNP Q03818
E	49	GLY	-	EXPRESSION TAG	UNP Q03818
F	49	GLY	-	EXPRESSION TAG	UNP Q03818

- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	9	Total	O	0	0
			9	9		
2	B	17	Total	O	0	0
			17	17		
2	C	7	Total	O	0	0
			7	7		

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	D	12	Total 12	O 12	0	0
2	E	12	Total 12	O 12	0	0
2	F	25	Total 25	O 25	0	0

3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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.

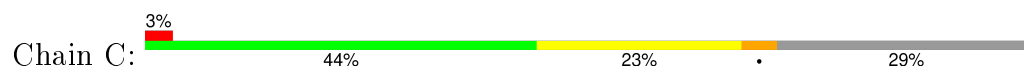
• Molecule 1: Autophagy protein 16



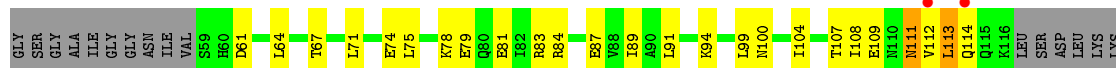
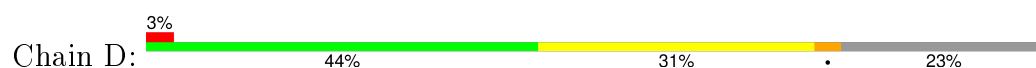
• Molecule 1: Autophagy protein 16



• Molecule 1: Autophagy protein 16



• Molecule 1: Autophagy protein 16

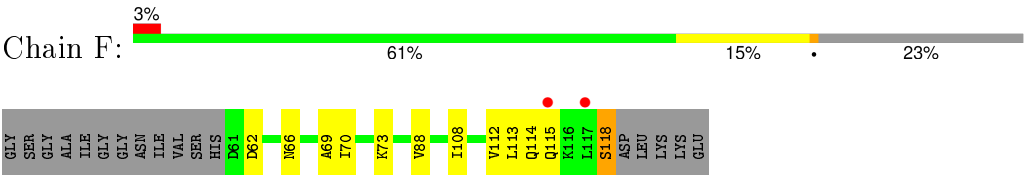


GLU

• Molecule 1: Autophagy protein 16



● Molecule 1: Autophagy protein 16



4 Data and refinement statistics

Property	Value	Source
Space group	P 43 21 2	Depositor
Cell constants a, b, c, α , β , γ	127.89Å 127.89Å 77.81Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	46.08 – 2.50 46.08 – 2.50	Depositor EDS
% Data completeness (in resolution range)	98.4 (46.08-2.50) 98.4 (46.08-2.50)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	12.90 (at 2.51Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.272 , 0.299 0.272 , 0.296	Depositor DCC
R_{free} test set	1094 reflections (4.86%)	DCC
Wilson B-factor (Å ²)	47.2	Xtriage
Anisotropy	0.320	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 37.1	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	0 of 22802 reflections	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	2744	wwPDB-VP
Average B, all atoms (Å ²)	56.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.89% 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, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

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	0.35	0/454	0.50	0/611
1	B	0.33	0/426	0.55	0/573
1	C	0.34	0/414	0.56	0/557
1	D	0.37	0/450	0.53	0/605
1	E	0.35	0/451	0.53	0/608
1	F	0.39	0/462	0.52	0/619
All	All	0.36	0/2657	0.53	0/3573

There are no bond length outliers.

There are no bond angle outliers.

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	455	0	464	11	0
1	B	427	0	440	8	0
1	C	414	0	422	17	0
1	D	451	0	459	16	1
1	E	452	0	449	8	0
1	F	463	0	488	7	0
2	A	9	0	0	0	0
2	B	17	0	0	3	0
2	C	7	0	0	0	0
2	D	12	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	E	12	0	0	2	0
2	F	25	0	0	0	0
All	All	2744	0	2722	57	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 11.

