



wwPDB X-ray Structure Validation Summary Report ⓘ

May 16, 2020 – 08:25 pm BST

PDB ID : 2HXY
Title : Crystal structure of human apo-eIF4AIII
Authors : Johansen, J.S.; Andersen, G.R.
Deposited on : 2006-08-04
Resolution : 3.30 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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 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	:	2.11
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)	:	2.11

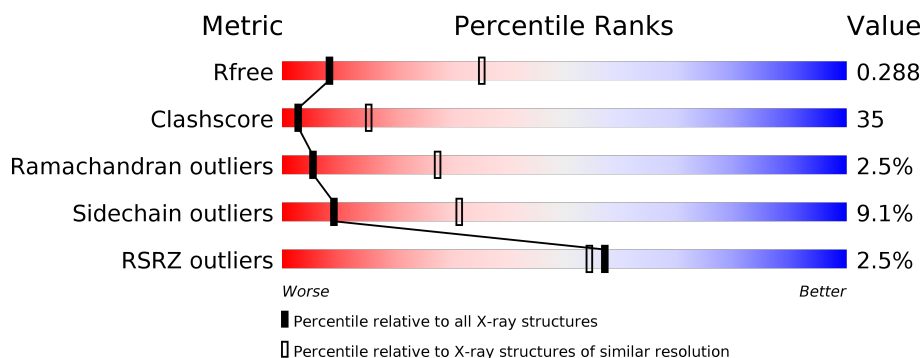
1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.30 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	1149 (3.34-3.26)
Clashscore	141614	1205 (3.34-3.26)
Ramachandran outliers	138981	1183 (3.34-3.26)
Sidechain outliers	138945	1182 (3.34-3.26)
RSRZ outliers	127900	1115 (3.34-3.26)

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 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	391	<div> <div>2%</div> <div> <div></div> <div>47%</div> <div>41%</div> <div>8%</div> <div>••</div> </div> </div>
1	B	391	<div> <div>2%</div> <div> <div></div> <div>45%</div> <div>44%</div> <div>7%</div> <div>•</div> </div> </div>
1	C	391	<div> <div>2%</div> <div> <div></div> <div>46%</div> <div>43%</div> <div>7%</div> <div>•</div> </div> </div>
1	D	391	<div> <div>3%</div> <div> <div></div> <div>45%</div> <div>44%</div> <div>7%</div> <div>•</div> </div> </div>

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 12080 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 Probable ATP-dependent RNA helicase DDX48.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	376	Total	C	N	O	S	0	0	0
			3020	1909	531	562	18			
1	B	376	Total	C	N	O	S	0	0	0
			3020	1909	531	562	18			
1	C	376	Total	C	N	O	S	0	0	0
			3020	1909	531	562	18			
1	D	376	Total	C	N	O	S	0	0	0
			3020	1909	531	562	18			

