

Full wwPDB X-ray Structure Validation Report (i)

May 30, 2020 - 01:47 am BST

PDB ID : 6RK6

Title : Characterization of an intertidal zone metagenome oligoribonuclease and the

role of the intermolecular disulfide bond for homodimer formation and nuclease

activity.

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Deposited on : 2019-04-30

Resolution : 3.15 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 (i) symbol.

The following versions of software and data (see references (1)) were used in the production of this report:

MolProbity : 4.02b-467 Xtriage (Phenix) : 1.13

EDS : 2.11

buster-report : 1.1.7 (2018)

Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)

 $Refmac \quad : \quad 5.8.0158$

CCP4 : 7.0.044 (Gargrove)

Ideal geometry (proteins) : Engh & Huber (2001) Ideal geometry (DNA, RNA) : Parkinson et al. (1996)

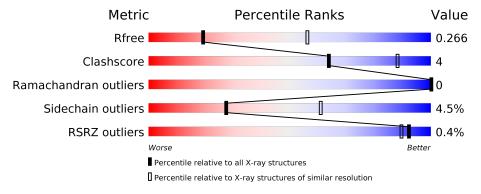
Validation Pipeline (wwPDB-VP) : 2.11

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure: X- $RAY\ DIFFRACTION$

The reported resolution of this entry is 3.15 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	$\begin{array}{c} \text{Whole archive} \\ (\#\text{Entries}) \end{array}$	$\begin{array}{c} {\rm Similar resolution} \\ (\#{\rm Entries, resolution range(\AA)}) \end{array}$
R_{free}	130704	1665 (3.20-3.12)
Clashscore	141614	1804 (3.20-3.12)
Ramachandran outliers	138981	1770 (3.20-3.12)
Sidechain outliers	138945	1769 (3.20-3.12)
RSRZ outliers	127900	1616 (3.20-3.12)

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 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 <=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	184	83%	13%	•	-
1	В	184	81%	15%	•	-
1	С	184	85%	10%	•	-



2 Entry composition (i)

There are 2 unique types of molecules in this entry. The entry contains 4379 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 Oligoribonuclease.

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace		
1	Λ	177	Total	С	N	О	S	0	0	0
1	Λ	111	1456	917	259	273	7	0		
1	D	177	Total	С	N	О	S	0	0	0
1		177	1456	917	259	273	7	0	U	
1	С	178	Total	С	N	О	S	0	0	0
	110	1464	921	260	276	7	0	0	U	

• Molecule 2 is MANGANESE (II) ION (three-letter code: MN) (formula: Mn) (labeled as "Ligand of Interest" by author).

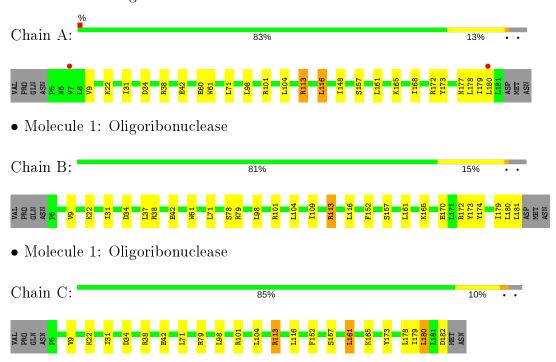
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	В	1	Total Mn 1 1	0	0
2	A	1	Total Mn 1 1	0	0
2	С	1	Total Mn 1 1	0	0



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 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: Oligoribonuclease





4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants	108.32Å 108.32Å 101.33Å	Depositor
a, b, c, α , β , γ	90.00° 90.00° 120.00°	
Resolution (Å)	44.58 - 3.15	Depositor
Resolution (A)	44.58 - 3.15	EDS
% Data completeness	99.8 (44.58-3.15)	Depositor
(in resolution range)	91.8 (44.58-3.15)	EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) > 1$	0.91 (at 3.12Å)	Xtriage
Refinement program	PHENIX (1.12_2829: ???)	Depositor
D D	0.245 , 0.266	Depositor
R, R_{free}	0.245 , 0.266	DCC
R_{free} test set	926 reflections (7.56%)	wwPDB-VP
Wilson B-factor (Å ²)	69.7	Xtriage
Anisotropy	0.333	Xtriage
Bulk solvent $k_{sol}(e/Å^3)$, $B_{sol}(Å^2)$	0.30,21.5	EDS
L-test for twinning ²	$< L >=0.49, < L^2>=0.32$	Xtriage
Estimated twinning fraction	0.037 for -h,-k,l	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	4379	wwPDB-VP
Average B, all atoms (Å ²)	82.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 11.68% of the height of the origin peak. No significant pseudotranslation is detected.

