



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 18, 2026 – 10:47 AM UTC

PDB ID : 1F60 / pdb_00001f60
Title : CRYSTAL STRUCTURE OF THE YEAST ELONGATION FACTOR COM-
PLEX EEF1A:EEF1BA
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Deposited on : 2000-06-19
Resolution : 1.67 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

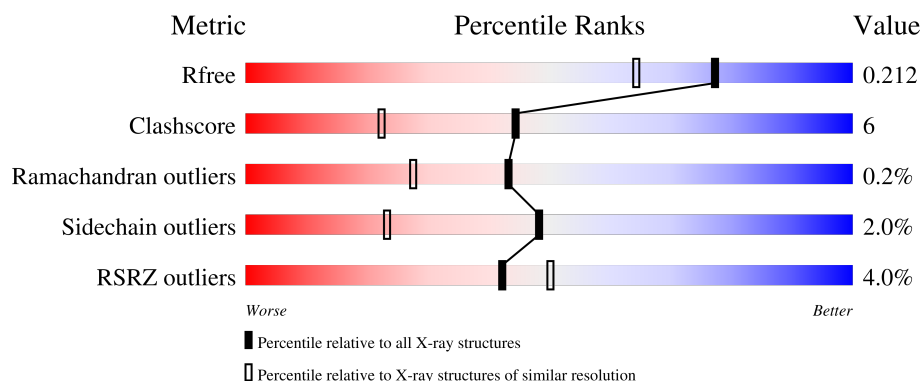
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.67 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	1054 (1.68-1.68)
Clashscore	190562	1078 (1.68-1.68)
Ramachandran outliers	187476	1068 (1.68-1.68)
Sidechain outliers	187428	1067 (1.68-1.68)
RSRZ outliers	180081	1055 (1.68-1.68)

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	458	5% 81% 14% • •
2	B	94	82% 14% •

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 4824 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 ELONGATION FACTOR EEF1A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	440	Total	C	N	O	S	0	0	0
			3388	2152	592	630	14			

- Molecule 2 is a protein called ELONGATION FACTOR EEF1BA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	90	Total	C	N	O	S	0	0	0
			703	441	112	146	4			

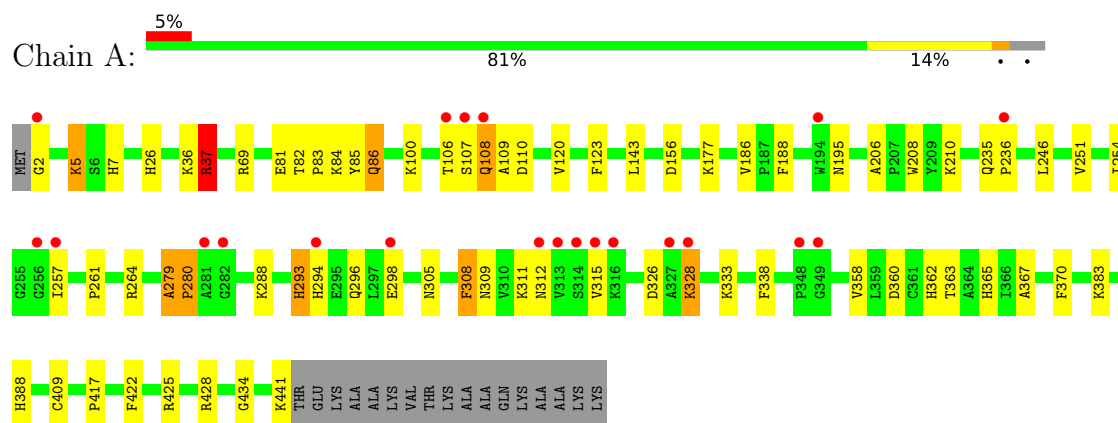
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	567	Total	O	0	0
			567	567		
3	B	166	Total	O	0	0
			166	166		

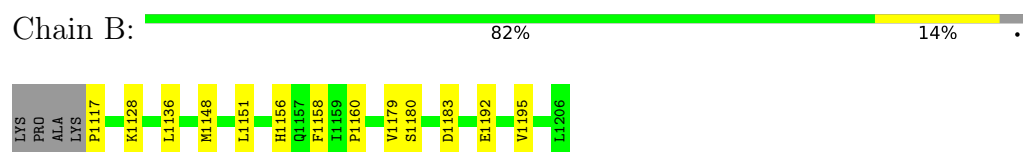
3 Residue-property plots

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: ELONGATION FACTOR EEF1A



• Molecule 2: ELONGATION FACTOR EEF1BA



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	63.85Å 91.81Å 92.98Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 1.67 20.00 – 1.67	Depositor EDS
% Data completeness (in resolution range)	99.3 (20.00-1.67) 99.1 (20.00-1.67)	Depositor EDS
R_{merge}	0.04	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	6.79 (at 1.67Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.187 , 0.221 0.179 , 0.212	Depositor DCC
R_{free} test set	3217 reflections (5.06%)	wwPDB-VP
Wilson B-factor (Å ²)	14.4	Xtriage
Anisotropy	0.334	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 51.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.004 for -h,l,k	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	4824	wwPDB-VP
Average B, all atoms (Å ²)	19.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.47% 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

