



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 7, 2016 – 04:48 AM EDT

PDB ID : 5K4B
Title : Structure of eukaryotic translation initiation factor 3 subunit D (eIF3d) cap binding domain from *Nasonia vitripennis*, Crystal form 1
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Deposited on : 2016-05-20
Resolution : 1.40 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

<http://wwpdb.org/validation/2016/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4.02b-467
Mogul	:	unknown
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20027939
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	rb-20027939

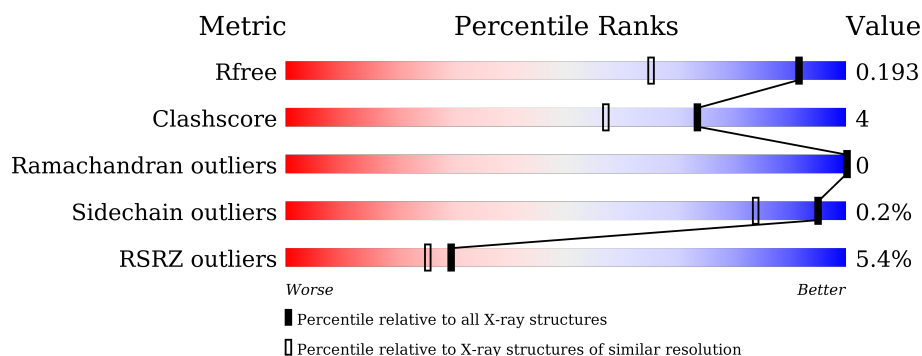
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.40 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}	91344	1199 (1.40-1.40)
Clashscore	102246	1295 (1.40-1.40)
Ramachandran outliers	100387	1259 (1.40-1.40)
Sidechain outliers	100360	1258 (1.40-1.40)
RSRZ outliers	91569	1198 (1.40-1.40)

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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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	368	<div> <div>5%</div> <div> <div></div> <div>92%</div> <div>6% •</div> </div> </div>
1	B	368	<div> <div>6%</div> <div> <div></div> <div>87%</div> <div>11% •</div> </div> </div>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 6750 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 D.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	361	Total	C	N	O	S	0	0	0
			2917	1834	505	563	15			
1	B	361	Total	C	N	O	S	0	0	0
			2920	1836	505	564	15			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	170	GLY	-	expression tag	UNP K7IM66
A	171	SER	-	expression tag	UNP K7IM66
B	170	GLY	-	expression tag	UNP K7IM66
B	171	SER	-	expression tag	UNP K7IM66

- Molecule 2 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	1	Total	Cl	0	0
			1	1		
2	A	2	Total	Cl	0	0
			2	2		

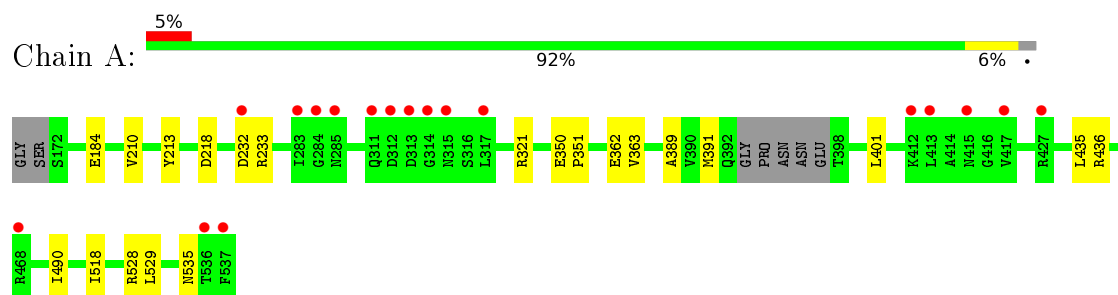
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	510	Total	O	0	0
			510	510		
3	B	400	Total	O	0	0
			400	400		

