



wwPDB/EMDataBank EM Map/Model Validation Summary Report ⓘ

Aug 20, 2017 – 07:21 PM EDT

PDB ID : 4CRM
EMDB ID: : EMD-2598
Title : Cryo-EM of a pre-recycling complex with eRF1 and ABCE1
Authors : Preis, A.; Heuer, A.; Barrio-Garcia, C.; Hauser, A.; Eyler, D.; Berninghausen, O.; Green, R.; Becker, T.; Beckmann, R.
Deposited on : unknown
Resolution : 8.75 Å(reported)
Based on PDB ID : 1DT9

This is a wwPDB/EMDataBank EM Map/Model Validation Summary Report
for a publicly released PDB/EMDB entry.

We welcome your comments at validation@mail.wwpdb.org
A user guide is available at
<http://wwpdb.org/validation/2016/EMValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

MolProbity : 4.02b-467
Mogul : 1.7.2 (RC1), CSD as538be (2017)
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : rb-20029824

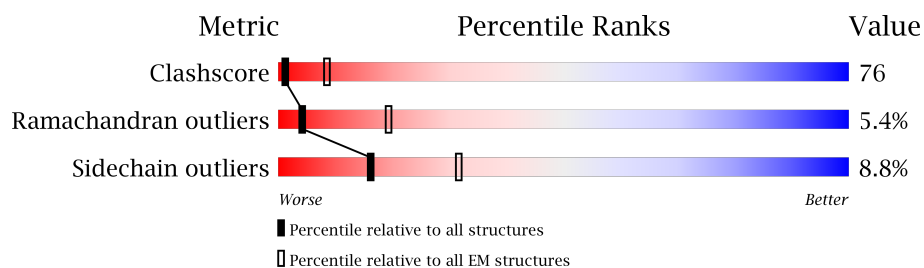
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 8.75 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	125131	1336
Ramachandran outliers	121729	1120
Sidechain outliers	121581	1026

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 ≥ 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	P	608	<div> <div style="width: 31%; background-color: green;"></div> <div style="width: 63%; background-color: yellow;"></div> <div style="width: 5%; background-color: orange;"></div> </div> <div>31% 63% 5%</div>
2	X	282	<div> <div style="width: 44%; background-color: green;"></div> <div style="width: 32%; background-color: yellow;"></div> <div style="width: 16%; background-color: orange;"></div> <div style="width: 9%; background-color: red;"></div> </div> <div>44% 32% 16% 9%</div>

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	ATP	P	1609	-	-	X	-
4	SF4	P	1610	-	-	X	-
4	SF4	P	1611	-	-	X	-
5	ADP	P	1612	-	-	X	-

2 Entry composition [i](#)

There are 7 unique types of molecules in this entry. The entry contains 7090 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 TRANSLATION INITIATION FACTOR RLI1.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	P	608	Total	C	N	O	S	0	0
			4804	3065	831	884	24		

- Molecule 2 is a protein called EUKARYOTIC PEPTIDE CHAIN RELEASE FACTOR SUB-UNIT 1.

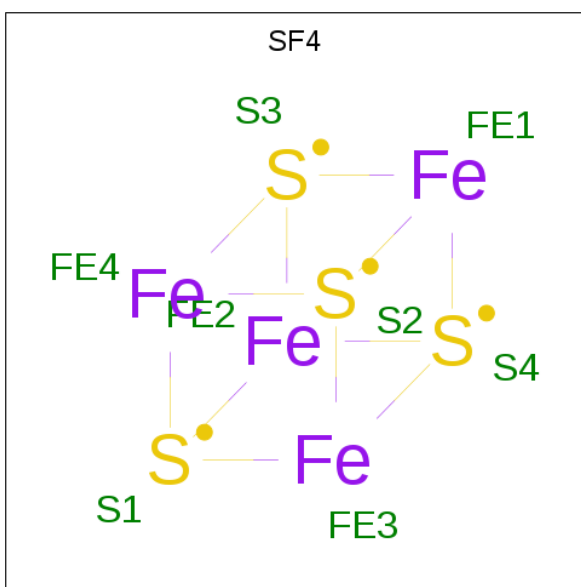
Mol	Chain	Residues	Atoms					AltConf	Trace
2	X	282	Total	C	N	O	S	0	0
			2210	1406	366	433	5		

- Molecule 3 is ADENOSINE-5'-TRIPHOSPHATE (three-letter code: ATP) (formula: $C_{10}H_{16}N_5O_{13}P_3$).



