W O R L D W I D E
wwPDB
P R O T E I N   D A T A   B A N K

wwPDB X-ray Structure Validation Summary Report (i)

Oct 14, 2023 – 09:52 PM EDT

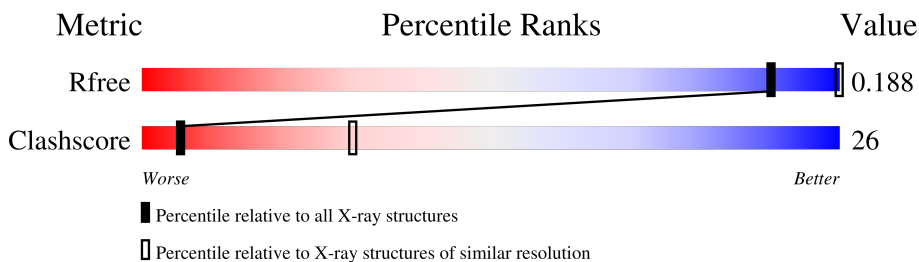|		
PDB ID	:	7U3R
Title	:	[2T7+10] Self-assembling tensegrity triangle with two turns of DNA and the sticky end attachment of a one-turn linker per axis, with R3 symmetry
Authors	:	Woloszyn, K.; Vecchioni, S.; Lu, B.; Seeman, N.C.; Sha, R.; Ohayon, Y.P.
Deposited on	:	2022-02-28
Resolution	:	6.27 Å(reported)

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 6.27 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	Percentile Ranks	Value
Rfree		0.188
Clashscore		26

Worse   *Better*

▌ Percentile relative to all X-ray structures

▯ Percentile relative to X-ray structures of similar resolution

Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R$_{free}$	130704	1007 (8.50-3.88)
Clashscore	141614	1056 (8.50-3.90)

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	C	8	38% · 62%
2	A	21	19% · 81%
3	Y	10	10% · 80% · 10%
4	X	10	100%
5	B	4	100%
6	D	6	17% · 83%
7	E	3	100%

WORLDWIDE
wwPDB
PROTEIN DATA BANK

2   Entry composition (i)

There are 7 unique types of molecules in this entry. The entry contains 1262 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 DNA chain called DNA (5'-D(*TP*CP*TP*GP*AP*TP*GP*T)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	C	8	Total	C	N	O	P	0	0	0
			161	79	26	49	7			

- Molecule 2 is a DNA chain called DNA (5'-D(*AP*CP*GP*CP*AP*GP*CP*CP*TP*GP* TP*AP*CP*GP*GP*AP*CP*AP*TP*CP*A)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	A	21	Total	C	N	O	P	0	0	0
			426	203	82	121	20			

- Molecule 3 is a DNA chain called DNA (5'-D(*GP*AP*TP*GP*CP*TP*GP*AP*GP*T)-3 ').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	Y	10	Total	C	N	O	P	0	0	0
			206	99	39	59	9			

- Molecule 4 is a DNA chain called DNA (5'-D(P*GP*TP*AP*CP*TP*CP*AP*GP*CP*A)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	X	10	Total	C	N	O	P	0	0	0
			204	97	38	59	10			

- Molecule 5 is a DNA chain called DNA (5'-D(P*CP*CP*GP*T)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
5	B	4	Total	C	N	O	P	0	0	0
			80	38	13	25	4			

- Molecule 6 is a DNA chain called DNA (5'-D(P*GP*GP*CP*TP*GP*C)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
6	D	6	Total 124	C 58	N 23	O 37	P 6	0	0	0

- Molecule 7 is a DNA chain called DNA (5'-D(P*AP*CP*A)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
7	E	3	Total 61	C 29	N 13	O 16	P 3	0	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.
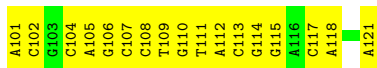
• Molecule 1: DNA (5'-D(*TP*CP*TP*GP*AP*TP*GP*T)-3')

Chain C:

