



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 7, 2024 – 05:29 pm GMT

PDB ID : 6FC2
Title : Crystal structure of the eIF4E-Eap1p complex from *Saccharomyces cerevisiae*
Authors : Gruener, S.; Valkov, E.
Deposited on : 2017-12-20
Resolution : 1.92 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 : 4.02b-467
Xtriage (Phenix) : 1.13
EDS : 2.36
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) : 2.36

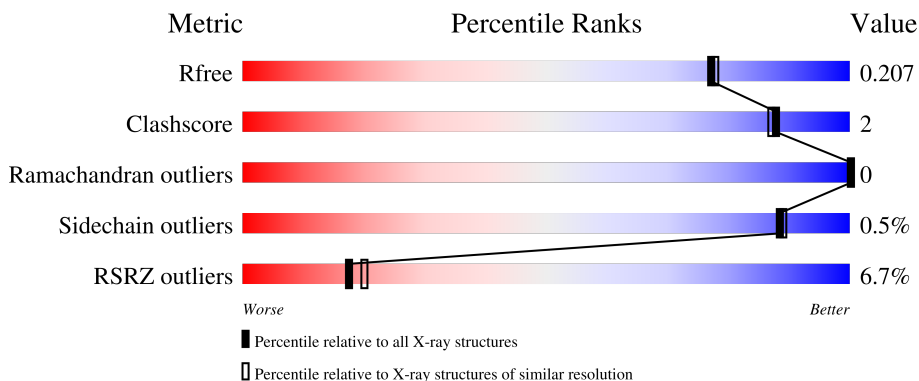
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.92 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	7937 (1.94-1.90)
Clashscore	141614	8644 (1.94-1.90)
Ramachandran outliers	138981	8530 (1.94-1.90)
Sidechain outliers	138945	8530 (1.94-1.90)
RSRZ outliers	127900	7793 (1.94-1.90)

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	183	 7% 89% 9%
1	C	183	 4% 92%
2	B	64	 6% 94% 6%
2	D	64	 14% 91% 6%

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 7904 atoms, of which 3800 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 4E.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	H	N	O	S			
1	A	178	2833	923	1397	243	269	1	0	0	0
1	C	177	2840	924	1399	246	271		0	1	0

There are 14 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	31	GLY	-	expression tag	UNP P07260
A	32	PRO	-	expression tag	UNP P07260
A	33	HIS	-	expression tag	UNP P07260
A	34	MET	-	expression tag	UNP P07260
A	42	ALA	LYS	engineered mutation	UNP P07260
A	168	ALA	LYS	engineered mutation	UNP P07260
A	187	ALA	LYS	engineered mutation	UNP P07260
C	31	GLY	-	expression tag	UNP P07260
C	32	PRO	-	expression tag	UNP P07260
C	33	HIS	-	expression tag	UNP P07260
C	34	MET	-	expression tag	UNP P07260
C	42	ALA	LYS	engineered mutation	UNP P07260
C	168	ALA	LYS	engineered mutation	UNP P07260
C	187	ALA	LYS	engineered mutation	UNP P07260

- Molecule 2 is a protein called Protein EAP1.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	H	N	O	S			
2	B	60	1006	325	502	83	93	3	0	0	0
2	D	60	1006	325	502	83	93	3	0	1	0

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	87	GLY	-	expression tag	UNP P36041
B	88	PRO	-	expression tag	UNP P36041
B	89	HIS	-	expression tag	UNP P36041
B	90	MET	-	expression tag	UNP P36041
D	87	GLY	-	expression tag	UNP P36041
D	88	PRO	-	expression tag	UNP P36041
D	89	HIS	-	expression tag	UNP P36041
D	90	MET	-	expression tag	UNP P36041

- Molecule 3 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total Ca 1 1	0	0

