



## wwPDB EM Map/Model Validation Report ⓘ

Apr 10, 2016 – 01:41 PM BST

PDB ID : 3J2I  
EMDB ID: : EMD-5513  
Title : Structure of late pre-60S ribosomal subunits with nuclear export factor Arx1 bound at the peptide exit tunnel  
Authors : Bradatsch, B.; Leidig, C.; Granneman, S.; Gnaedig, M.; Tollervey, D.; Boettcher, B.; Beckmann, R.; Hurt, E.  
Deposited on : 2012-10-04  
Resolution : 11.90 Å(reported)

This is a wwPDB EM Map/Model Validation Report for a publicly released PDB/EMDB entry.  
For rigid body fitted models, validation errors reported here could stem from errors in the original structure(s) used in the fitting.

We welcome your comments at [validation@mail.wwpdb.org](mailto:validation@mail.wwpdb.org)

A user guide is available at

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

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MolProbity : 4.02b-467  
Mogul : unknown  
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)  
Validation Pipeline (wwPDB-VP) : trunk27241

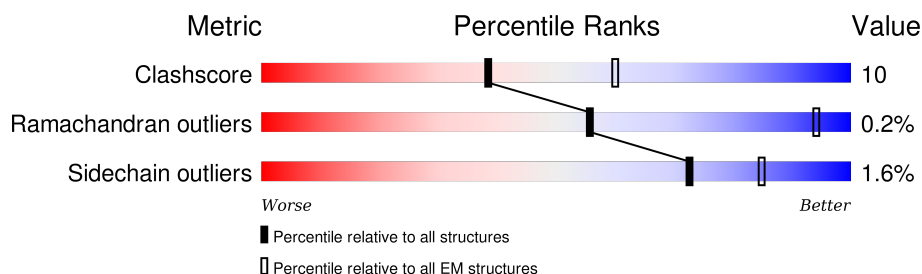
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*ELECTRON MICROSCOPY*

The reported resolution of this entry is 11.90 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)	EM structures (#Entries)
Clashscore	114402	924
Ramachandran outliers	111179	726
Sidechain outliers	111093	686

The table below summarises the geometric issues observed across the polymeric chains. The red, orange, yellow and green segments on the bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 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  $\leq 5\%$ .

Mol	Chain	Length	Quality of chain
1	B	245	
2	A	394	

## 2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 4529 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, 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 6.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	B	224	Total	C	N	O	S	0	0
			1692	1051	293	341	7		

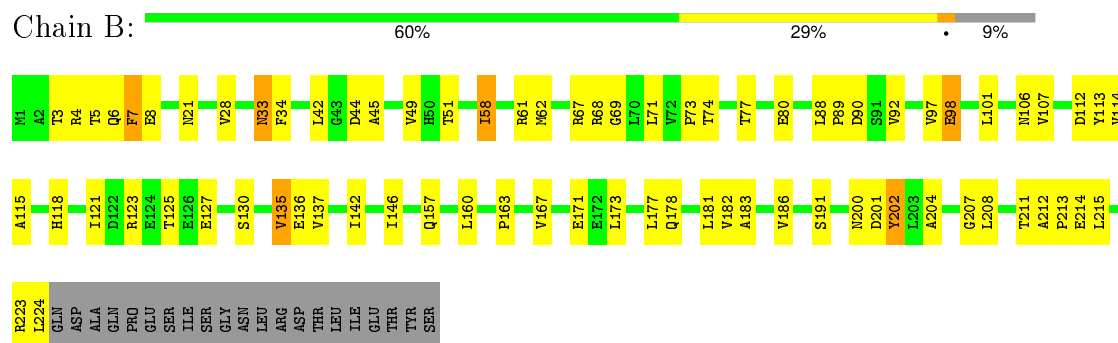
- Molecule 2 is a protein called Proliferation-associated protein 2G4.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	A	354	Total	C	N	O	S	17	0
			2837	1792	483	543	19		

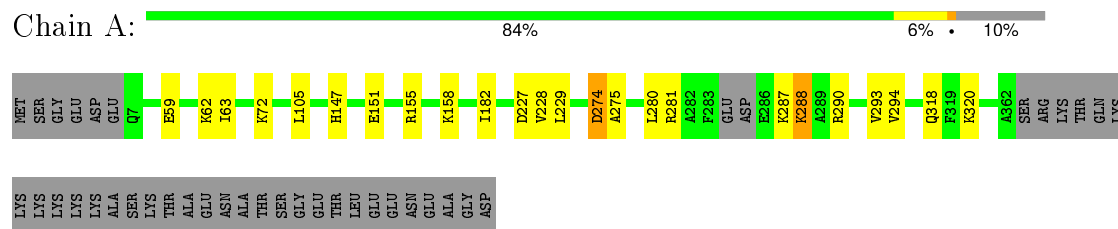
### 3 Residue-property plots

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. 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. 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 6