²Theoretical values of <|L|>, $< L^2>$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.



 $^{^{1}}$ Intensities estimated from amplitudes.

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: MN

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		
MIOI	Chain	RMSZ	# Z >5	RMSZ	# Z > 5	
1	A	0.26	0/1484	0.50	0/2006	
1	В	0.26	0/1484	0.49	0/2006	
1	С	0.26	0/1492	0.46	0/2017	
All	All	0.26	0/4460	0.48	0/6029	

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	С	0	1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	С	180	LEU	Peptide

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	1456	0	1454	16	0
1	В	1456	0	1455	16	0
1	С	1464	0	1458	10	0
2	A	1	0	0	0	0
2	В	1	0	0	0	0
2	С	1	0	0	0	0
All	All	4379	0	4367	37	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 4.

All (37) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic	Clash
Atom-1	Atom-2	${f distance}\;({f \AA})$	$overlap(\AA)$
1:B:180:LEU:HG	1:B:181:LEU:HG	1.51	0.92
1:B:104:LEU:HB2	1:B:113:ARG:HH21	1.54	0.73
1:C:104:LEU:HB2	1:C:113:ARG:HH21	1.54	0.72
1:A:104:LEU:HB2	1:A:113:ARG:HH21	1.54	0.70
1:A:177:ASN:ND2	1:B:79:ARG:H	1.89	0.69
1:A:177:ASN:HD22	1:B:79:ARG:H	1.44	0.64
1:C:179:ILE:HG22	1:C:180:LEU:H	1.66	0.61
1:A:179:ILE:HG22	1:A:180:LEU:H	1.66	0.60
1:B:179:ILE:HG22	1:B:180:LEU:H	1.66	0.60
1:A:42:GLU:OE2	1:A:165:LYS:NZ	2.42	0.53
1:B:42:GLU:OE2	1:B:165:LYS:NZ	2.42	0.53
1:C:42:GLU:OE2	1:C:165:LYS:NZ	2.42	0.52
1:C:34:ASP:OD2	1:C:38:ARG:NH1	2.42	0.52
1:A:34:ASP:OD2	1:A:38:ARG:NH1	2.42	0.51
1:B:37:LEU:O	1:B:172:ARG:NH2	2.43	0.51
1:B:34:ASP:OD2	1:B:38:ARG:NH1	2.42	0.50
1:A:179:ILE:HG23	1:C:178:LEU:HG	1.94	0.49
1:A:61:TRP:CD1	1:A:61:TRP:N	2.84	0.46
1:A:31:ILE:HD11	1:A:161:LEU:HG	1.98	0.46
1:C:31:ILE:HD11	1:C:161:LEU:HG	1.98	0.46
1:B:31:ILE:HD11	1:B:161:LEU:HG	1.98	0.44
1:A:148:ILE:O	1:A:173:TYR:OH	2.25	0.44
1:A:177:ASN:HD22	1:B:79:ARG:N	2.14	0.44
1:B:180:LEU:O	1:B:181:LEU:HB2	2.19	0.43
1:A:168:ILE:O	1:A:172:ARG:HG2	2.18	0.42
1:B:78:SER:O	1:B:79:ARG:HD3	2.20	0.42
1:A:98:LEU:HB2	1:A:101:ARG:HD2	2.02	0.42
1:A:178:LEU:HG	1:C:179:ILE:HG23	2.01	0.42