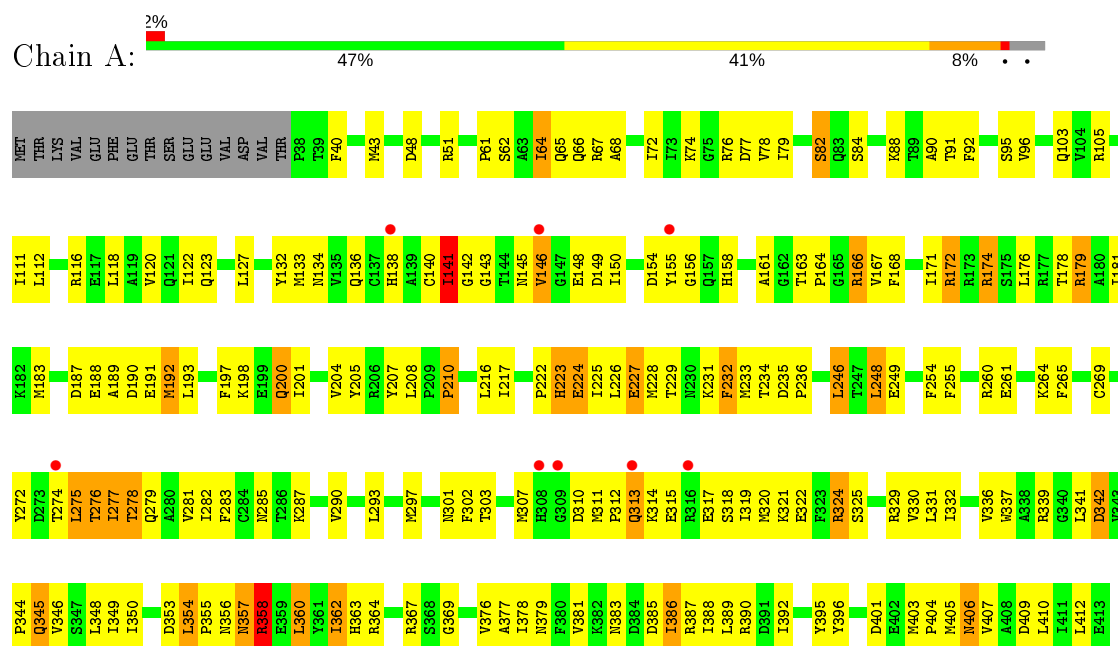
There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	412	LEU	-	CLONING ARTIFACT	UNP P38919
A	413	GLU	-	CLONING ARTIFACT	UNP P38919
B	412	LEU	-	CLONING ARTIFACT	UNP P38919
B	413	GLU	-	CLONING ARTIFACT	UNP P38919
C	412	LEU	-	CLONING ARTIFACT	UNP P38919
C	413	GLU	-	CLONING ARTIFACT	UNP P38919
D	412	LEU	-	CLONING ARTIFACT	UNP P38919
D	413	GLU	-	CLONING ARTIFACT	UNP P38919

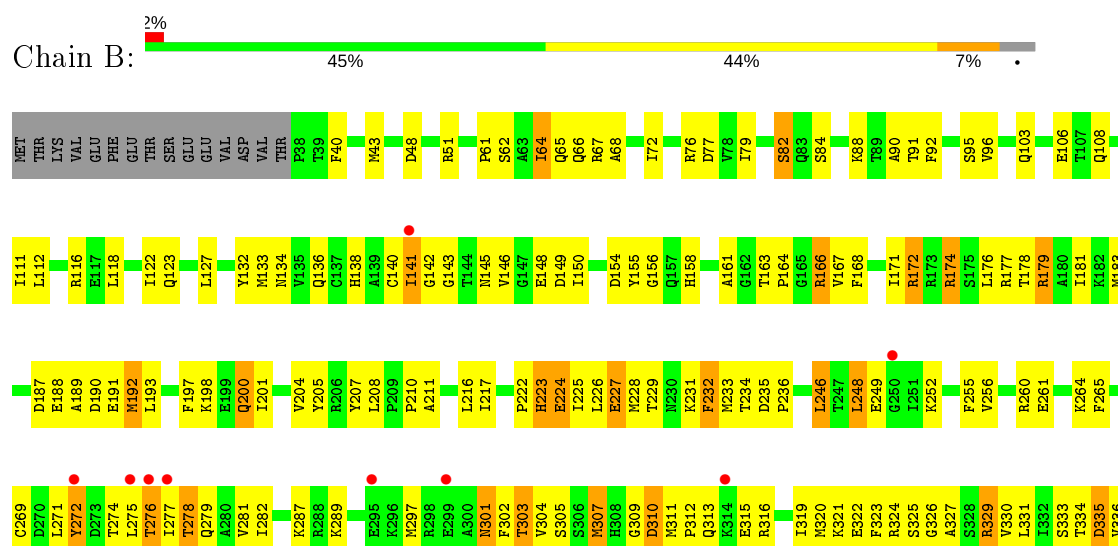
3 Residue-property plots

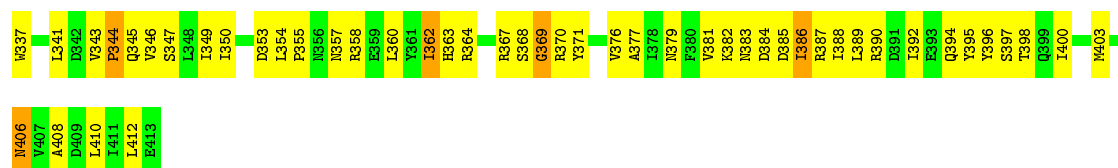
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are 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: Probable ATP-dependent RNA helicase DDX48