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	0.79	1/3458 (0.0%)	1.59	35/4674 (0.7%)
2	B	0.78	0/713	1.37	1/964 (0.1%)
All	All	0.79	1/4171 (0.0%)	1.55	36/5638 (0.6%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	2

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	388	HIS	CA-CB	5.59	1.56	1.52

All (36) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	279	ALA	CA-C-O	-14.80	101.32	119.67
1	A	296	GLN	OE1-CD-NE2	-9.68	112.92	122.60
1	A	279	ALA	CA-C-N	9.62	131.86	119.84
1	A	279	ALA	C-N-CA	9.62	131.86	119.84
1	A	280	PRO	N-CA-CB	8.99	112.69	103.25
1	A	280	PRO	CA-N-CD	-8.34	100.33	112.00
1	A	37	ARG	CD-NE-CZ	8.06	135.68	124.40
1	A	5	LYS	CB-CG-CD	7.84	129.33	111.30
1	A	120	VAL	N-CA-C	7.39	117.52	110.42
1	A	246	LEU	O-C-N	-7.25	116.16	121.23
1	A	296	GLN	CG-CD-NE2	7.11	127.07	116.40
1	A	279	ALA	N-CA-C	7.08	120.69	109.64
1	A	280	PRO	N-CD-CG	6.51	112.97	103.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	312	ASN	CA-CB-CG	-6.46	106.14	112.60
1	A	69	ARG	CD-NE-CZ	6.42	133.39	124.40
1	A	246	LEU	CA-C-O	6.41	127.28	121.06
1	A	338	PHE	CA-CB-CG	6.38	120.18	113.80
2	B	1195	VAL	N-CA-C	6.31	117.39	108.36
1	A	326	ASP	CA-CB-CG	6.22	118.82	112.60
1	A	109	ALA	N-CA-C	6.09	119.29	110.28
1	A	186	VAL	N-CA-C	6.01	114.59	107.73
1	A	308	PHE	CA-CB-CG	5.95	119.75	113.80
1	A	370	PHE	CA-CB-CG	5.69	119.49	113.80
1	A	206	ALA	O-C-N	-5.64	116.72	121.12
1	A	85	TYR	N-CA-C	5.53	118.09	109.52
1	A	309	ASN	CA-CB-CG	5.49	118.09	112.60
1	A	86	GLN	OE1-CD-NE2	-5.47	117.13	122.60
1	A	123	PHE	CA-CB-CG	-5.46	108.34	113.80
1	A	279	ALA	CB-CA-C	-5.38	103.12	109.47
1	A	298	GLU	N-CA-C	-5.31	105.94	112.90
1	A	235	GLN	OE1-CD-NE2	5.26	127.86	122.60
1	A	417	PRO	N-CA-C	5.26	117.12	110.70
1	A	309	ASN	CB-CG-ND2	5.21	124.21	116.40
1	A	188	PHE	CA-C-N	-5.16	116.01	123.28
1	A	188	PHE	C-N-CA	-5.16	116.01	123.28
1	A	5	LYS	N-CA-CB	-5.09	102.70	110.85

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	279	ALA	Mainchain,Peptide

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	3388	0	3440	42	0
2	B	703	0	688	8	0
3	A	567	0	0	11	1

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	166	0	0	4	0
All	All	4824	0	4128	49	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 6.