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Atom-1	Atom-2	Interatomic	Clash
Atom-1	Atom-2	$\operatorname{distance}\left(ext{\AA} ight)$	overlap(A)
1:C:98:LEU:HB2	1:C:101:ARG:HD2	2.02	0.41
1:B:98:LEU:HB2	1:B:101:ARG:HD2	2.02	0.41
1:A:116:LEU:HD13	1:A:116:LEU:HA	1.90	0.41
1:C:179:ILE:HG22	1:C:180:LEU:N	2.35	0.41
1:A:60:GLU:HB2	1:A:61:TRP:HD1	1.86	0.41
1:C:152:PHE:HB2	1:C:173:TYR:CD2	2.56	0.41
1:B:109:ILE:HD13	1:B:109:ILE:HA	1.97	0.41
1:B:152:PHE:HB2	1:B:173:TYR:CD2	2.56	0.40
1:B:170:GLU:HG2	1:B:174:TYR:CE2	2.57	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	ed Favoured Allowed		Outliers	Perce	$_{ m ntiles}$
1	A	175/184~(95%)	172 (98%)	3 (2%)	0	100	100
1	В	175/184 (95%)	173 (99%)	2 (1%)	0	100	100
1	С	176/184 (96%)	172 (98%)	4 (2%)	0	100	100
All	All	526/552 (95%)	517 (98%)	9 (2%)	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	Perce	entiles
1	A	$162/169 \; (96\%)$	156 (96%)	6 (4%)	34	66
1	В	162/169 (96%)	155 (96%)	7 (4%)	29	62
1	С	163/169 (96%)	154 (94%)	9 (6%)	21	54
All	All	487/507 (96%)	465 (96%)	22 (4%)	27	61

All (22) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	9	VAL
1	A	22	LYS
1	A	71	LEU
1	A	113	ARG
1	A	116	LEU
1	A	157	SER
1	В	9	VAL
1	В	22	LYS
1	В	61	TRP
1	В	71	LEU
1	В	113	ARG
1	В	116	LEU
1	В	157	SER
1	С	9	VAL
1	С	22	LYS
1	С	71	LEU
1	С	79	ARG
1	С	113	ARG
1	С	116	LEU
1	C C C C C C	157	SER
1	С	161	LEU
1	С	182	ASP

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

Mol	Chain	${f Res}$	\mathbf{Type}
1	A	177	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 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 (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 (i)

6.1 Protein, DNA and RNA chains (i)

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, 95^{th} 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 $>$	$\# \mathrm{RSRZ}{>}2$	$\mathbf{OWAB}(\mathrm{\AA}^2)$	Q<0.9
1	A	177/184 (96%)	-0.01	2 (1%) 80 70	62, 84, 99, 106	0
1	В	177/184 (96%)	0.01	0 100 100	63, 83, 99, 110	1 (0%)
1	С	178/184 (96%)	-0.05	0 100 100	60, 79, 99, 106	0
All	All	532/552 (96%)	-0.01	2 (0%) 92 89	60, 82, 99, 110	1 (0%)

All (2) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	7	VAL	2.2
1	A	180	LEU	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. The B-factors column lists the minimum, median, 95^{th} 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	$ extbf{B-factors}(extbf{A}^2)$	Q<0.9
2	MN	A	201	1/1	0.81	0.08	99,99,99,99	0