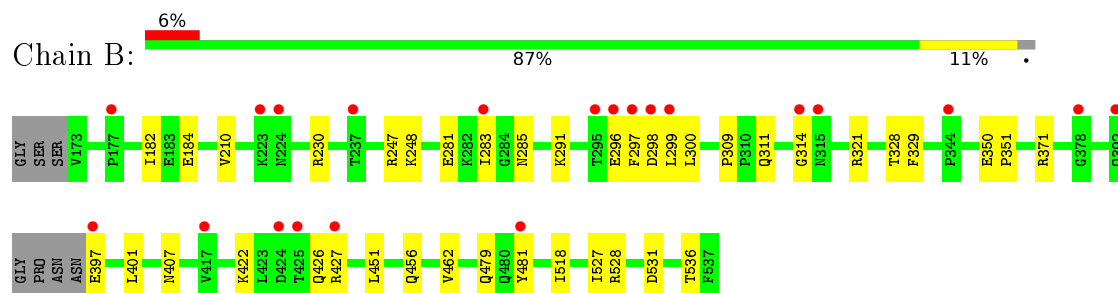
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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 ($\text{RSRZ} > 2$). 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 D



- Molecule 1: Eukaryotic translation initiation factor 3 subunit D



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	61.83Å 62.91Å 192.11Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.03 – 1.40 48.03 – 1.40	Depositor EDS
% Data completeness (in resolution range)	99.9 (48.03-1.40) 99.9 (48.03-1.40)	Depositor EDS
R_{merge}	0.03	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.50 (at 1.40Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.9_1692)	Depositor
R, R_{free}	0.174 , 0.193 0.175 , 0.193	Depositor DCC
R_{free} test set	2357 reflections (1.59%)	DCC
Wilson B-factor (Å ²)	15.4	Xtriage
Anisotropy	0.397	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 40.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.27$	Xtriage
Estimated twinning fraction	0.035 for k,h,-l	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	6750	wwPDB-VP
Average B, all atoms (Å ²)	23.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.05% 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: CL

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	0.39	0/2973	0.59	0/4027
1	B	0.32	0/2976	0.57	0/4031
All	All	0.36	0/5949	0.58	0/8058

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	2917	0	2874	18	0
1	B	2920	0	2875	33	0
2	A	2	0	0	0	0
2	B	1	0	0	0	0
3	A	510	0	0	8	0
3	B	400	0	0	14	1
All	All	6750	0	5749	47	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 4.

All (47) 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:218:ASP:OD1	3:A:701:HOH:O	1.84	0.96
1:A:233:ARG:NH1	3:A:702:HOH:O	2.01	0.90
1:B:397:GLU:OE1	3:B:701:HOH:O	1.90	0.90
1:B:299:LEU:O	3:B:702:HOH:O	1.94	0.85
1:A:436:ARG:NH1	3:A:703:HOH:O	2.18	0.77
1:A:213:TYR:OH	3:A:701:HOH:O	1.98	0.73
1:B:518:ILE:HD11	1:B:527:ILE:HD12	1.68	0.73
1:B:296:GLU:OE1	3:B:704:HOH:O	2.04	0.73
1:B:248:LYS:NZ	3:B:706:HOH:O	2.21	0.73
1:B:297:PHE:N	3:B:707:HOH:O	2.23	0.72
1:A:184:GLU:HG2	1:A:528:ARG:HG2	1.75	0.69
1:B:481:TYR:OH	3:B:705:HOH:O	2.14	0.64
1:B:296:GLU:HB3	1:B:299:LEU:HG	1.81	0.62
1:B:397:GLU:N	3:B:701:HOH:O	2.31	0.62
1:B:184:GLU:HG2	1:B:528:ARG:HG2	1.81	0.62
1:A:210:VAL:HG22	1:A:389:ALA:HA	1.84	0.60
1:A:362:GLU:OE1	1:B:247:ARG:NH1	2.32	0.59
1:B:300:LEU:N	3:B:703:HOH:O	2.32	0.58
1:A:535:ASN:ND2	3:A:713:HOH:O	2.37	0.57
1:B:422:LYS:HD2	1:B:426:GLN:OE1	2.03	0.57
1:B:314:GLY:HA2	3:B:713:HOH:O	2.06	0.55
1:B:328:THR:OG1	3:B:703:HOH:O	2.01	0.55
1:A:232:ASP:O	1:B:247:ARG:NH2	2.41	0.54
1:A:321:ARG:NH1	1:B:536:THR:O	2.40	0.54
1:B:230:ARG:HG2	1:B:329:PHE:HZ	1.73	0.53
1:A:363:VAL:HG22	3:A:988:HOH:O	2.09	0.52
1:B:350:GLU:HG3	1:B:351:PRO:HD2	1.90	0.51
1:A:435:LEU:HD21	1:A:490:ILE:HA	1.93	0.51
1:B:285:ASN:HB3	3:B:768:HOH:O	2.11	0.50
1:A:350:GLU:HG3	1:A:351:PRO:HD2	1.93	0.50
1:B:321:ARG:NH2	3:B:720:HOH:O	2.40	0.50
1:A:436:ARG:NH2	3:A:709:HOH:O	2.30	0.49
1:B:401:LEU:HD22	1:B:456:GLN:HB2	1.93	0.49
1:B:291:LYS:NZ	1:B:298:ASP:OD1	2.44	0.49
1:A:362:GLU:HB3	1:B:247:ARG:HD3	1.95	0.48
1:B:427:ARG:CZ	1:B:479:GLN:HG2	2.45	0.47
1:A:321:ARG:HG3	3:A:740:HOH:O	2.16	0.46
1:B:230:ARG:HG2	1:B:329:PHE:CZ	2.51	0.45
1:B:427:ARG:HG2	3:B:714:HOH:O	2.17	0.45
1:A:391:MET:HB2	1:A:401:LEU:HD11	2.00	0.44
1:B:281:GLU:HG2	1:B:283:ILE:HG13	2.00	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:182:ILE:HD11	1:B:531:ASP:HB2	2.00	0.43
1:B:371:ARG:NE	1:B:451:LEU:O	2.49	0.42
1:B:407:ASN:HA	1:B:462:VAL:O	2.19	0.42
1:A:518:ILE:HG23	1:A:529:LEU:HD13	2.02	0.41
1:B:309:PRO:O	1:B:311:GLN:HG2	2.21	0.41
1:B:300:LEU:HA	3:B:702:HOH:O	2.21	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:B:836:HOH:O	3:B:1060:HOH:O[4_556]	2.10	0.10

