



wwPDB X-ray Structure Validation Summary Report ⓘ

Feb 1, 2016 – 07:58 PM GMT

PDB ID : 4QJH
Title : Crystal Structure of the Twister Ribozyme with the Nucleotide 5'- to the
Cleavage Site Ordered at 4.1 Å Resolution
Authors : Eiler, D.R.; Wang, J.; Steitz, T.A.
Deposited on : 2014-06-03
Resolution : 3.88 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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 : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
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) : trunk26865

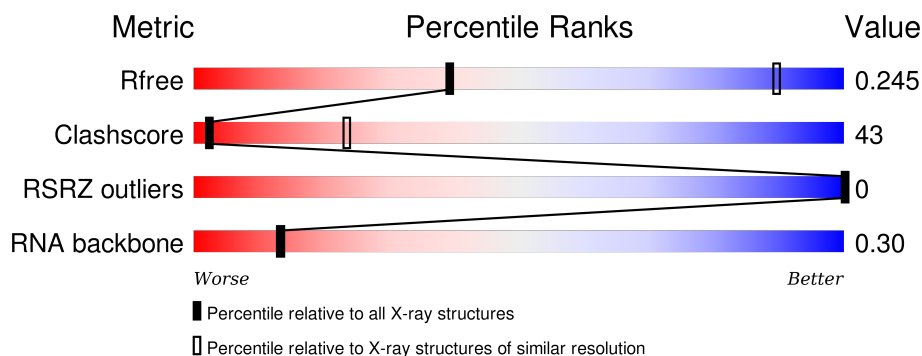
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.88 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	1005 (4.24-3.52)
Clashscore	102246	1026 (4.20-3.56)
RSRZ outliers	91569	1009 (4.24-3.52)
RNA backbone	2183	1071 (4.84-2.80)

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	26	<div> <div>12%</div> <div>62%</div> <div>27%</div> </div>
1	C	26	<div> <div>15%</div> <div>54%</div> <div>19%</div> <div>12%</div> </div>
1	E	26	<div> <div>•</div> <div>69%</div> <div>27%</div> </div>
2	B	48	<div> <div>8%</div> <div>46%</div> <div>46%</div> </div>
2	D	48	<div> <div>8%</div> <div>46%</div> <div>42%</div> <div>•</div> </div>
2	F	48	<div> <div>8%</div> <div>52%</div> <div>40%</div> </div>

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 4774 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 Twister Ribozyme.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	26	Total	C	N	O	P	0	0	0
			544	245	92	182	25			
1	C	26	Total	C	N	O	P	0	0	0
			544	245	92	182	25			
1	E	26	Total	C	N	O	P	0	0	0
			544	245	92	182	25			

- Molecule 2 is a RNA chain called Twister Ribozyme.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	48	Total	C	N	O	P	0	0	0
			1044	463	199	332	50			
2	D	48	Total	C	N	O	P	0	0	0
			1044	463	199	332	50			
2	F	48	Total	C	N	O	P	0	0	0
			1044	463	199	332	50			

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

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	1	Total	Mg	0	0
			1	1		
3	C	3	Total	Mg	0	0
			3	3		


- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	C	5	Total	O	0	0
			5	5		
4	E	1	Total	O	0	0
			1	1		

3 Residue-property plots

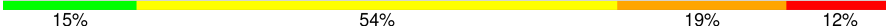
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according 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 ($RSRZ > 2$). 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: Twister Ribozyme

Chain A: 



• Molecule 1: Twister Ribozyme

Chain C: 



• Molecule 1: Twister Ribozyme

Chain E: 



• Molecule 2: Twister Ribozyme

Chain B: 



• Molecule 2: Twister Ribozyme

Chain D: 



• Molecule 2: Twister Ribozyme

Chain F: 

G27	G28	C29	A30	A31	U32	A33	A34	A35	G36	C37	G38	G39	U40	U41	A42	C43	A44	A45	G46	C47	C48	C49	G50	C51	A52	A53	A54	A55	A56	U57	A58	G59	C60	A61	G62	A63	G64	U65	A66	A67	U68	G69	U70	C71	G72	C73	G74
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4 Data and refinement statistics

Property	Value	Source
Space group	P 61 2 2	Depositor
Cell constants a, b, c, α , β , γ	228.77Å 228.77Å 101.06Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	198.12 – 3.88 49.77 – 4.16	Depositor EDS
% Data completeness (in resolution range)	99.4 (198.12-3.88) 100.0 (49.77-4.16)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.00 (at 4.14Å)	Xtriage
Refinement program	REFMAC 5.8.0073	Depositor
R, R_{free}	0.185 , 0.232 0.192 , 0.245	Depositor DCC
R_{free} test set	600 reflections (5.19%)	DCC
Wilson B-factor (Å ²)	270.3	Xtriage
Anisotropy	0.168	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.28 , 146.5	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 12154 reflections	Xtriage
F_o, F_c correlation	0.98	EDS
Total number of atoms	4774	wwPDB-VP
Average B, all atoms (Å ²)	147.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.83% 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, 0.375 respectively for untwinned datasets, and 0.333, 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: GTP, MG

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	0.31	0/606	0.72	0/941
1	C	0.50	1/606 (0.2%)	0.96	4/941 (0.4%)
1	E	0.30	0/606	0.74	0/941
2	B	0.33	0/1135	0.69	0/1768
2	D	0.44	0/1135	0.82	3/1768 (0.2%)
2	F	0.32	0/1135	0.74	0/1768
All	All	0.37	1/5223 (0.0%)	0.77	7/8127 (0.1%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	10	A	O3'-P	-6.13	1.53	1.61