- Molecule 2: Proliferation-associated protein 2G4



## 4 Experimental information

Property	Value	Source
Reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	Depositor
Number of images	100000	Depositor
Resolution determination method	Not provided	Depositor
CTF correction method	subvolumes	Depositor
Microscope	FEI TITAN KRIOS	Depositor
Voltage (kV)	200	Depositor
Electron dose ( $e^-/\text{\AA}^2$ )	Not provided	Depositor
Minimum defocus (nm)	Not provided	Depositor
Maximum defocus (nm)	Not provided	Depositor
Magnification	Not provided	Depositor
Image detector	CCD	Depositor

## 5 Model quality [i](#)

### 5.1 Standard geometry [i](#)

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	$\# Z  > 2$	RMSZ	$\# Z  > 2$
1	B	0.45	1/1713 (0.1%)	0.70	1/2331 (0.0%)
2	A	0.56	0/2911	0.64	2/3913 (0.1%)
All	All	0.52	1/4624 (0.0%)	0.66	3/6244 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	202	TYR	CB-CG	-6.09	1.42	1.51

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	281	ARG	NE-CZ-NH1	6.46	123.53	120.30
2	A	281	ARG	NE-CZ-NH2	-6.02	117.29	120.30
1	B	7	PHE	N-CA-C	-5.04	97.40	111.00

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	B	1692	0	1689	73	0
2	A	2837	0	2868	14	0
All	All	4529	0	4557	87	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 (87) 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:B:5:THR:HA	1:B:208:LEU:HD11	1.46	0.96
2:A:147[B]:HIS:HD2	2:A:151:GLU:OE2	1.56	0.88
2:A:288[A]:LYS:H	2:A:288[A]:LYS:HE3	1.37	0.87
1:B:5:THR:CA	1:B:208:LEU:HD11	2.04	0.86
1:B:5:THR:CA	1:B:208:LEU:CD1	2.52	0.86
1:B:21:ASN:HD21	1:B:201:ASP:H	1.21	0.85
1:B:5:THR:HG22	1:B:6:GLN:H	1.42	0.84
1:B:118:HIS:HE2	1:B:146:ILE:HD12	1.42	0.83
1:B:107:VAL:HG13	1:B:118:HIS:HB2	1.61	0.80
1:B:5:THR:C	1:B:208:LEU:CD1	2.51	0.78
2:A:147[B]:HIS:CD2	2:A:151:GLU:OE2	2.39	0.74
1:B:49:VAL:HG21	1:B:88:LEU:HD12	1.70	0.74
1:B:89:PRO:O	1:B:92:VAL:HG12	1.88	0.74
1:B:212:ALA:HB3	1:B:213:PRO:HD3	1.70	0.73
1:B:5:THR:C	1:B:208:LEU:HD13	2.09	0.73
1:B:118:HIS:NE2	1:B:146:ILE:HD12	2.05	0.72
1:B:200:ASN:ND2	1:B:202:TYR:CZ	2.58	0.72
1:B:5:THR:N	1:B:208:LEU:HD12	2.05	0.71
1:B:6:GLN:HG3	1:B:208:LEU:HD13	1.76	0.68
1:B:113:TYR:C	1:B:135:VAL:HG22	2.14	0.67
1:B:167:VAL:O	1:B:171:GLU:HG3	1.97	0.65
2:A:182:ILE:HD12	2:A:229:LEU:HB3	1.79	0.65
1:B:178:GLN:HG2	1:B:178:GLN:O	1.96	0.64
1:B:5:THR:HG21	1:B:34:PHE:HA	1.80	0.63
1:B:177:LEU:HD12	1:B:181:LEU:HD11	1.83	0.61
1:B:68:ARG:NH2	1:B:112:ASP:O	2.33	0.61
1:B:5:THR:N	1:B:208:LEU:CD1	2.64	0.60
1:B:77:THR:OG1	1:B:80:GLU:HG3	2.02	0.60
2:A:59:GLU:HA	2:A:62:LYS:HE3	1.84	0.59
1:B:4:ARG:HB3	1:B:208:LEU:HA	1.83	0.59
1:B:67:ARG:HG2	1:B:112:ASP:OD1	2.01	0.59
1:B:163:PRO:HD2	1:B:191:SER:HB2	1.85	0.58
1:B:74:THR:OG1	1:B:98:GLU:HG2	2.04	0.58
1:B:61:ARG:HD3	1:B:106:ASN:HD21	1.68	0.58
1:B:5:THR:HG22	1:B:6:GLN:N	2.17	0.56
1:B:49:VAL:HG21	1:B:88:LEU:CD1	2.35	0.55
1:B:58:ILE:HG13	1:B:62:MET:HG2	1.89	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:71:LEU:HD22	1:B:97:VAL:CG2	2.38	0.53
1:B:113:TYR:O	1:B:135:VAL:HG22	2.09	0.52
1:B:71:LEU:HD22	1:B:97:VAL:HG21	1.90	0.52
1:B:5:THR:O	1:B:208:LEU:HD13	2.09	0.52
1:B:42:LEU:O	1:B:45:ALA:HB3	2.10	0.51