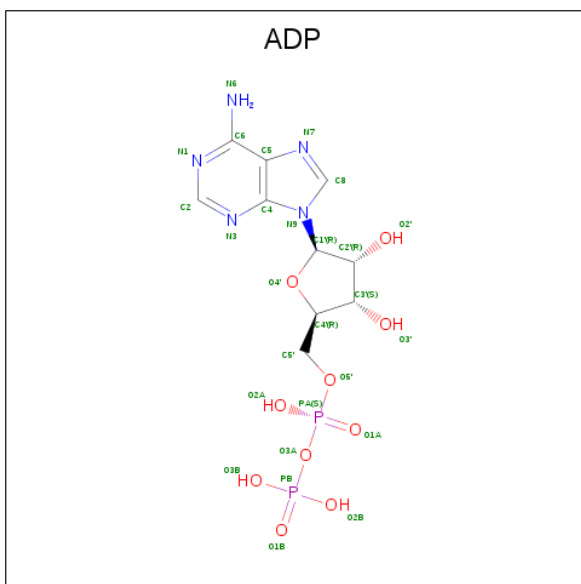
Mol	Chain	Residues	Atoms					AltConf
3	P	1	Total	C	N	O	P	0
			31	10	5	13	3	

- Molecule 4 is IRON/SULFUR CLUSTER (three-letter code: SF4) (formula: Fe_4S_4).



Mol	Chain	Residues	Atoms			AltConf
4	P	1	Total 16	Fe 8	S 8	0
4	P	1	Total 16	Fe 8	S 8	0

- Molecule 5 is ADENOSINE-5'-DIPHOSPHATE (three-letter code: ADP) (formula: $\text{C}_{10}\text{H}_{15}\text{N}_5\text{O}_{10}\text{P}_2$).



Mol	Chain	Residues	Atoms					AltConf
5	P	1	Total	C	N	O	P	0
			27	10	5	10	2	

- Molecule 6 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		AltConf
6	P	1	Total	Mg	0
			1	1	

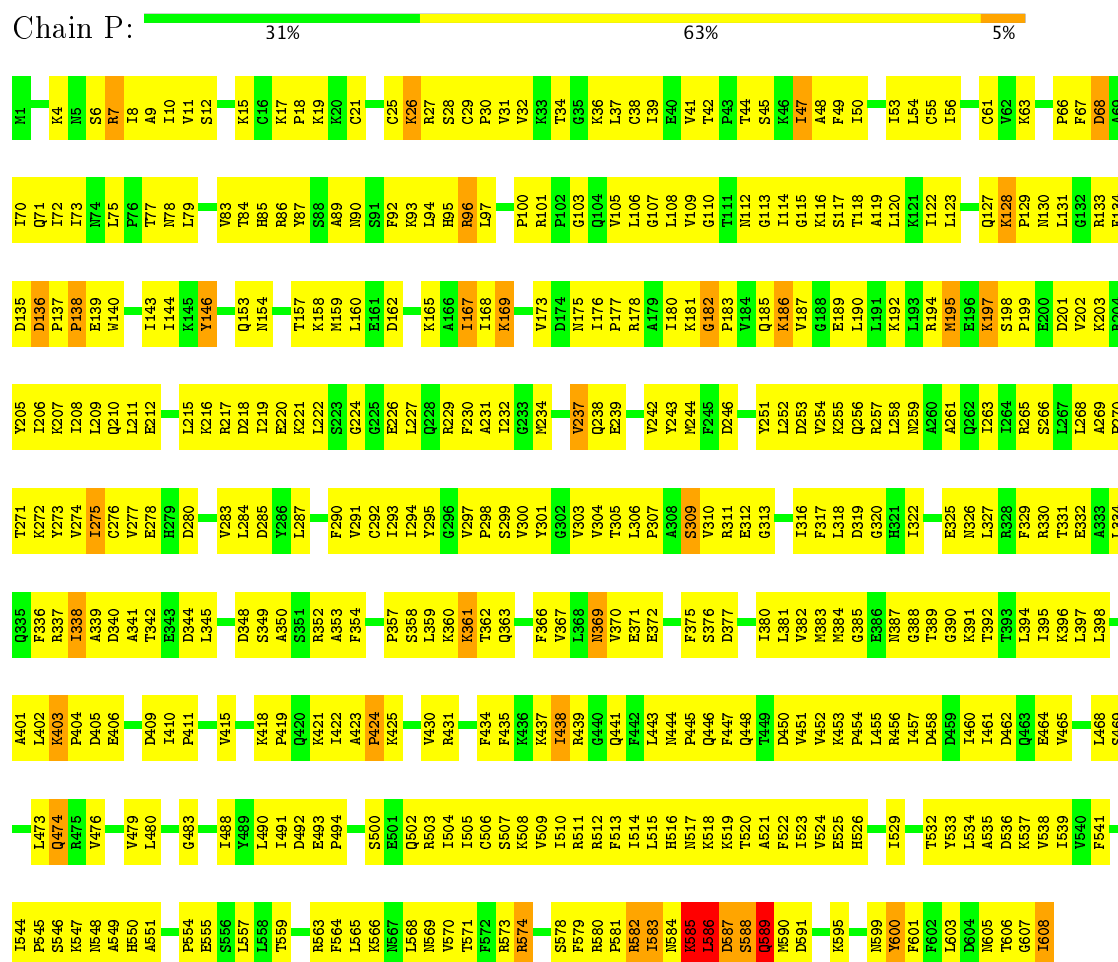
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		AltConf
7	P	1	Total	O	0
			1	1	

