



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 14, 2026 – 04:12 PM UTC

PDB ID : 4K51 / pdb_00004k51
Title : Crystal Structure of the PCI domain of eIF3a
Authors : Khoshnevis, S.; Neumann, P.; Ficner, R.
Deposited on : 2013-04-12
Resolution : 2.65 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

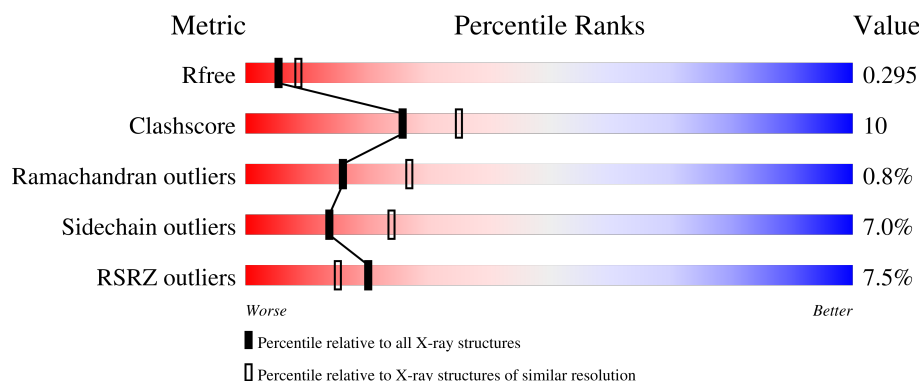
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.65 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	1110 (2.66-2.66)
Clashscore	190562	1141 (2.66-2.66)
Ramachandran outliers	187476	1126 (2.66-2.66)
Sidechain outliers	187428	1126 (2.66-2.66)
RSRZ outliers	180081	1110 (2.66-2.66)

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	224	
1	B	224	

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 3032 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 Eukaryotic translation initiation factor 3 subunit A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	213	Total	C	N	O	S	0	0	0
			1739	1127	273	336	3			
1	B	159	Total	C	N	O	S	0	0	0
			1293	844	199	249	1			

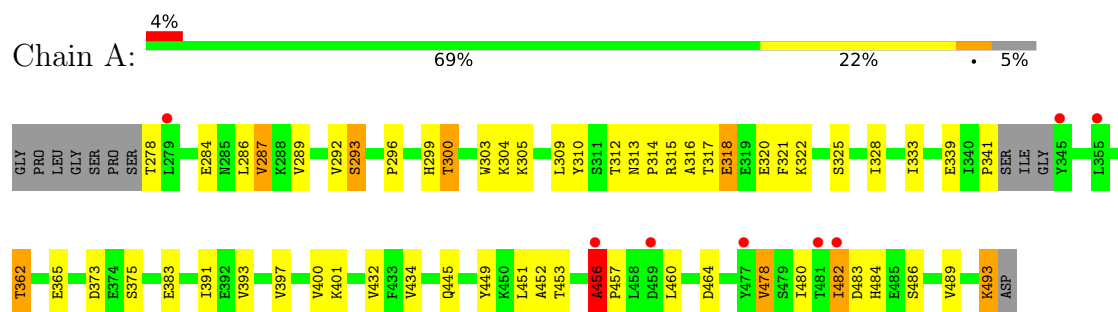
There are 10 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	271	GLY	-	expression tag	UNP P38249
A	272	PRO	-	expression tag	UNP P38249
A	273	LEU	-	expression tag	UNP P38249
A	274	GLY	-	expression tag	UNP P38249
A	275	SER	-	expression tag	UNP P38249
B	271	GLY	-	expression tag	UNP P38249
B	272	PRO	-	expression tag	UNP P38249
B	273	LEU	-	expression tag	UNP P38249
B	274	GLY	-	expression tag	UNP P38249
B	275	SER	-	expression tag	UNP P38249

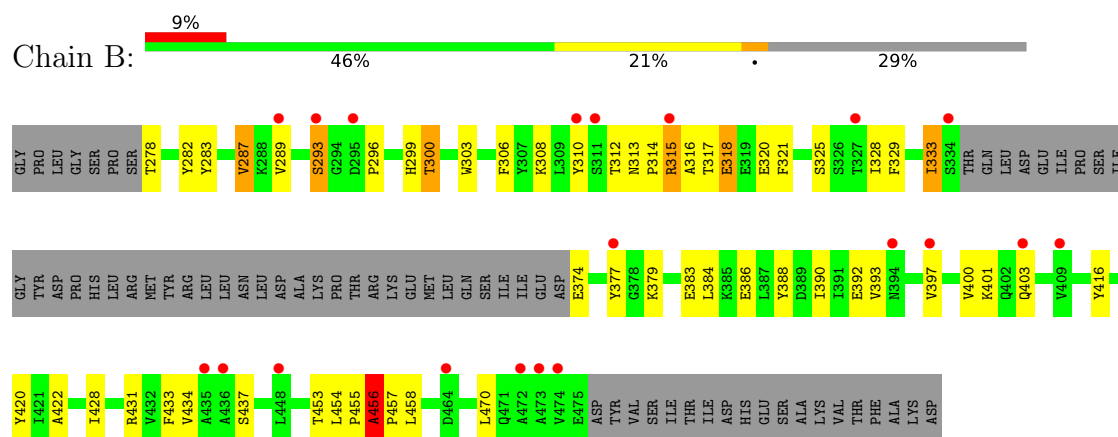
3 Residue-property plots

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: Eukaryotic translation initiation factor 3 subunit A



