



Full wwPDB X-ray Structure Validation Report ⓘ

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PDB ID : 7XTP
Title : eIF4E in Complex with a Disulphide-Free Autonomous VH Domain
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Deposited on : 2022-05-17
Resolution : 1.83 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
Mogul : ?? (??), CSD ??CSD?? (????)
Xtriage (Phenix) : 1.13
EDS : 2.29
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) : 2.29

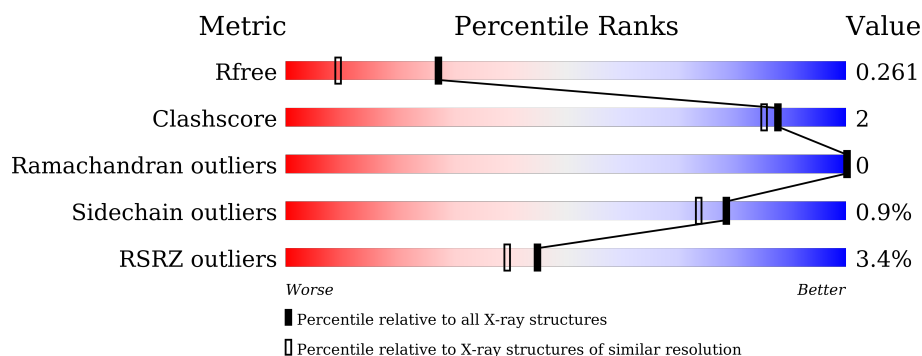
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.83 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	7484 (1.84-1.80)
Clashscore	141614	8401 (1.84-1.80)
Ramachandran outliers	138981	8290 (1.84-1.80)
Sidechain outliers	138945	8290 (1.84-1.80)
RSRZ outliers	127900	7371 (1.84-1.80)

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	217	<div> <div>6%</div> <div> <div></div> <div>75%</div> <div>5%</div> <div>20%</div> </div> </div>
1	B	217	<div> <div>3%</div> <div> <div></div> <div>73%</div> <div>7%</div> <div>20%</div> </div> </div>
2	C	154	<div> <div></div> <div> <div></div> <div>84%</div> <div>.</div> <div>13%</div> </div> </div>
2	D	154	<div> <div></div> <div> <div></div> <div>81%</div> <div>5%</div> <div>14%</div> </div> </div>

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 5357 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.

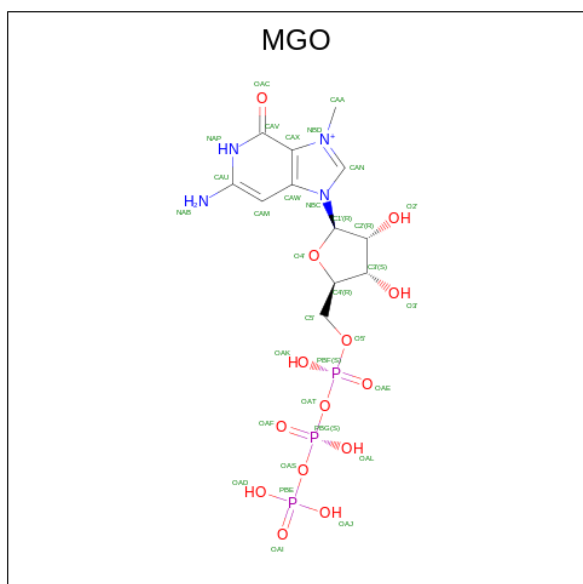
- Molecule 1 is a protein called Eukaryotic translation initiation factor 4E.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	173	Total	C	N	O	S	0	2	0
			1447	926	255	260	6			
1	B	174	Total	C	N	O	S	0	1	0
			1447	923	254	264	6			

- Molecule 2 is a protein called VH-S4ss.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	134	Total	C	N	O	S	0	5	0
			1012	628	174	205	5			
2	D	132	Total	C	N	O	S	0	6	0
			1005	627	172	201	5			

- Molecule 3 is [[(2R,3S,4R,5R)-5-(6-AMINO-3-METHYL-4-OXO-5H-IMIDAZO[4,5-C]PYRIDIN-1-YL)-3,4-DIHYDROXY-OXOLAN-2-YL]METHOXY-HYDROXY-PHOSPHORYL] PHOSPHONO HYDROGEN PHOSPHATE (three-letter code: MGO) (formula: C₁₂H₂₀N₄O₁₄P₃).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	P	0	0
			29	12	4	11	2		
3	B	1	Total	C	N	O	P	0	0
			29	12	4	11	2		