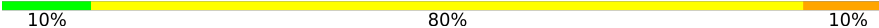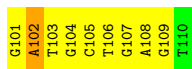• Molecule 2: DNA (5'-D(*AP*CP*GP*CP*AP*GP*CP*CP*TP*GP*TP*AP*CP*GP*GP*AP*CP*AP*TP*CP*A)-3')
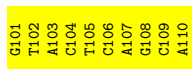
Chain A:

• Molecule 3: DNA (5'-D(*GP*AP*TP*GP*CP*TP*GP*AP*GP*T)-3')

Chain Y:

• Molecule 4: DNA (5'-D(P*GP*TP*AP*CP*TP*CP*AP*GP*CP*A)-3')

Chain X:

• Molecule 5: DNA (5'-D(P*CP*CP*GP*T)-3')

Chain B:

• Molecule 6: DNA (5'-D(P*GP*GP*CP*TP*GP*C)-3')

Chain D:

G109 G110 C111 T112 G113 C114

• Molecule 7: DNA (5'-D(P*AP*CP*A)-3')

Chain E: 100%

A105 C106 A107

4 Data and refinement statistics ⓘ

Property	Value	Source
Space group	H 3	Depositor
Cell constants a, b, c, α, β, γ	157.55Å 157.55Å 133.47Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	36.41 – 6.27 95.41 – 6.27	Depositor EDS
% Data completeness (in resolution range)	66.6 (36.41-6.27) 58.7 (95.41-6.27)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$<I/\sigma(I)>$ [1]	0.50 (at 6.19Å)	Xtriage
Refinement program	PHENIX 1.18.2_3874	Depositor
R, R_{free}	0.118 , 0.187 0.118 , 0.188	Depositor DCC
R_{free} test set	116 reflections (6.32%)	wwPDB-VP
Wilson B-factor (Å^2)	327.1	Xtriage
Anisotropy	1.047	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.32 , 500.0	EDS
L-test for twinning[2]	$<|L|>$ = 0.46, $<L^2>$ = 0.29	Xtriage
Estimated twinning fraction	0.064 for h,-h-k,-l	Xtriage
F_o,F_c correlation	0.81	EDS
Total number of atoms	1262	wwPDB-VP
Average B, all atoms (Å^2)	663.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 11.69% of the height of the origin peak. No significant pseudotranslation is detected.*

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[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   ⓘ

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	C	0.59	0/179	1.19	0/275
2	A	0.70	0/478	0.96	0/735
3	Y	0.81	0/231	1.17	1/356 (0.3%)
4	X	0.75	0/228	0.96	0/349
5	B	0.54	0/88	1.05	0/133
6	D	0.75	0/138	0.94	0/211
7	E	0.56	0/68	0.82	0/102
All	All	0.70	0/1410	1.03	1/2161 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
3	Y	102	DA	O4'-C1'-N9	5.45	111.82	108.00

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	C	161	0	94	4	0
2	A	426	0	236	23	2
3	Y	206	0	115	10	0
4	X	204	0	113	6	2
5	B	80	0	46	5	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
6	D	124	0	68	6	0
7	E	61	0	34	3	0
All	All	1262	0	706	50	2

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 26.

The worst 5 of 50 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)
3:Y:103:DT:H2"	3:Y:104:DG:H5"	1.64	0.79
4:X:104:DC:H2"	4:X:105:DT:H5"	1.66	0.77
2:A:101:DA:H5"	2:A:102:DC:H5	1.51	0.74
2:A:106:DG:H1	6:D:111:DC:H42	1.39	0.69
2:A:101:DA:H5"	2:A:102:DC:C5	2.29	0.65

All (2) 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:A:102:DC:O2	4:X:101:DG:N2[4_455]	1.39	0.81
2:A:102:DC:O2	4:X:101:DG:C2[4_455]	2.07	0.13

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

There are no protein molecules in this entry.

5.3.2 Protein sidechains ⓘ

There are no protein molecules in this entry.

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates ⓘ

There are no monosaccharides in this entry.

5.6 Ligand geometry ⓘ

There are no ligands in this entry.

5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

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.