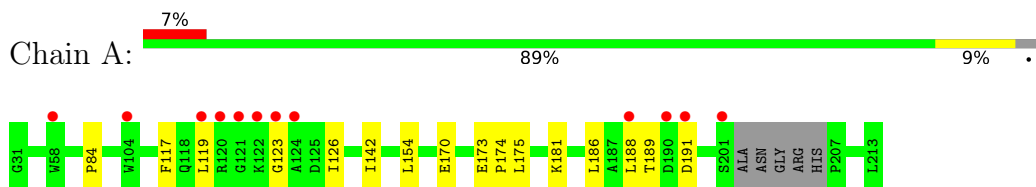
- Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	72	Total O 72 72	0	0
4	B	37	Total O 37 37	0	0
4	C	90	Total O 90 90	0	0
4	D	19	Total O 19 19	0	0

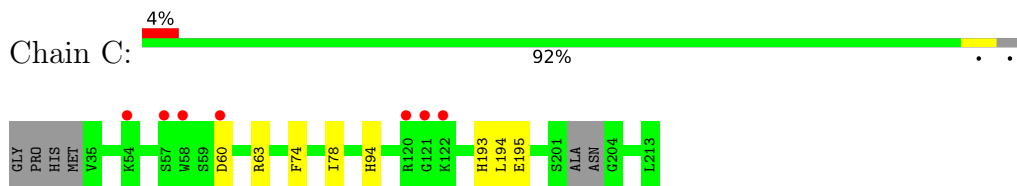
3 Residue-property plots [i](#)

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.

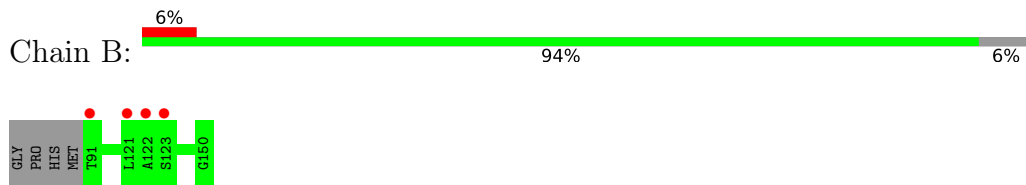
- Molecule 1: Eukaryotic translation initiation factor 4E



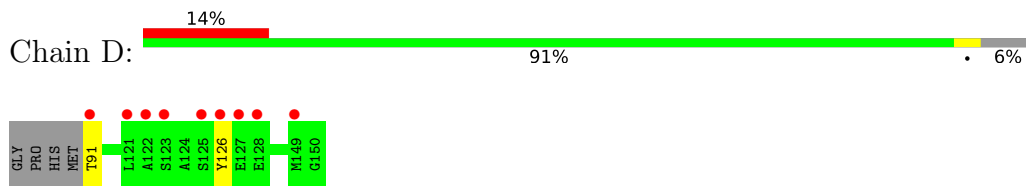
- Molecule 1: Eukaryotic translation initiation factor 4E



- Molecule 2: Protein EAP1



- Molecule 2: Protein EAP1



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	87.57Å 95.12Å 56.01Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.26 – 1.92 48.26 – 1.92	Depositor EDS
% Data completeness (in resolution range)	100.0 (48.26-1.92) 100.0 (48.26-1.92)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.10	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.57 (at 1.92Å)	Xtrriage
Refinement program	PHENIX (1.11_2567: ???)	Depositor
R, R_{free}	0.173 , 0.206 0.173 , 0.207	Depositor DCC
R_{free} test set	1786 reflections (4.90%)	wwPDB-VP
Wilson B-factor (Å ²)	32.2	Xtrriage
Anisotropy	0.325	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.41 , 51.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	7904	wwPDB-VP
Average B, all atoms (Å ²)	49.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.37% 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](#)

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

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	0.32	0/1478	0.51	0/2009
1	C	0.31	0/1482	0.52	0/2014
2	B	0.30	0/518	0.47	0/699
2	D	0.29	0/518	0.49	0/699
All	All	0.31	0/3996	0.51	0/5421

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 [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	1436	1397	1396	11	0
1	C	1441	1399	1397	3	0
2	B	504	502	502	0	0
2	D	504	502	495	2	0
3	A	1	0	0	0	0
4	A	72	0	0	1	0
4	B	37	0	0	0	0
4	C	90	0	0	0	0
4	D	19	0	0	1	0
All	All	4104	3800	3790	15	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 2.

