



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 10:52 PM UTC

PDB ID : 1KK0 / pdb_00001kk0
Title : Structure of the large gamma subunit of initiation factor eIF2 from *Pyrococcus abyssi*
Authors : Schmitt, E.; Blanquet, S.; Mechulam, Y.
Deposited on : 2001-12-06
Resolution : 1.95 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)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

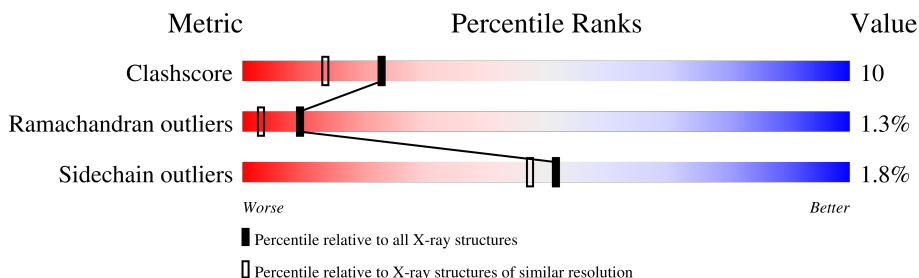
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.95 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))
Clashscore	190562	3612 (1.96-1.96)
Ramachandran outliers	187476	3587 (1.96-1.96)
Sidechain outliers	187428	3587 (1.96-1.96)

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 $\leq 5\%$.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	410	 79% 15% ...

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 3239 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 eIF2gamma.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	399	Total	C	N	O	S	Se	0	0	0
			3040	1929	537	560	6	8			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	15	MSE	MET	modified residue	UNP Q9V1G0
A	97	MSE	MET	modified residue	UNP Q9V1G0
A	100	MSE	MET	modified residue	UNP Q9V1G0
A	107	MSE	MET	modified residue	UNP Q9V1G0
A	130	MSE	MET	modified residue	UNP Q9V1G0
A	211	MSE	MET	modified residue	UNP Q9V1G0
A	307	MSE	MET	modified residue	UNP Q9V1G0
A	363	MSE	MET	modified residue	UNP Q9V1G0

- Molecule 2 is ZINC ION (CCD ID: ZN) (formula: Zn).

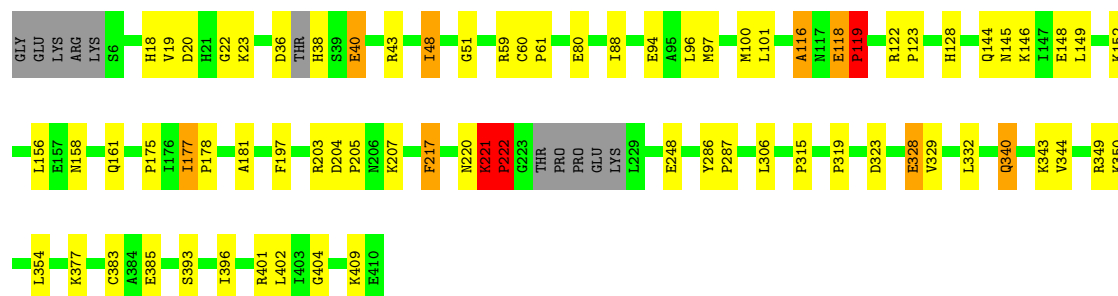
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Zn	0	0
			1	1		

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	198	Total	O	0	0
			198	198		

Note EDS was not executed.

- Chain A:  79% 15% ...



4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	50.87Å 86.42Å 58.21Å 90.00° 109.75° 90.00°	Depositor
Resolution (Å)	47.00 – 1.95	Depositor
% Data completeness (in resolution range)	(Not available) (47.00-1.95)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	0.04	Depositor
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.219 , 0.245	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	3239	wwPDB-VP
Average B, all atoms (Å ²)	38.0	wwPDB-VP

5 Model quality ⓘ

5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: ZN

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	0.38	0/3084	1.00	12/4160 (0.3%)

There are no bond length outliers.