All (57) 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:113:LEU:HB2	1:B:113:LEU:HD23	1.56	0.85
1:D:91:LEU:HA	1:D:94:LYS:HE3	1.64	0.78
1:C:61:ASP:O	1:C:65:LEU:HG	1.89	0.72
1:B:100:ASN:ND2	2:B:127:HOH:O	2.28	0.66
1:E:76:LYS:O	1:E:80:GLN:HG3	1.99	0.62
1:C:106:GLY:O	1:C:110:ASN:HB2	2.00	0.62
1:C:94:LYS:O	1:C:97:GLU:HB2	1.99	0.61
1:E:108:ILE:O	1:E:112:VAL:HG23	2.00	0.61
1:D:84:ARG:HA	1:D:87:GLU:OE1	2.01	0.61
1:A:91:LEU:HD23	1:D:104:ILE:HG23	1.83	0.61
1:F:108:ILE:O	1:F:112:VAL:HG23	2.05	0.57
1:C:69:ALA:O	1:C:73:LYS:HG2	2.07	0.55
1:E:61:ASP:O	1:E:65:LEU:HB2	2.06	0.55
1:D:108:ILE:O	1:D:112:VAL:HG13	2.09	0.53
1:C:92:LYS:HE3	1:D:89:ILE:HG23	1.90	0.53
1:C:104:ILE:CD1	1:F:88:VAL:HG22	2.39	0.53
1:B:98:ARG:HH11	1:B:98:ARG:HG2	1.74	0.52
1:E:80:GLN:NE2	2:E:9:HOH:O	2.44	0.51
1:E:98:ARG:NH2	2:E:8:HOH:O	2.40	0.51
1:E:113:LEU:HB2	1:F:113:LEU:HD13	1.92	0.51
1:F:114:GLN:O	1:F:118:SER:HB2	2.11	0.50
1:A:75:LEU:O	1:A:79:GLU:HG3	2.12	0.49
1:D:100:ASN:O	1:D:104:ILE:HG13	2.12	0.49
1:C:89:ILE:O	1:C:93:ASN:ND2	2.46	0.48
1:C:100:ASN:ND2	1:D:99:LEU:HD21	2.28	0.48
1:A:108:ILE:O	1:A:112:VAL:HG23	2.14	0.47
1:B:79:GLU:HB2	2:B:31:HOH:O	2.14	0.47
1:D:91:LEU:HD11	1:E:103:LEU:HD23	1.96	0.47
1:D:83:ARG:HG3	1:D:83:ARG:HH11	1.79	0.47
1:C:97:GLU:OE1	1:C:97:GLU:HA	2.15	0.47
1:C:70:ILE:HG13	1:C:71:LEU:N	2.30	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:66:ASN:O	1:F:70:ILE:HG13	2.15	0.47
1:D:75:LEU:O	1:D:79:GLU:HG3	2.15	0.46
1:B:83:ARG:HH11	1:B:83:ARG:HG3	1.80	0.46
1:A:113:LEU:CB	1:B:113:LEU:HD23	2.37	0.46
1:C:88:VAL:CG1	1:C:92:LYS:HE2	2.47	0.45
1:A:87:GLU:HA	1:A:87:GLU:OE2	2.17	0.45
1:E:81:GLU:OE1	1:E:81:GLU:HA	2.17	0.45
1:C:70:ILE:O	1:C:73:LYS:HB2	2.17	0.44
1:F:69:ALA:O	1:F:73:LYS:HG3	2.17	0.44
1:C:93:ASN:O	1:C:97:GLU:HG2	2.17	0.44
1:D:109:GLU:O	1:D:112:VAL:HG22	2.18	0.44
1:D:112:VAL:HG23	1:D:113:LEU:N	2.32	0.44
1:D:107:THR:O	1:D:111:ASN:HB2	2.17	0.44
1:C:68:LEU:HD23	1:C:68:LEU:O	2.18	0.44
1:B:109:GLU:HB2	2:B:125:HOH:O	2.18	0.42
1:A:68:LEU:C	1:A:68:LEU:HD23	2.39	0.42
1:D:74:GLU:OE2	1:D:78:LYS:HE2	2.20	0.42
1:D:67:THR:O	1:D:71:LEU:HG	2.19	0.42
1:A:91:LEU:HA	1:A:91:LEU:HD12	1.88	0.41
1:A:80:GLN:HG3	1:A:83:ARG:NH1	2.35	0.41
1:C:87:GLU:O	1:C:91:LEU:HG	2.20	0.41
1:A:98:ARG:O	1:A:102:GLU:HG3	2.20	0.41
1:C:100:ASN:HD21	1:D:99:LEU:HD21	1.85	0.41
1:B:98:ARG:NH1	1:B:98:ARG:HG2	2.35	0.41
1:C:104:ILE:HD11	1:F:88:VAL:HG22	2.01	0.41
1:A:76:LYS:HD2	1:A:76:LYS:HA	1.88	0.41