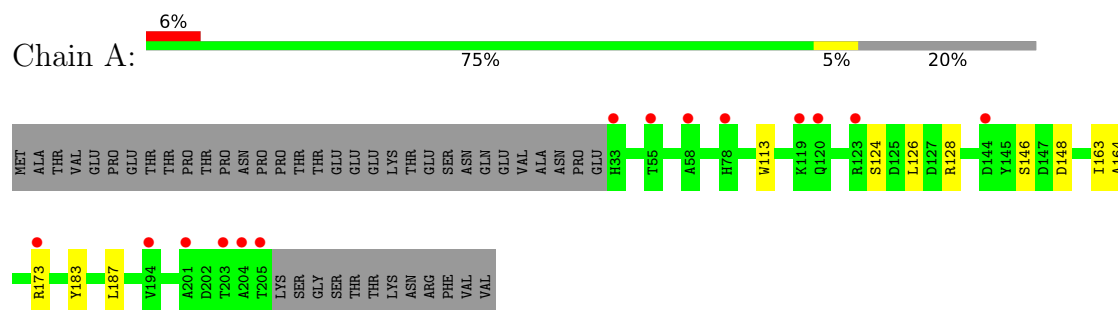
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	88	Total	O	0	0
			88	88		
4	B	102	Total	O	0	0
			102	102		
4	C	95	Total	O	0	0
			95	95		
4	D	103	Total	O	0	0
			103	103		

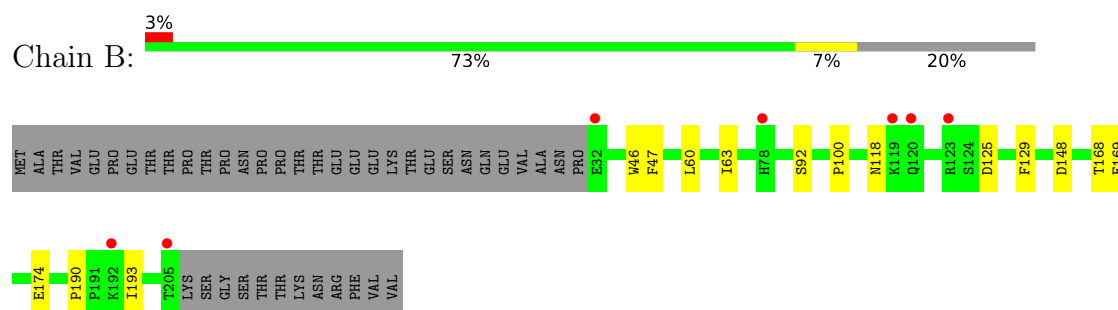
3 Residue-property plots [i](#)

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.

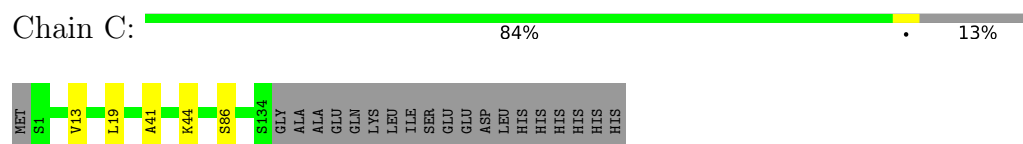
- Molecule 1: Eukaryotic translation initiation factor 4E



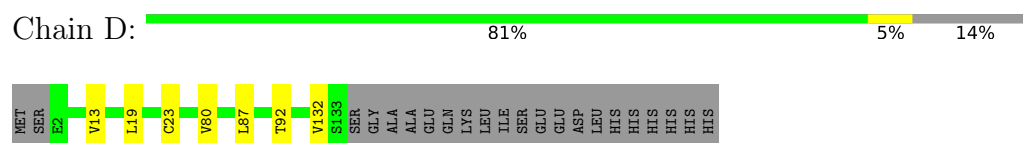
- Molecule 1: Eukaryotic translation initiation factor 4E



- Molecule 2: VH-S4ss



- Molecule 2: VH-S4ss



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	75.62Å 57.21Å 78.96Å 90.00° 93.95° 90.00°	Depositor
Resolution (Å)	46.33 – 1.83 46.29 – 1.83	Depositor EDS
% Data completeness (in resolution range)	99.1 (46.33-1.83) 99.1 (46.29-1.83)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.66 (at 1.83Å)	Xtriage
Refinement program	REFMAC 5.8.0267	Depositor
R, R_{free}	0.226 , 0.259 0.231 , 0.261	Depositor DCC
R_{free} test set	3060 reflections (5.16%)	wwPDB-VP
Wilson B-factor (Å ²)	19.6	Xtriage
Anisotropy	0.331	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 35.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.000 for l,-k,h	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	5357	wwPDB-VP
Average B, all atoms (Å ²)	24.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 42.13 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 2.1339e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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: MGO

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	0.62	0/1491	0.69	0/2018
1	B	0.63	0/1485	0.68	0/2011
2	C	0.67	0/1041	0.73	0/1411
2	D	0.66	0/1037	0.73	0/1407
All	All	0.64	0/5054	0.70	0/6847