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	357/368 (97%)	346 (97%)	11 (3%)	0	100	100
1	B	357/368 (97%)	345 (97%)	12 (3%)	0	100	100
All	All	714/736 (97%)	691 (97%)	23 (3%)	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	326/331 (98%)	326 (100%)	0	100	100
1	B	326/331 (98%)	325 (100%)	1 (0%)	94	83
All	All	652/662 (98%)	651 (100%)	1 (0%)	95	84

All (1) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	210	VAL

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

Mol	Chain	Res	Type
1	A	535	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

Of 3 ligands modelled in this entry, 3 are 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

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	361/368 (98%)	0.42	18 (4%) 32 29	8, 17, 38, 76	0
1	B	361/368 (98%)	0.39	21 (5%) 26 23	12, 21, 43, 73	0
All	All	722/736 (98%)	0.40	39 (5%) 29 25	8, 19, 41, 76	0

All (39) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	313	ASP	6.4
1	B	296	GLU	5.7
1	A	285	ASN	5.3
1	B	315	ASN	4.9
1	B	177	PRO	4.6
1	B	299	LEU	4.5
1	A	314	GLY	4.1
1	B	297	PHE	3.9
1	A	537	PHE	3.4
1	A	315	ASN	3.4
1	A	284	GLY	3.3
1	A	417	VAL	3.2
1	B	425	THR	3.0
1	A	312	ASP	3.0
1	A	283	ILE	2.9
1	B	417	VAL	2.9
1	B	223	LYS	2.9
1	A	232	ASP	2.8
1	A	415	ASN	2.8
1	A	536	THR	2.7
1	A	317	LEU	2.7
1	A	413	LEU	2.7
1	A	468	ARG	2.7
1	B	237	THR	2.7

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Mol	Chain	Res	Type	RSRZ
1	A	311	GLN	2.5
1	B	392	GLN	2.5
1	B	397	GLU	2.4
1	B	224	ASN	2.4
1	A	412	LYS	2.3
1	B	344	PRO	2.3
1	B	481	TYR	2.3
1	A	427	ARG	2.2
1	B	283	ILE	2.1
1	B	295	THR	2.1
1	B	298	ASP	2.1
1	B	424	ASP	2.1
1	B	427	ARG	2.0
1	B	314	GLY	2.0
1	B	378	GLY	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 carbohydrates 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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(Å ²)	Q<0.9
2	CL	A	601	1/1	0.97	0.06	-1.39	54,54,54,54	0
2	CL	B	601	1/1	0.98	0.05	-1.50	22,22,22,22	0
2	CL	A	602	1/1	1.00	0.05	-	17,17,17,17	0

6.5 Other polymers [i](#)

There are no such residues in this entry.