All (15) 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:123:GLY:N	4:A:401:HOH:O	2.24	0.67
1:C:193:HIS:ND1	1:C:195[B]:GLU:OE2	2.32	0.55
1:A:186:LEU:HB2	1:A:188:LEU:CD2	2.45	0.46
1:A:170:GLU:HA	1:A:175:LEU:HD11	1.97	0.45
1:A:126:ILE:HD11	1:A:154:LEU:HD11	1.98	0.45
1:A:84:PRO:HB3	1:A:154:LEU:HD21	1.99	0.44
2:D:91:THR:N	4:D:201:HOH:O	2.50	0.44
1:A:173:GLU:HB2	1:A:174:PRO:HD3	1.99	0.43
1:A:189:THR:HB	1:A:191:ASP:OD2	2.18	0.43
1:A:142:ILE:HG22	1:A:181:LYS:HB3	2.01	0.42
1:A:117:PHE:CZ	1:A:188:LEU:HD11	2.54	0.42
1:C:60:ASP:O	1:C:63:ARG:NH1	2.53	0.41
1:C:74:PHE:CZ	1:C:78:ILE:HG13	2.57	0.40
1:A:123:GLY:HA2	2:D:126:TYR:CD2	2.56	0.40
1:A:117:PHE:CE2	1:A:119:LEU:HD23	2.57	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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	174/183 (95%)	171 (98%)	3 (2%)	0	100	100
1	C	174/183 (95%)	173 (99%)	1 (1%)	0	100	100
2	B	58/64 (91%)	58 (100%)	0	0	100	100
2	D	58/64 (91%)	58 (100%)	0	0	100	100
All	All	464/494 (94%)	460 (99%)	4 (1%)	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	156/159 (98%)	156 (100%)	0	100	100
1	C	156/159 (98%)	154 (99%)	2 (1%)	69	66
2	B	56/59 (95%)	56 (100%)	0	100	100
2	D	56/59 (95%)	56 (100%)	0	100	100
All	All	424/436 (97%)	422 (100%)	2 (0%)	88	89

All (2) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	94	HIS
1	C	194	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	184	GLN
1	C	40	ASN
1	C	118	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 monosaccharides 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 [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	178/183 (97%)	0.18	12 (6%) 17 20	22, 38, 82, 153	0
1	C	177/183 (96%)	0.15	7 (3%) 38 41	24, 33, 95, 140	0
2	B	60/64 (93%)	0.39	4 (6%) 17 20	25, 44, 70, 139	0
2	D	60/64 (93%)	1.00	9 (15%) 2 2	32, 50, 125, 142	0
All	All	475/494 (96%)	0.30	32 (6%) 17 20	22, 38, 96, 153	0

All (32) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	D	122	ALA	13.1
1	C	120	ARG	6.3
1	A	58	TRP	6.2
1	A	123	GLY	5.8
2	D	126	TYR	5.5
1	C	58	TRP	5.4
2	D	125	SER	5.2
2	B	122	ALA	5.0
2	B	91	THR	4.8
2	D	123	SER	4.3
1	C	122	LYS	4.1
1	A	119	LEU	3.7
1	A	122	LYS	3.6
1	C	60	ASP	3.5
2	D	91	THR	3.4
2	D	128	GLU	3.3
1	A	120	ARG	3.3
2	D	121	LEU	3.2
1	A	121	GLY	3.0
1	A	190	ASP	3.0
1	A	124	ALA	2.9

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Mol	Chain	Res	Type	RSRZ
1	A	191	ASP	2.8
2	D	127	GLU	2.7
1	A	104	TRP	2.6
2	B	121	LEU	2.6
1	A	188	LEU	2.5
1	C	121	GLY	2.4
1	C	54	LYS	2.2
1	A	201	SER	2.1
1	C	57	SER	2.0
2	D	149	MET	2.0
2	B	123	SER	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 monosaccharides 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(Å ²)	Q<0.9
3	CA	A	301	1/1	0.99	0.18	25,25,25,25	0

6.5 Other polymers [i](#)

There are no such residues in this entry.