All (49) 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)
2:B:1148:MET:HE2	2:B:1151:LEU:HD22	1.29	1.09
1:A:2:GLY:HA3	3:A:2698:HOH:O	1.71	0.90
1:A:425:ARG:NH2	1:A:428:ARG:H	1.76	0.83
1:A:108:GLN:HG3	3:A:2474:HOH:O	1.82	0.80
1:A:294:HIS:CE1	2:B:1160:PRO:HG2	2.17	0.79
1:A:261:PRO:HD2	1:A:308:PHE:O	1.96	0.66
1:A:288:LYS:HE2	1:A:311:LYS:HG2	1.79	0.64
1:A:362:HIS:CD2	1:A:363:THR:H	2.16	0.64
2:B:1148:MET:CE	2:B:1151:LEU:HD22	2.18	0.63
1:A:251:VAL:HG12	1:A:315:VAL:HG12	1.81	0.61
1:A:333:LYS:HE3	1:A:441:LYS:HD3	1.82	0.61
1:A:37:ARG:HG2	1:A:37:ARG:HH11	1.66	0.60
1:A:36:LYS:HE2	3:A:2545:HOH:O	2.01	0.59
1:A:106:THR:HB	1:A:264:ARG:NH1	2.18	0.58
1:A:107:SER:H	1:A:264:ARG:HH12	1.52	0.57
2:B:1156:HIS:HD2	3:B:2171:HOH:O	1.87	0.56
1:A:365:HIS:HE1	3:A:2333:HOH:O	1.87	0.56
1:A:360:ASP:OD1	1:A:365:HIS:HD2	1.89	0.55
1:A:328:LYS:HD2	3:A:2329:HOH:O	2.06	0.54
1:A:26:HIS:HD2	3:A:2060:HOH:O	1.89	0.54
1:A:26:HIS:HE1	3:A:2042:HOH:O	1.91	0.52
1:A:2:GLY:CA	3:A:2698:HOH:O	2.42	0.52
1:A:156:ASP:OD2	1:A:195:ASN:ND2	2.41	0.52
1:A:362:HIS:HD2	1:A:363:THR:H	1.56	0.51
1:A:293:HIS:HD2	3:B:2515:HOH:O	1.94	0.50
1:A:358:VAL:HA	1:A:367:ALA:HA	1.94	0.50
1:A:362:HIS:HE1	1:A:409:CYS:O	1.95	0.50
1:A:208:TRP:O	1:A:210:LYS:HD2	2.12	0.49
1:A:108:GLN:OE1	1:A:143:LEU:HD13	2.15	0.47
1:A:5:LYS:HE2	1:A:81:GLU:OE1	2.14	0.46
2:B:1136:LEU:HB3	2:B:1158:PHE:CE1	2.52	0.45
2:B:1117:PRO:N	3:B:2582:HOH:O	2.48	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:1180:SER:HB3	2:B:1183:ASP:HB2	1.98	0.45
2:B:1148:MET:CE	2:B:1179:VAL:HG13	2.47	0.45
1:A:26:HIS:CE1	3:A:2042:HOH:O	2.69	0.44
1:A:425:ARG:NH2	1:A:428:ARG:N	2.56	0.43
1:A:254:ILE:CG2	1:A:257:ILE:HD12	2.49	0.43
1:A:84:LYS:HE3	3:A:2306:HOH:O	2.18	0.43
1:A:110:ASP:OD2	1:A:236:PRO:HG3	2.19	0.42
1:A:422:PHE:CZ	1:A:434:GLY:HA3	2.54	0.42
1:A:425:ARG:HG3	3:B:2621:HOH:O	2.19	0.42
1:A:7:HIS:CE1	1:A:86:GLN:OE1	2.73	0.42
1:A:107:SER:H	1:A:264:ARG:NH1	2.17	0.41
1:A:82:THR:HB	1:A:83:PRO:HD2	2.01	0.41
1:A:5:LYS:HE3	1:A:82:THR:O	2.21	0.41
1:A:177:LYS:HE2	3:A:2561:HOH:O	2.19	0.41
1:A:37:ARG:HD3	1:A:37:ARG:HA	1.73	0.41
1:A:425:ARG:C	1:A:425:ARG:HD3	2.46	0.40
1:A:82:THR:HB	1:A:83:PRO:CD	2.51	0.40

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)
3:A:2165:HOH:O	3:A:2307:HOH:O[1_655]	2.11	0.09

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	438/458 (96%)	426 (97%)	11 (2%)	1 (0%)	43	27
2	B	88/94 (94%)	88 (100%)	0	0	100	100
All	All	526/552 (95%)	514 (98%)	11 (2%)	1 (0%)	43	27

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	280	PRO

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	367/379 (97%)	360 (98%)	7 (2%)	50	25
2	B	79/82 (96%)	77 (98%)	2 (2%)	42	16
All	All	446/461 (97%)	437 (98%)	9 (2%)	48	23

All (9) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	37	ARG
1	A	100	LYS
1	A	108	GLN
1	A	293	HIS
1	A	305	ASN
1	A	328	LYS
1	A	383	LYS
2	B	1128	LYS
2	B	1192	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	7	HIS
1	A	26	HIS
1	A	86	GLN
1	A	293	HIS
1	A	294	HIS
1	A	296	GLN
1	A	362	HIS

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Mol	Chain	Res	Type
1	A	365	HIS
1	A	388	HIS
2	B	1142	ASN
2	B	1156	HIS
2	B	1169	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 [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	440/458 (96%)	-0.07	21 (4%) 35 44	8, 15, 35, 45	0
2	B	90/94 (95%)	-0.46	0 100 100	9, 14, 23, 29	0
All	All	530/552 (96%)	-0.14	21 (3%) 42 50	8, 15, 33, 45	0

All (21) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	106	THR	4.8
1	A	294	HIS	4.7
1	A	315	VAL	4.5
1	A	107	SER	4.1
1	A	194	TRP	3.1
1	A	313	VAL	2.9
1	A	328	LYS	2.8
1	A	348	PRO	2.8
1	A	349	GLY	2.7
1	A	316	LYS	2.7
1	A	282	GLY	2.7
1	A	298	GLU	2.6
1	A	236	PRO	2.6
1	A	312	ASN	2.5
1	A	256	GLY	2.4
1	A	327	ALA	2.4
1	A	2	GLY	2.3
1	A	281	ALA	2.3
1	A	108	GLN	2.1
1	A	257	ILE	2.0
1	A	314	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 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.