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	40	GLU	N-CA-C	-14.16	94.32	113.30
1	A	221	LYS	CA-C-N	10.41	132.85	119.84
1	A	221	LYS	C-N-CA	10.41	132.85	119.84
1	A	222	PRO	N-CA-C	8.79	130.58	112.47
1	A	48	ILE	N-CA-C	-8.74	95.69	108.71
1	A	221	LYS	N-CA-C	7.82	127.10	109.81
1	A	328	GLU	N-CA-C	-7.05	98.47	109.25
1	A	118	GLU	N-CA-C	5.77	118.63	109.40
1	A	116	ALA	N-CA-C	-5.67	105.85	112.89
1	A	329	VAL	N-CA-C	5.54	116.96	108.71
1	A	323	ASP	N-CA-C	-5.49	106.75	113.50
1	A	177	ILE	N-CA-C	5.19	113.16	107.60

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	3040	0	3133	64	0
2	A	1	0	0	0	0
3	A	198	0	0	2	0
All	All	3239	0	3133	64	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 (64) 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:220:ASN:C	1:A:222:PRO:HD2	1.80	1.04
1:A:97:MSE:HE3	1:A:128:HIS:HD2	1.25	1.01
1:A:220:ASN:O	1:A:222:PRO:HD2	1.69	0.92
1:A:97:MSE:HE1	1:A:128:HIS:HA	1.58	0.84
1:A:328:GLU:HB2	1:A:409:LYS:HE3	1.65	0.77
1:A:145:ASN:ND2	1:A:146:LYS:H	1.83	0.77
1:A:319:PRO:HG2	1:A:383:CYS:SG	2.26	0.76
1:A:61:PRO:HG3	1:A:80:GLU:HG3	1.68	0.75
1:A:97:MSE:HE3	1:A:128:HIS:CD2	2.17	0.72
1:A:332:LEU:HG	1:A:404:GLY:HA2	1.74	0.69
1:A:96:LEU:HG	1:A:100:MSE:HE2	1.76	0.68
1:A:97:MSE:HE2	1:A:100:MSE:HE3	1.75	0.67
1:A:97:MSE:CE	1:A:128:HIS:HD2	2.04	0.67
1:A:396:ILE:HD12	1:A:401:ARG:HD3	1.81	0.62
1:A:220:ASN:C	1:A:222:PRO:CD	2.66	0.61
1:A:40:GLU:HA	1:A:43:ARG:NE	2.16	0.59
1:A:116:ALA:HB2	1:A:144:GLN:HG2	1.87	0.57
1:A:22:GLY:HA3	1:A:145:ASN:OD1	2.07	0.55
1:A:119:PRO:C	1:A:161:GLN:HE22	2.15	0.55
1:A:97:MSE:HE2	1:A:100:MSE:CE	2.37	0.54
1:A:101:LEU:HD12	1:A:402:LEU:HD21	1.89	0.54
1:A:221:LYS:N	1:A:222:PRO:CD	2.69	0.54
1:A:148:GLU:OE2	1:A:149:LEU:HG	2.08	0.53
1:A:340:GLN:CD	1:A:340:GLN:N	2.66	0.53
1:A:96:LEU:HG	1:A:100:MSE:CE	2.38	0.53
1:A:119:PRO:O	1:A:161:GLN:NE2	2.41	0.53
1:A:340:GLN:HG3	1:A:343:LYS:CE	2.40	0.52
1:A:221:LYS:N	1:A:222:PRO:HD2	2.22	0.52
1:A:18:HIS:C	1:A:20:ASP:H	2.17	0.52
1:A:204:ASP:HB3	1:A:207:LYS:HE3	1.91	0.51
1:A:145:ASN:ND2	1:A:146:LYS:N	2.58	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:217:PHE:N	1:A:217:PHE:CD1	2.79	0.50
1:A:340:GLN:HA	1:A:343:LYS:HE2	1.92	0.50
1:A:286:TYR:HB3	1:A:287:PRO:CD	2.43	0.49
1:A:203:ARG:O	1:A:205:PRO:HD3	2.12	0.49
1:A:116:ALA:HA	1:A:158:ASN:ND2	2.28	0.48
1:A:61:PRO:CG	1:A:80:GLU:HG3	2.42	0.47