- Molecule 1: Eukaryotic translation initiation factor 3 subunit A



4 Data and refinement statistics

Property	Value	Source
Space group	I 21 3	Depositor
Cell constants a, b, c, α , β , γ	137.14Å 137.14Å 137.14Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.49 – 2.65 48.49 – 2.65	Depositor EDS
% Data completeness (in resolution range)	99.5 (48.49-2.65) 99.5 (48.49-2.65)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.79 (at 2.65Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8_1069)	Depositor
R, R_{free}	0.249 , 0.286 0.256 , 0.295	Depositor DCC
R_{free} test set	630 reflections (4.98%)	wwPDB-VP
Wilson B-factor (Å ²)	64.6	Xtriage
Anisotropy	0.000	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 81.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	0.053 for -l,-k,-h	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	3032	wwPDB-VP
Average B, all atoms (Å ²)	118.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 13.11% 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	0.45	0/1775	0.92	3/2408 (0.1%)
1	B	0.44	0/1320	1.06	7/1791 (0.4%)
All	All	0.45	0/3095	0.98	10/4199 (0.2%)

There are no bond length outliers.

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	456	ALA	CA-C-N	-11.95	107.32	120.94
1	B	456	ALA	C-N-CA	-11.95	107.32	120.94
1	A	456	ALA	CA-C-N	-11.20	108.17	120.94
1	A	456	ALA	C-N-CA	-11.20	108.17	120.94
1	B	400	VAL	N-CA-C	7.47	117.59	110.42
1	A	400	VAL	N-CA-C	6.69	116.84	110.42
1	B	315	ARG	N-CA-C	-6.02	106.29	113.21
1	B	333	ILE	N-CA-C	-5.19	105.44	110.42
1	B	422	ALA	CA-C-N	-5.08	113.80	119.19
1	B	422	ALA	C-N-CA	-5.08	113.80	119.19

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	1739	0	1746	31	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	B	1293	0	1299	27	0
All	All	3032	0	3045	58	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 10.