- Molecule 1: Probable ATP-dependent RNA helicase DDX48





4 Data and refinement statistics

Property	Value	Source
Space group	P 2 ₁ 2 ₁ 2	Depositor
Cell constants a, b, c, α , β , γ	150.22Å 238.05Å 79.52Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 3.30 49.00 – 3.30	Depositor EDS
% Data completeness (in resolution range)	(Not available) (50.00-3.30) 99.7 (49.00-3.30)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.15	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.49 (at 3.33Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.275 , 0.308 0.262 , 0.288	Depositor DCC
R_{free} test set	1335 reflections (3.06%)	wwPDB-VP
Wilson B-factor (Å ²)	71.3	Xtriage
Anisotropy	0.435	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 69.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.28$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.89	EDS
Total number of atoms	12080	wwPDB-VP
Average B, all atoms (Å ²)	60.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 13.20% 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	0.54	0/3068	0.82	0/4138
1	B	0.53	0/3068	0.83	2/4138 (0.0%)
1	C	0.51	0/3068	0.81	0/4138
1	D	0.51	0/3068	0.80	1/4138 (0.0%)
All	All	0.52	0/12272	0.81	3/16552 (0.0%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	272	TYR	CB-CG-CD1	-8.62	115.83	121.00
1	D	272	TYR	CB-CG-CD1	-6.91	116.85	121.00
1	B	272	TYR	CB-CG-CD2	6.06	124.64	121.00

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	3020	0	3075	213	0
1	B	3020	0	3075	199	1
1	C	3020	0	3075	217	0
1	D	3020	0	3075	237	0
All	All	12080	0	12300	857	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 35.

The worst 5 of 857 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:D:272:TYR:CD1	1:D:275:LEU:HD13	1.81	1.14
1:C:207:TYR:HA	1:D:210:PRO:HG3	1.32	1.10
1:A:312:PRO:HG2	1:A:315:GLU:HB3	1.30	1.09
1:D:276:THR:HG22	1:D:329:ARG:HB3	1.10	1.09
1:D:346:VAL:HG22	1:D:347:SER:H	1.18	1.09

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:B:177:ARG:NH2	1:B:326:GLY:O[2_555]	2.10	0.10

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	374/391 (96%)	309 (83%)	55 (15%)	10 (3%)	5	26
1	B	374/391 (96%)	297 (79%)	68 (18%)	9 (2%)	6	28
1	C	374/391 (96%)	303 (81%)	62 (17%)	9 (2%)	6	28
1	D	374/391 (96%)	300 (80%)	64 (17%)	10 (3%)	5	26
All	All	1496/1564 (96%)	1209 (81%)	249 (17%)	38 (2%)	5	27

5 of 38 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	141	ILE

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Mol	Chain	Res	Type
1	A	406	ASN
1	B	141	ILE
1	B	310	ASP
1	C	141	ILE

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	331/346 (96%)	299 (90%)	32 (10%)	8	29
1	B	331/346 (96%)	304 (92%)	27 (8%)	11	36
1	C	331/346 (96%)	298 (90%)	33 (10%)	7	27
1	D	331/346 (96%)	302 (91%)	29 (9%)	10	33
All	All	1324/1384 (96%)	1203 (91%)	121 (9%)	9	31

5 of 121 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	325	SER
1	C	192	MET
1	D	301	ASN
1	B	362	ILE
1	C	84	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 29 such sidechains are listed below:

Mol	Chain	Res	Type
1	C	103	GLN
1	C	123	GLN
1	D	285	ASN
1	C	108	GLN
1	C	200	GLN

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 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	376/391 (96%)	-0.07	8 (2%) 63 62	4, 45, 113, 151	0
1	B	376/391 (96%)	-0.09	9 (2%) 59 56	6, 52, 129, 173	0
1	C	376/391 (96%)	-0.10	8 (2%) 63 62	3, 49, 126, 165	0
1	D	376/391 (96%)	0.06	13 (3%) 44 42	9, 62, 135, 156	0
All	All	1504/1564 (96%)	-0.05	38 (2%) 57 54	3, 51, 127, 173	0

The worst 5 of 38 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	299	GLU	6.1
1	B	272	TYR	3.9
1	A	313	GLN	3.6
1	C	299	GLU	3.3
1	C	310	ASP	3.3

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.