Full wwPDB X-ray Structure Validation Report (i)

Dec 4, 2022 – 03:14 pm GMT

PDB ID	:	6GYV
Title	:	Lariat-capping ribozyme (circular permutation form)
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Deposited on	:	2018-07-02
Resolution	:	2.50 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 types of validation reports are described at
http://www.wwpdb.org/validation/2017/FAQs#types.

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

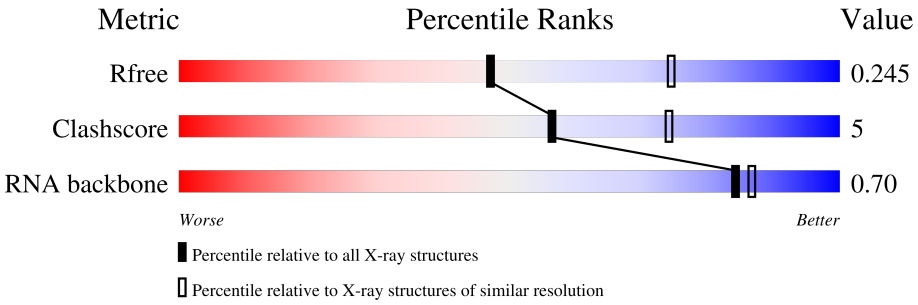
MolProbity	:	4.02b-467
Mogul	:	1.8.4, CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.31.3
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0267
CCP4	:	7.1.010 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.31.3

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 2.50 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	4661 (2.50-2.50)
Clashscore	141614	5346 (2.50-2.50)
RNA backbone	3102	1008 (2.84-2.16)

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 <=5%

Mol	Chain	Length	Quality of chain
1	A	192	72% 21% 5% •

2 Entry composition

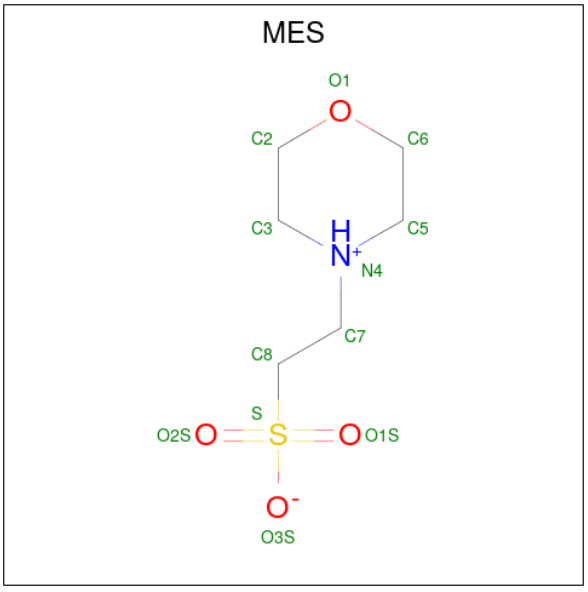
There are 5 unique types of molecules in this entry. The entry contains 4210 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 RNA chain called Lariat-capping ribozyme.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	P			
1	A	190	4096	1831	740	1334	191	0	2	0

- Molecule 2 is 2-(N-MORPHOLINO)-ETHANESULFONIC ACID (three-letter code: MES) (formula: $C_6H_{13}NO_4S$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	N	O	S		
2	A	1	12	6	1	4	1	0	0

- Molecule 3 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	Mg		
3	A	2	2	2	0	0

- Molecule 4 is SODIUM ION (three-letter code: NA) (formula: Na).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total Na 1 1	0	0

- Molecule 5 is water.

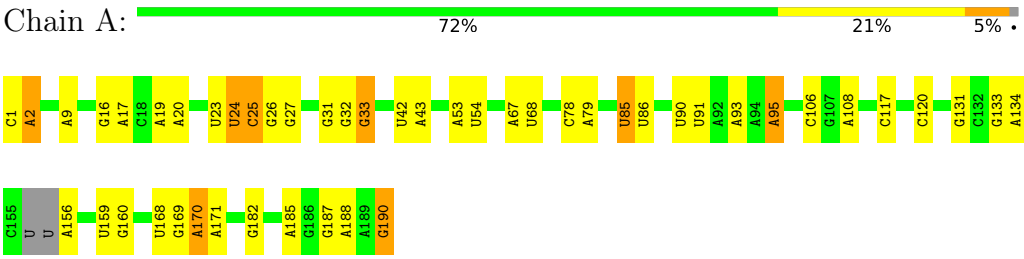
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	99	Total O 99 99	0	0

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. 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: Lariat-capping ribozyme

Chain A:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α, β, γ	57.63Å 85.71Å 108.53Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	47.82 – 2.50 47.82 – 2.50	Depositor EDS
% Data completeness (in resolution range)	96.5 (47.82-2.50) 96.0 (47.82-2.50)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.61 (at 2.51Å)	Xtriage
Refinement program	PHENIX 1.11rc1_2513	Depositor
R, R_{free}	0.198 , 0.245 0.198 , 0.245	Depositor DCC
R_{free} test set	1044 reflections (5.66%)	wwPDB-VP
Wilson B-factor (Å^2)	46.6	Xtriage
Anisotropy	0.183	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	(Not available) , (Not available)	EDS
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.94	EDS
Total number of atoms	4210	wwPDB-VP
Average B, all atoms (Å^2)	45.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.49% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical 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.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: MES, MG, NA

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	0.19	0/4581	0.75	3/7137 (0.0%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	85	U	C2-N1-C1'	6.82	125.88	117.70
1	A	85	U	N1-C2-O2	5.94	126.95	122.80
1	A	85	U	N3-C2-O2	-5.71	118.21	122.20

There are no chirality outliers.