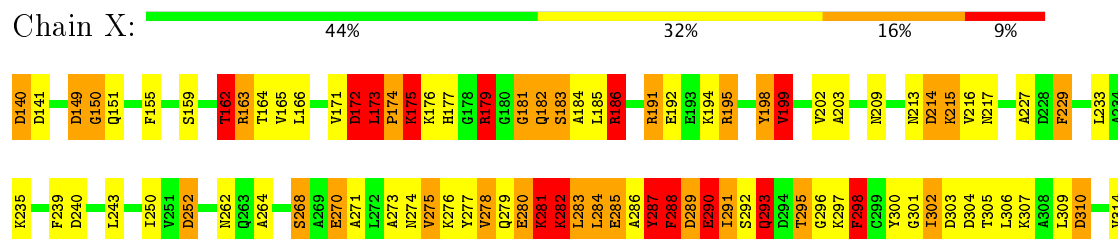
3 Residue-property plots

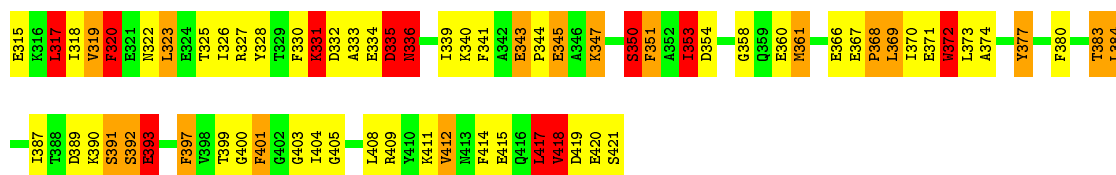
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry. Residues are color-coded according 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: TRANSLATION INITIATION FACTOR RLI1



• Molecule 2: EUKARYOTIC PEPTIDE CHAIN RELEASE FACTOR SUBUNIT 1





4 Experimental information

Property	Value	Source
Reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C1	Depositor
Number of particles used	39309	Depositor
Resolution determination method	Not provided	Depositor
CTF correction method	ON VOLUMES (SPIDER)	Depositor
Microscope	FEI MORGAGNI	Depositor
Voltage (kV)	200	Depositor
Electron dose ($e^-/\text{\AA}^2$)	20	Depositor
Minimum defocus (nm)	Not provided	Depositor
Maximum defocus (nm)	Not provided	Depositor
Magnification	147136	Depositor
Image detector	TVIPS TEMCAM-F416 (4k x 4k)	Depositor

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: ATP, MG, SF4, ADP

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	P	1.02	8/4893 (0.2%)	1.21	21/6603 (0.3%)
2	X	1.09	10/2246 (0.4%)	1.57	53/3021 (1.8%)
All	All	1.04	18/7139 (0.3%)	1.34	74/9624 (0.8%)

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	P	0	2
2	X	1	17
All	All	1	19

The worst 5 of 18 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	X	150	GLY	C-N	18.11	1.75	1.34
1	P	608	ILE	C-O	-12.06	1.00	1.23
2	X	421	SER	C-OXT	-12.06	1.00	1.23
2	X	421	SER	C-O	-12.06	1.00	1.23
1	P	608	ILE	C-OXT	-12.03	1.00	1.23