1:A:349:ARG:O	1:A:350:LYS:HB2	2.14	0.47
1:A:36:ASP:O	1:A:38:HIS:N	2.48	0.47
1:A:248:GLU:OE2	1:A:315:PRO:HG2	2.14	0.47
1:A:354:LEU:HB3	1:A:393:SER:HB2	1.98	0.46
1:A:94:GLU:HA	1:A:94:GLU:OE1	2.14	0.46
1:A:377:LYS:HE2	3:A:1138:HOH:O	2.15	0.46
1:A:122:ARG:CG	1:A:123:PRO:HD2	2.46	0.46
1:A:319:PRO:HB3	1:A:385:GLU:HG2	1.98	0.46
1:A:60:CYS:HA	1:A:61:PRO:HD3	1.83	0.46
1:A:340:GLN:HG3	1:A:343:LYS:HE3	1.98	0.46
1:A:122:ARG:HG3	1:A:123:PRO:HD2	1.98	0.45
1:A:396:ILE:HG22	1:A:396:ILE:O	2.16	0.45
1:A:18:HIS:O	1:A:20:ASP:N	2.49	0.45
1:A:23:LYS:HE3	3:A:1010:HOH:O	2.17	0.45
1:A:51:GLY:O	1:A:88:ILE:HA	2.17	0.45
1:A:286:TYR:HB3	1:A:287:PRO:HD2	1.98	0.45
1:A:175:PRO:HG3	1:A:197:PHE:CZ	2.53	0.43
1:A:96:LEU:CG	1:A:100:MSE:HE2	2.46	0.43
1:A:306:LEU:HD23	1:A:306:LEU:O	2.18	0.43
1:A:152:LYS:NZ	1:A:156:LEU:HD11	2.34	0.43
1:A:118:GLU:HA	1:A:119:PRO:HD3	1.86	0.42
1:A:61:PRO:HG3	1:A:80:GLU:CG	2.42	0.41
1:A:20:ASP:OD1	1:A:20:ASP:C	2.62	0.41
1:A:145:ASN:HD21	1:A:181:ALA:H	1.68	0.41
1:A:349:ARG:HE	1:A:349:ARG:HB2	1.75	0.41
1:A:18:HIS:C	1:A:20:ASP:N	2.79	0.41
1:A:177:ILE:HA	1:A:178:PRO:HD3	1.94	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	393/410 (96%)	373 (95%)	15 (4%)	5 (1%)	9 3

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	19	VAL
1	A	119	PRO
1	A	221	LYS
1	A	222	PRO
1	A	344	VAL

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	327/334 (98%)	321 (98%)	6 (2%)	51 47

All (6) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	48	ILE
1	A	59	ARG
1	A	119	PRO
1	A	217	PHE
1	A	222	PRO
1	A	340	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	62	ASN
1	A	117	ASN
1	A	128	HIS
1	A	144	GLN
1	A	145	ASN
1	A	158	ASN
1	A	161	GLN
1	A	276	GLN
1	A	379	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](#)

Of 1 ligands modelled in this entry, 1 is monoatomic - leaving 0 for Mogul analysis.

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 ⓘ

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

EDS was not executed - this section is therefore empty.

6.2 Non-standard residues in protein, DNA, RNA chains

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates

EDS was not executed - this section is therefore empty.

6.4 Ligands

EDS was not executed - this section is therefore empty.

6.5 Other polymers

EDS was not executed - this section is therefore empty.