There are no planarity outliers.

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	A	4096	0	2073	31	0
2	A	12	0	12	1	0
3	A	2	0	0	0	0
4	A	1	0	0	0	0
5	A	99	0	0	0	0
All	All	4210	0	2085	31	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 5.

All (31) 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:154:U:H3	1:A:156:A:H61	1.43	0.67
1:A:95:A:OP2	1:A:131:G:O2'	2.15	0.64
1:A:16:G:N2	1:A:33:G:H22	1.98	0.61
1:A:16:G:H1	1:A:33:G:H1	1.52	0.58
1:A:24:U:H4'	1:A:25:C:OP2	2.04	0.57
1:A:16:G:H2'	1:A:17:A:C8	2.41	0.55
1:A:182:G:N2	1:A:185:A:OP2	2.32	0.53
1:A:42:U:H2'	1:A:43:A:C8	2.45	0.51
1:A:16:G:H22	1:A:33:G:N2	2.10	0.50
1:A:1:C:H3'	1:A:2:A:H5"	1.94	0.49
1:A:16:G:H2'	1:A:17:A:H8	1.77	0.48
1:A:53:A:H2'	1:A:54:U:C6	2.49	0.47
1:A:32:G:H2'	1:A:33:G:O4'	2.15	0.47
1:A:159:U:H5"	1:A:160:G:H5'	1.96	0.47
1:A:16:G:N2	1:A:33:G:N2	2.65	0.45
1:A:16:G:H22	1:A:33:G:H22	1.62	0.45
1:A:19:A:H2'	1:A:20:A:C8	2.52	0.45
1:A:170:A:O2'	1:A:171[B]:A:OP2	2.28	0.45
1:A:31:G:H2'	1:A:32:G:C8	2.53	0.44
1:A:90:U:H2'	1:A:91:U:C6	2.52	0.44
1:A:78:C:H2'	1:A:79:A:C8	2.53	0.44
1:A:187:G:H2'	1:A:188:A:H8	1.82	0.43
1:A:136:A:H2'	1:A:137:G:H5"	2.00	0.43
1:A:16:G:H22	1:A:33:G:H1	1.66	0.43
1:A:9:A:N1	1:A:170:A:O2'	2.52	0.43
1:A:168:U:OP2	2:A:201:MES:H62	2.20	0.42
1:A:187:G:H2'	1:A:188:A:C8	2.56	0.41
1:A:17:A:C2	1:A:33:G:N2	2.89	0.41
1:A:106:C:O2'	1:A:108:A:OP2	2.24	0.41
1:A:171[B]:A:H8	1:A:190:G:N7	2.18	0.40
1:A:133:G:H4'	1:A:134:A:H5"	2.02	0.40

There are no symmetry-related clashes.

5.3 Torsion angles (i)

5.3.1 Protein backbone (i)

There are no protein molecules in this entry.

5.3.2 Protein sidechains (i)

There are no protein molecules in this entry.

5.3.3 RNA (i)

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
1	A	186/192 (96%)	18 (9%)	3 (1%)

All (18) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
1	A	2	A
1	A	24	U
1	A	25	C
1	A	26	G
1	A	27	G
1	A	33	G
1	A	67	A
1	A	68	U
1	A	86	U
1	A	93	A
1	A	95	A
1	A	117	C
1	A	120	C
1	A	136	A
1	A	137	G
1	A	169	G
1	A	170	A
1	A	190	G

All (3) RNA pucker outliers are listed below:

Mol	Chain	Res	Type
1	A	23	U
1	A	24	U
1	A	85	U

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)

Of 4 ligands modelled in this entry, 3 are monoatomic - leaving 1 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	$\#\|Z\| > 2$
2	MES	A	201	4	12,12,12	2.19	1 (8%)	14,16,16	2.32	8 (57%)

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
2	MES	A	201	4	-	5/6/14/14	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	201	MES	C8-S	-7.31	1.67	1.77

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	A	201	MES	C5-N4-C3	4.53	119.04	108.83
2	A	201	MES	C7-N4-C3	3.24	119.52	111.23
2	A	201	MES	C2-C3-N4	-3.00	105.56	110.10

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	A	201	MES	O1S-S-C8	2.83	110.32	106.92
2	A	201	MES	C7-N4-C5	2.73	118.20	111.23
2	A	201	MES	O2S-S-C8	2.39	109.79	106.92
2	A	201	MES	O3S-S-C8	2.06	109.10	105.77
2	A	201	MES	C6-C5-N4	-2.03	107.03	110.10

There are no chirality outliers.

All (5) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	201	MES	C8-C7-N4-C3
2	A	201	MES	C8-C7-N4-C5
2	A	201	MES	C7-C8-S-O3S
2	A	201	MES	C7-C8-S-O1S
2	A	201	MES	C7-C8-S-O2S

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	201	MES	1	0

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

Unable to reproduce the depositors R factor - this section is therefore empty.

6.2 Non-standard residues in protein, DNA, RNA chains

Unable to reproduce the depositors R factor - this section is therefore empty.

6.3 Carbohydrates

Unable to reproduce the depositors R factor - this section is therefore empty.

6.4 Ligands

Unable to reproduce the depositors R factor - this section is therefore empty.

6.5 Other polymers

Unable to reproduce the depositors R factor - this section is therefore empty.