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	1447	0	1421	4	0
1	B	1447	0	1402	7	0
2	C	1012	0	1000	3	0
2	D	1005	0	998	7	0
3	A	29	0	16	0	0
3	B	29	0	16	0	0
4	A	88	0	0	0	0
4	B	102	0	0	0	0
4	C	95	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	D	103	0	0	0	0
All	All	5357	0	4853	20	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 (20) 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:D:92[A]:THR:HG22	2:D:132:VAL:H	1.30	0.94
2:D:23[A]:CYS:HB3	2:D:80[A]:VAL:HG22	1.56	0.86
2:D:23[A]:CYS:HB3	2:D:80[A]:VAL:CG2	2.12	0.78
1:B:190:PRO:HD2	1:B:193:ILE:HD12	1.80	0.63
2:D:13:VAL:HG21	2:D:87:LEU:HD13	1.93	0.49
2:D:23[A]:CYS:CB	2:D:80[A]:VAL:HG22	2.37	0.49
1:B:174:GLU:HG3	2:C:86:SER:HB3	1.95	0.49
2:D:13:VAL:HG11	2:D:19:LEU:HG	1.97	0.46
1:A:183:TYR:CZ	1:A:187:LEU:HD11	2.50	0.46
1:A:113:TRP:O	1:A:164:ALA:HA	2.15	0.45
1:B:47:PHE:HB2	1:B:63:ILE:HD11	1.98	0.45
1:A:124:SER:O	1:A:128[A]:ARG:NH2	2.50	0.45
2:D:23[A]:CYS:HB3	2:D:80[A]:VAL:HG23	1.98	0.44
2:C:13:VAL:HG11	2:C:19:LEU:HG	1.99	0.44
1:B:46:TRP:HB2	1:B:92:SER:HB2	2.00	0.43
1:B:125:ASP:HB3	1:B:129:PHE:CE2	2.54	0.42
1:B:168:THR:HG23	1:B:169:GLU:HG2	2.03	0.41
2:C:41:ALA:HB3	2:C:44:LYS:HG3	2.04	0.40
1:B:60:LEU:HD21	1:B:100:PRO:HG2	2.03	0.40
1:A:126:LEU:HD12	1:A:163:ILE:HD11	2.04	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	173/217 (80%)	168 (97%)	5 (3%)	0	100	100
1	B	173/217 (80%)	168 (97%)	5 (3%)	0	100	100
2	C	137/154 (89%)	135 (98%)	2 (2%)	0	100	100
2	D	136/154 (88%)	135 (99%)	1 (1%)	0	100	100
All	All	619/742 (83%)	606 (98%)	13 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	156/195 (80%)	153 (98%)	3 (2%)	57	45
1	B	155/195 (80%)	153 (99%)	2 (1%)	69	61
2	C	116/128 (91%)	116 (100%)	0	100	100
2	D	115/128 (90%)	115 (100%)	0	100	100
All	All	542/646 (84%)	537 (99%)	5 (1%)	78	74

All (5) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	146	SER
1	A	148	ASP
1	A	173	ARG
1	B	118	ASN
1	B	148	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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](#)

2 ligands are modelled in this entry.

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	173/217 (79%)	0.32	14 (8%) 12 9	15, 26, 45, 58	0
1	B	174/217 (80%)	0.27	7 (4%) 38 32	15, 24, 41, 49	0
2	C	134/154 (87%)	-0.04	0 100 100	15, 20, 29, 40	0
2	D	132/154 (85%)	0.03	0 100 100	14, 18, 27, 38	0
All	All	613/742 (82%)	0.17	21 (3%) 45 39	14, 22, 40, 58	0

All (21) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	201	ALA	4.9
1	A	123	ARG	4.2
1	A	205	THR	4.2
1	A	204	ALA	3.5
1	B	192	LYS	3.2
1	A	173	ARG	3.1
1	A	194	VAL	3.1
1	A	119	LYS	2.6
1	B	78	HIS	2.5
1	A	33	HIS	2.5
1	B	123	ARG	2.5
1	A	78	HIS	2.4
1	B	32	GLU	2.4
1	B	205	THR	2.4
1	B	120	GLN	2.4
1	B	119	LYS	2.3
1	A	58	ALA	2.2
1	A	55	THR	2.2
1	A	144	ASP	2.1
1	A	120	GLN	2.1
1	A	203	THR	2.1

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(\AA^2)	Q<0.9
3	MGO	B	301	29/33	0.92	0.09	25,30,36,36	0
3	MGO	A	301	29/33	0.95	0.09	25,27,29,29	0

6.5 Other polymers [i](#)

There are no such residues in this entry.