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)
1:D:114:GLN:OE1	1:D:114:GLN:OE1[7_556]	2.03	0.17

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	57/75 (76%)	57 (100%)	0	0	100	100
1	B	52/75 (69%)	52 (100%)	0	0	100	100
1	C	51/75 (68%)	51 (100%)	0	0	100	100
1	D	56/75 (75%)	56 (100%)	0	0	100	100
1	E	58/75 (77%)	55 (95%)	3 (5%)	0	100	100
1	F	56/75 (75%)	55 (98%)	1 (2%)	0	100	100
All	All	330/450 (73%)	326 (99%)	4 (1%)	0	100	100

There are no Ramachandran outliers to report.

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	49/66 (74%)	46 (94%)	3 (6%)	23	42
1	B	48/66 (73%)	45 (94%)	3 (6%)	22	40
1	C	46/66 (70%)	42 (91%)	4 (9%)	13	24
1	D	49/66 (74%)	44 (90%)	5 (10%)	9	17
1	E	48/66 (73%)	45 (94%)	3 (6%)	22	40
1	F	53/66 (80%)	50 (94%)	3 (6%)	25	46
All	All	293/396 (74%)	272 (93%)	21 (7%)	18	33

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

Mol	Chain	Res	Type
1	A	62	ASP
1	A	91	LEU
1	A	109	GLU
1	B	62	ASP

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Mol	Chain	Res	Type
1	B	65	LEU
1	B	100	ASN
1	C	61	ASP
1	C	68	LEU
1	C	71	LEU
1	C	112	VAL
1	D	61	ASP
1	D	64	LEU
1	D	81	GLU
1	D	111	ASN
1	D	113	LEU
1	E	65	LEU
1	E	75	LEU
1	E	87	GLU
1	F	62	ASP
1	F	115	GLN
1	F	118	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	100	ASN
1	B	72	GLN
1	B	80	GLN
1	B	100	ASN
1	B	114	GLN
1	C	66	ASN
1	C	100	ASN
1	C	110	ASN
1	D	100	ASN
1	E	80	GLN
1	E	95	ASN
1	F	100	ASN
1	F	115	GLN

5.3.3 RNA ⓘ

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 carbohydrates 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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	59/75 (78%)	0.39	1 (1%) 73 76	34, 51, 99, 104	0
1	B	54/75 (72%)	0.44	3 (5%) 28 31	41, 53, 102, 108	0
1	C	53/75 (70%)	0.42	2 (3%) 44 49	47, 58, 97, 102	0
1	D	58/75 (77%)	0.41	2 (3%) 49 54	33, 52, 90, 93	0
1	E	60/75 (80%)	0.25	2 (3%) 50 55	33, 47, 94, 98	0
1	F	58/75 (77%)	0.41	2 (3%) 49 54	29, 44, 92, 95	0
All	All	342/450 (76%)	0.39	12 (3%) 48 53	29, 52, 95, 108	0

All (12) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	F	117	LEU	5.6
1	B	113	LEU	5.4
1	B	112	VAL	4.5
1	D	112	VAL	4.1
1	A	112	VAL	3.5
1	D	114	GLN	2.9
1	E	117	LEU	2.9
1	B	108	ILE	2.4
1	C	111	ASN	2.3
1	E	113	LEU	2.3
1	F	115	GLN	2.1
1	C	112	VAL	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 carbohydrates 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.