All (58) 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:318:GLU:OE1	1:B:379:LYS:NZ	2.27	0.65
1:B:299:HIS:NE2	1:B:420:TYR:OH	2.34	0.61
1:B:401:LYS:HG2	1:B:455:PRO:HD3	1.83	0.60
1:B:289:VAL:O	1:B:293:SER:OG	2.19	0.59
1:A:482:ILE:HG23	1:A:483:ASP:H	1.66	0.59
1:B:433:PHE:O	1:B:437:SER:N	2.33	0.59
1:B:314:PRO:C	1:B:316:ALA:H	2.12	0.56
1:B:283:TYR:HB3	1:B:306:PHE:HB2	1.89	0.55
1:B:303:TRP:HB3	1:B:328:ILE:HG12	1.91	0.52
1:A:303:TRP:HB3	1:A:328:ILE:HG12	1.92	0.51
1:A:478:VAL:HG13	1:A:480:ILE:HG23	1.92	0.51
1:B:296:PRO:O	1:B:300:THR:HG22	2.12	0.49
1:B:390:ILE:HD13	1:B:403:GLN:HB3	1.93	0.49
1:B:310:TYR:OH	1:B:320:GLU:HG3	2.12	0.49
1:A:289:VAL:O	1:A:293:SER:OG	2.29	0.48
1:B:386:GLU:O	1:B:390:ILE:HG13	2.13	0.48
1:A:383:GLU:OE1	1:A:383:GLU:N	2.37	0.48
1:B:283:TYR:O	1:B:287:VAL:HG13	2.14	0.47
1:A:321:PHE:O	1:A:325:SER:OG	2.29	0.47
1:B:384:LEU:HD11	1:B:416:TYR:CE1	2.49	0.47
1:A:362:THR:HG22	1:A:365:GLU:H	1.80	0.47
1:A:456:ALA:CB	1:A:457:PRO:HD3	2.45	0.47
1:B:383:GLU:OE1	1:B:383:GLU:N	2.43	0.46
1:A:373:ASP:OD1	1:A:375:SER:OG	2.22	0.46
1:B:401:LYS:HE3	1:B:453:THR:O	2.15	0.46
1:A:432:VAL:HG11	1:A:452:ALA:HA	1.98	0.46
1:A:305:LYS:HD3	1:A:305:LYS:HA	1.77	0.46
1:B:313:ASN:OD1	1:B:315:ARG:HG2	2.15	0.46
1:A:449:TYR:O	1:A:453:THR:OG1	2.29	0.46
1:A:456:ALA:HB1	1:A:457:PRO:HD3	1.98	0.45
1:A:296:PRO:O	1:A:300:THR:HG22	2.15	0.45
1:B:388:TYR:O	1:B:392:GLU:HB2	2.17	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:318:GLU:O	1:A:321:PHE:HB2	2.17	0.45
1:A:445:GLN:HG2	1:A:489:VAL:HG23	1.98	0.45
1:B:454:LEU:HB2	1:B:458:LEU:HB2	1.99	0.45
1:A:314:PRO:C	1:A:316:ALA:H	2.25	0.44
1:B:310:TYR:CZ	1:B:316:ALA:HA	2.53	0.44
1:A:284:GLU:O	1:A:287:VAL:HG22	2.18	0.44
1:A:296:PRO:O	1:A:299:HIS:HB3	2.18	0.44
1:A:310:TYR:OH	1:A:320:GLU:HG3	2.17	0.44
1:A:333:ILE:CG2	1:A:391:ILE:HD11	2.48	0.43
1:A:309:LEU:O	1:A:312:THR:OG1	2.29	0.43
1:A:401:LYS:NZ	1:A:451:LEU:O	2.48	0.43
1:B:374:GLU:HA	1:B:377:TYR:HB3	2.01	0.42
1:A:322:LYS:HE2	1:A:322:LYS:HB3	1.93	0.42
1:A:482:ILE:O	1:A:484:HIS:N	2.45	0.42
1:A:460:LEU:HD13	1:A:464:ASP:HB3	2.01	0.42
1:B:329:PHE:CE2	1:B:333:ILE:HD11	2.55	0.42
1:A:300:THR:O	1:A:304:LYS:HG3	2.20	0.41
1:B:456:ALA:CB	1:B:457:PRO:HD3	2.50	0.41
1:B:470:LEU:HD23	1:B:470:LEU:HA	1.92	0.41
1:A:313:ASN:OD1	1:A:315:ARG:HG2	2.21	0.41
1:A:286:LEU:HD23	1:A:286:LEU:HA	1.92	0.41
1:B:321:PHE:O	1:B:325:SER:OG	2.27	0.41
1:A:339:GLU:O	1:A:341:PRO:HD3	2.22	0.40
1:A:493:LYS:H	1:A:493:LYS:HG2	1.48	0.40
1:B:308:LYS:O	1:B:312:THR:HG23	2.21	0.40
1:B:428:ILE:O	1:B:431:ARG:HB2	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	209/224 (93%)	193 (92%)	14 (7%)	2 (1%)	12	20
1	B	155/224 (69%)	149 (96%)	5 (3%)	1 (1%)	21	34
All	All	364/448 (81%)	342 (94%)	19 (5%)	3 (1%)	16	27

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	456	ALA
1	B	456	ALA
1	A	482	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	196/204 (96%)	182 (93%)	14 (7%)	13	23
1	B	145/204 (71%)	135 (93%)	10 (7%)	14	24
All	All	341/408 (84%)	317 (93%)	24 (7%)	14	24

All (24) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	278	THR
1	A	287	VAL
1	A	292	VAL
1	A	293	SER
1	A	300	THR
1	A	317	THR
1	A	318	GLU
1	A	362	THR
1	A	393	VAL
1	A	397	VAL
1	A	434	VAL
1	A	478	VAL
1	A	486	SER

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Mol	Chain	Res	Type
1	A	493	LYS
1	B	278	THR
1	B	282	TYR
1	B	287	VAL
1	B	293	SER
1	B	300	THR
1	B	317	THR
1	B	318	GLU
1	B	393	VAL
1	B	397	VAL
1	B	434	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	403	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 oligosaccharides 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	213/224 (95%)	0.51	8 (3%) 44 36	47, 78, 125, 167	0
1	B	159/224 (70%)	1.23	20 (12%) 8 6	92, 146, 215, 242	158 (99%)
All	All	372/448 (83%)	0.82	28 (7%) 20 15	47, 102, 194, 242	158 (42%)

All (28) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	448	LEU	4.1
1	A	459	ASP	3.8
1	A	481	THR	3.7
1	B	472	ALA	3.5
1	A	279	LEU	3.3
1	B	435	ALA	3.0
1	A	482	ILE	2.9
1	B	394	ASN	2.7
1	A	355	LEU	2.7
1	B	377	TYR	2.6
1	B	289	VAL	2.5
1	B	327	THR	2.5
1	B	403	GLN	2.5
1	B	334	SER	2.4
1	B	436	ALA	2.4
1	B	397	VAL	2.3
1	B	409	VAL	2.3
1	B	473	ALA	2.2
1	B	293	SER	2.2
1	B	295	ASP	2.2
1	A	345	TYR	2.2
1	A	456	ALA	2.1
1	B	315	ARG	2.1
1	A	477	TYR	2.1

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Mol	Chain	Res	Type	RSRZ
1	B	310	TYR	2.1
1	B	464	ASP	2.1
1	B	311	SER	2.0
1	B	474	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 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.