1:B:44:ASP:O	1:B:44:ASP:OD2	2.29	0.51
1:B:211:THR:OG1	1:B:214:GLU:HG3	2.12	0.50
1:B:28:VAL:HG13	1:B:51:THR:C	2.32	0.50
1:B:208:LEU:N	1:B:208:LEU:HD12	2.26	0.50
1:B:5:THR:O	1:B:208:LEU:CD1	2.60	0.50
2:A:280:LEU:HD21	2:A:293:VAL:HG21	1.94	0.49
1:B:107:VAL:HG13	1:B:118:HIS:CB	2.37	0.49
1:B:68:ARG:O	1:B:92:VAL:HA	2.13	0.49
1:B:90:ASP:OD2	1:B:90:ASP:O	2.31	0.48
1:B:142:ILE:N	1:B:142:ILE:HD12	2.29	0.48
2:A:105:LEU:HD23	2:A:105:LEU:C	2.35	0.47
1:B:7:PHE:O	1:B:8:GLU:HB2	2.15	0.47
1:B:33:ASN:ND2	1:B:33:ASN:O	2.46	0.47
1:B:118:HIS:HE2	1:B:146:ILE:CD1	2.23	0.46
2:A:280:LEU:HD21	2:A:293:VAL:CG2	2.46	0.46
1:B:160:LEU:HD12	1:B:182:VAL:O	2.16	0.46
1:B:135:VAL:HG13	1:B:136:GLU:N	2.30	0.46
1:B:58:ILE:HG13	1:B:58:ILE:O	2.17	0.45
1:B:62:MET:O	1:B:73:PRO:HD3	2.16	0.45
2:A:290:ARG:O	2:A:294:VAL:HG23	2.16	0.45
1:B:121:ILE:HD11	1:B:125:THR:HG22	1.99	0.45
1:B:204:ALA:HB2	1:B:224:LEU:HD11	1.99	0.44
1:B:208:LEU:N	1:B:208:LEU:CD1	2.81	0.44
2:A:227:ASP:CG	2:A:320[A]:LYS:HE2	2.39	0.43
1:B:163:PRO:HA	1:B:183:ALA:HB1	1.99	0.43
1:B:157:GLN:HG3	1:B:223:ARG:HG3	1.99	0.43
1:B:121:ILE:HD11	1:B:125:THR:CG2	2.48	0.43
1:B:21:ASN:ND2	1:B:201:ASP:H	2.02	0.43
2:A:274:ASP:HB3	2:A:275:ALA:H	1.62	0.43
1:B:173:LEU:CB	1:B:181:LEU:HD12	2.49	0.43
1:B:163:PRO:CD	1:B:191:SER:HB2	2.49	0.42
1:B:101:LEU:HD13	1:B:118:HIS:CG	2.54	0.42
1:B:173:LEU:HB3	1:B:181:LEU:HD12	2.02	0.42
2:A:155:ARG:O	2:A:158[B]:LYS:HD2	2.20	0.42
1:B:215:LEU:HA	1:B:215:LEU:HD12	1.81	0.41
1:B:114:VAL:HG12	1:B:115:ALA:N	2.35	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:69:GLY:HA2	1:B:92:VAL:HG23	2.03	0.41
1:B:130:SER:HB2	1:B:137:VAL:HG23	2.02	0.41
2:A:63:ILE:O	2:A:72:LYS:NZ	2.50	0.41
1:B:208:LEU:CD1	1:B:208:LEU:H	2.34	0.41
1:B:212:ALA:N	1:B:213:PRO:CD	2.83	0.41
1:B:33:ASN:C	1:B:33:ASN:HD22	2.20	0.41
2:A:228:VAL:O	2:A:318:GLN:HA	2.21	0.41
1:B:123:ARG:O	1:B:127:GLU:HG3	2.22	0.40
1:B:5:THR:O	1:B:207:GLY:HA2	2.21	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 PDB entries followed by that with respect to all EM entries.

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	B	222/245 (91%)	208 (94%)	13 (6%)	1 (0%)	34	77
2	A	363/394 (92%)	356 (98%)	7 (2%)	0	100	100
All	All	585/639 (92%)	564 (96%)	20 (3%)	1 (0%)	56	86

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	58	ILE

### 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 PDB entries followed by that with respect to all EM entries.

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	B	192/211 (91%)	187 (97%)	5 (3%)	54	80
2	A	316/336 (94%)	313 (99%)	3 (1%)	84	93
All	All	508/547 (93%)	500 (98%)	8 (2%)	72	88

All (8) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	3	THR
1	B	33	ASN
1	B	98	GLU
1	B	135	VAL
1	B	186	VAL
2	A	274	ASP
2	A	287	LYS
2	A	288[A]	LYS

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

Mol	Chain	Res	Type
1	B	6	GLN
1	B	11	ASN
1	B	21	ASN
1	B	50	HIS
1	B	79	GLN
1	B	82	GLN
1	B	83	HIS
1	B	86	ASN
1	B	106	ASN
1	B	168	GLN
1	B	178	GLN
2	A	242	GLN

### 5.3.3 RNA ⓘ

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 carbohydrates 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.