The worst 5 of 74 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	X	149	ASP	O-C-N	-17.67	93.16	123.20
2	X	149	ASP	CA-C-N	14.81	145.82	116.20
2	X	298	PHE	CB-CG-CD1	-13.00	111.70	120.80
1	P	589	GLN	CA-CB-CG	12.85	141.68	113.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	P	588	SER	N-CA-CB	12.81	129.71	110.50

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
2	X	331	LYS	CA

5 of 19 planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	P	574	ARG	Sidechain
1	P	589	GLN	Mainchain
2	X	140	ASP	Peptide
2	X	141	ASP	Sidechain,Peptide
2	X	149	ASP	Sidechain

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	P	4804	0	4945	827	0
2	X	2210	0	2182	251	0
3	P	31	0	12	29	0
4	P	16	0	0	15	0
5	P	27	0	12	20	0
6	P	1	0	0	1	0
7	P	1	0	0	0	0
All	All	7090	0	7151	1076	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 76.

The worst 5 of 1076 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:X:302:ILE:HD11	2:X:325:THR:CB	1.26	1.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:P:589:GLN:C	1:P:589:GLN:CA	1.77	1.50
2:X:328:TYR:CD1	2:X:366:GLU:HG2	1.45	1.49
2:X:150:GLY:C	2:X:151:GLN:N	1.75	1.37
2:X:323:LEU:CD1	2:X:370:ILE:CG1	1.94	1.37

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	P	606/608 (100%)	563 (93%)	35 (6%)	8 (1%)	14	56
2	X	280/282 (99%)	197 (70%)	43 (15%)	40 (14%)	0	5
All	All	886/890 (100%)	760 (86%)	78 (9%)	48 (5%)	4	25

5 of 48 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	P	68	ASP
1	P	138	PRO
1	P	197	LYS
2	X	162	THR
2	X	172	ASP

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	P	537/537 (100%)	516 (96%)	21 (4%)	37	66
2	X	234/234 (100%)	187 (80%)	47 (20%)	1	9
All	All	771/771 (100%)	703 (91%)	68 (9%)	16	39

5 of 68 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	X	186	ARG
2	X	268	SER
2	X	387	ILE
2	X	194	LYS
2	X	215	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 17 such sidechains are listed below:

Mol	Chain	Res	Type
1	P	347	ASN
1	P	369	ASN
1	P	562	ASN
1	P	259	ASN
1	P	599	ASN

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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 5 ligands modelled in this entry, 1 is monoatomic - leaving 4 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. 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 $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
3	ATP	P	1609	6	27,33,33	2.06	8 (29%)	25,52,52	3.36	9 (36%)
4	SF4	P	1610	1	0,12,12	0.00	-	0,24,24	0.00	-
4	SF4	P	1611	1	0,12,12	0.00	-	0,24,24	0.00	-
5	ADP	P	1612	-	25,29,29	1.11	2 (8%)	24,45,45	0.91	1 (4%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the chemical component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	ATP	P	1609	6	-	0/18/38/38	0/3/3/3
4	SF4	P	1610	1	-	0/0/48/48	0/6/5/5
4	SF4	P	1611	1	-	0/0/48/48	0/6/5/5
5	ADP	P	1612	-	-	0/12/32/32	0/3/3/3

The worst 5 of 10 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	P	1609	ATP	O5'-C5'	-4.45	1.27	1.44
3	P	1609	ATP	PA-O5'	-2.79	1.47	1.59
3	P	1609	ATP	C3'-C4'	-2.15	1.47	1.53
3	P	1609	ATP	C8-N7	-2.07	1.30	1.34
3	P	1609	ATP	PB-O1B	-2.01	1.43	1.50

The worst 5 of 10 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	P	1609	ATP	O5'-PA-O1A	-6.34	83.69	109.25
3	P	1609	ATP	C5'-C4'-C3'	-5.17	95.58	115.29
3	P	1609	ATP	N3-C2-N1	-2.94	126.29	128.86
3	P	1609	ATP	O2A-PA-O5'	-2.19	97.82	108.14
3	P	1609	ATP	C4'-O4'-C1'	-2.12	107.51	109.77

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

4 monomers are involved in 64 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	P	1609	ATP	29	0
4	P	1610	SF4	7	0
4	P	1611	SF4	8	0
5	P	1612	ADP	20	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
2	X	1

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	X	150:GLY	C	151:GLN	N	1.75