The worst 5 of 7 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	10	A	O5'-P-OP2	-6.89	99.50	105.70
1	C	15	A	C2'-C3'-O3'	6.54	124.16	113.70
2	D	35	A	C2'-C3'-O3'	6.38	123.91	113.70
1	C	10	A	C2'-C3'-O3'	-5.86	96.61	109.50
2	D	53	A	C4'-C3'-O3'	5.30	123.60	113.00

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	A	544	0	280	23	0
1	C	544	0	280	39	0
1	E	544	0	280	31	0
2	B	1044	0	522	79	1
2	D	1044	0	522	90	0
2	F	1044	0	522	67	0
3	B	1	0	0	0	0
3	C	3	0	0	0	0
4	C	5	0	0	0	0
4	E	1	0	0	0	0
All	All	4774	0	2406	300	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 43.

The worst 5 of 300 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:D:40:U:H3'	2:D:41:U:C5'	1.67	1.22
2:D:40:U:H3'	2:D:41:U:H5'	1.21	1.10
2:F:40:U:H3'	2:F:41:U:H5'	1.12	1.09
1:C:6:C:H2'	1:C:7:A:H8	1.21	1.06
2:B:51:C:H3'	2:B:52:A:H5''	1.37	1.04

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)
2:B:53:A:OP1	2:B:53:A:OP1[7_556]	2.12	0.08

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	24/26 (92%)	13 (54%)	2 (8%)
1	C	24/26 (92%)	9 (37%)	4 (16%)
1	E	24/26 (92%)	11 (45%)	2 (8%)
2	B	47/48 (97%)	23 (48%)	7 (14%)
2	D	47/48 (97%)	22 (46%)	5 (10%)
2	F	46/48 (95%)	26 (56%)	7 (15%)
All	All	212/222 (95%)	104 (49%)	27 (12%)

5 of 104 RNA backbone outliers are listed below:

Mol	Chain	Res	Type
1	A	4	G
1	A	10	A
1	A	11	A
1	A	13	G
1	A	14	C

5 of 27 RNA pucker outliers are listed below:

Mol	Chain	Res	Type
1	C	16	G
2	D	45	A
2	F	57	U
2	D	27	GTP
2	B	35	A

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

3 non-standard protein/DNA/RNA residues are modelled in this entry.

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	GTP	B	27	1,2	25,34,34	1.23	2 (8%)	34,54,54	2.11	10 (29%)
2	GTP	D	27	1,2	25,34,34	1.28	3 (12%)	34,54,54	2.33	11 (32%)
2	GTP	F	27	1,2	25,34,34	1.28	3 (12%)	34,54,54	1.80	7 (20%)

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	GTP	B	27	1,2	-	0/18/38/38	0/3/3/3
2	GTP	D	27	1,2	-	0/18/38/38	0/3/3/3
2	GTP	F	27	1,2	-	0/18/38/38	0/3/3/3

The worst 5 of 8 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	27	GTP	O4'-C1'	2.59	1.44	1.41
2	F	27	GTP	O4'-C1'	2.70	1.44	1.41
2	B	27	GTP	C5-C4	3.10	1.47	1.40
2	D	27	GTP	C5-C4	3.33	1.48	1.40
2	F	27	GTP	C5-C4	3.49	1.48	1.40

The worst 5 of 28 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	27	GTP	C2'-C1'-N9	-6.22	104.80	114.29
2	D	27	GTP	C5-C6-N1	-4.50	117.44	123.59
2	F	27	GTP	PA-O3A-PB	-4.40	120.36	132.73
2	B	27	GTP	C5-C6-N1	-4.35	117.64	123.59
2	B	27	GTP	C1'-N9-C4	-4.15	120.67	126.94

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

3 monomers are involved in 10 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	27	GTP	2	0
2	D	27	GTP	3	0
2	F	27	GTP	5	0

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 4 ligands modelled in this entry, 4 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, 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	26/26 (100%)	-0.87	0 100 100	77, 171, 249, 269	0
1	C	26/26 (100%)	-0.56	0 100 100	49, 89, 163, 189	0
1	E	26/26 (100%)	-0.92	0 100 100	96, 185, 228, 236	0
2	B	47/48 (97%)	-0.85	0 100 100	81, 156, 240, 281	0
2	D	47/48 (97%)	-0.57	0 100 100	55, 88, 161, 194	0
2	F	47/48 (97%)	-1.01	0 100 100	98, 158, 227, 267	0
All	All	219/222 (98%)	-0.80	0 100 100	49, 138, 241, 281	0

There are no RSRZ outliers to report.

6.2 Non-standard residues in protein, DNA, RNA chains [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	GTP	F	27	32/32	0.91	0.10	-	166,202,295,306	0
2	GTP	B	27	32/32	0.94	0.08	-	221,265,323,330	0
2	GTP	D	27	32/32	0.93	0.08	-	168,200,215,223	0

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
3	MG	C	101	1/1	0.38	0.33	-	128,128,128,128	0
3	MG	C	103	1/1	0.97	0.28	-	43,43,43,43	0
3	MG	B	101	1/1	0.96	0.16	-	147,147,147,147	1
3	MG	C	102	1/1	0.57	1.12	-	19,19,19,19	0

6.5 Other polymers [i](#)

There are no such residues in this entry.