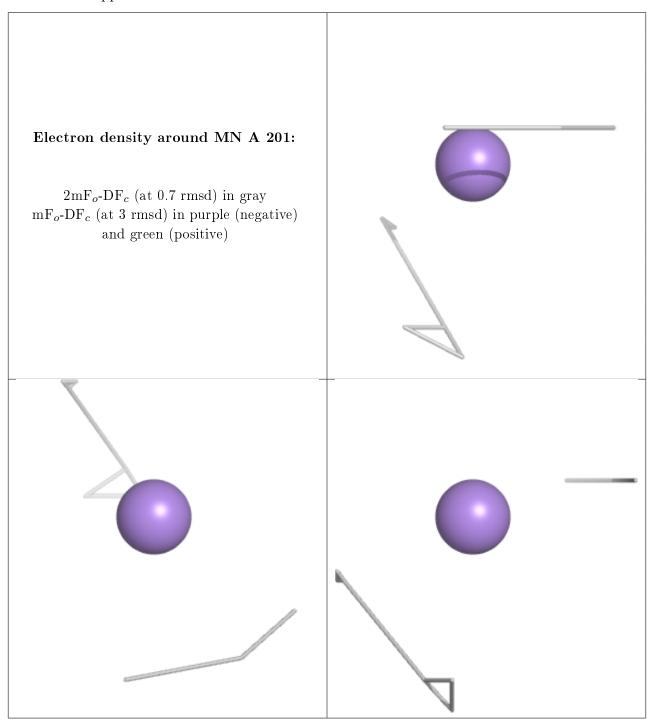
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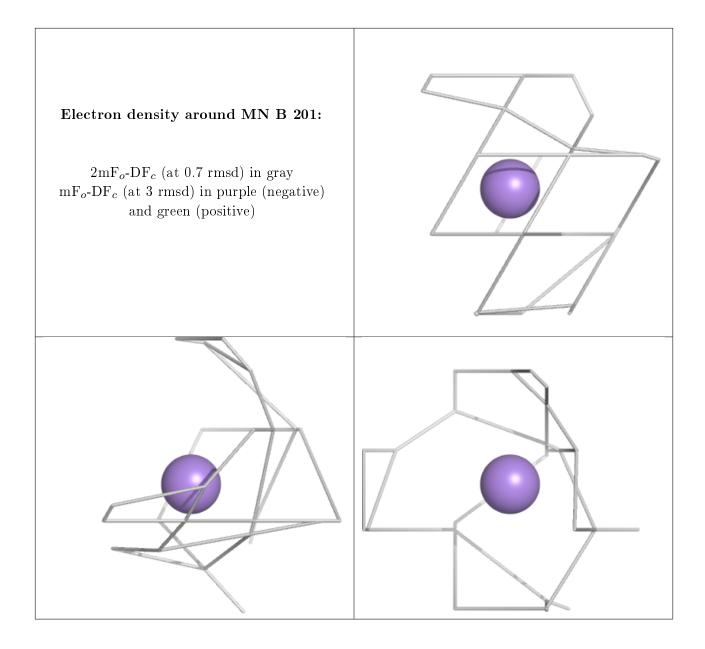
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	${f B\text{-factors}}({f \AA}^2)$	Q < 0.9
2	MN	В	201	1/1	0.81	0.22	115,115,115,115	0
2	MN	С	201	1/1	0.81	0.16	95,95,95,95	0

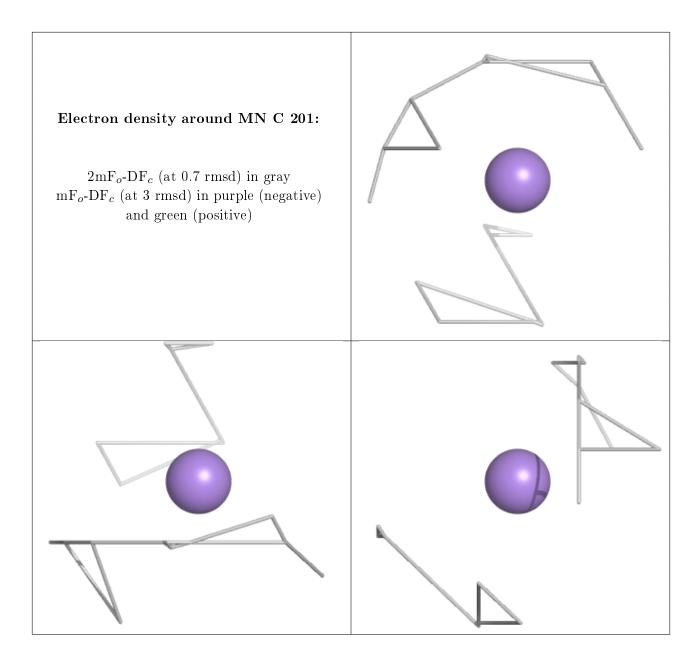
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.











6.5 Other polymers (i)

There are no such residues in this entry.

