

Full wwPDB X-ray Structure Validation Report (i)

May 23, 2022 – 10:02 AM EDT

PDB ID	:	7TZV
Title	:	Structure of DriD C-domain bound to 9mer ssDNA
Authors	:	Schumacher, M.A.; Laub, M.
Deposited on		
Resolution	:	1.65 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:

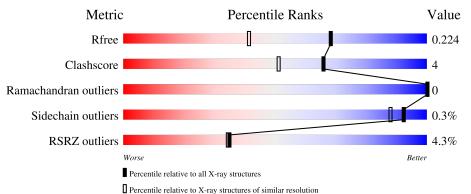
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.28.1
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.28.1

1 Overall quality at a glance (i)

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

The reported resolution of this entry is 1.65 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	$\begin{array}{c} \textbf{Whole archive} \\ (\#\textbf{Entries}) \end{array}$	${f Similar\ resolution}\ (\#{ m Entries,\ resolution\ range}({ m \AA}))$
R_{free}	130704	1827 (1.66-1.66)
Clashscore	141614	1931 (1.66-1.66)
Ramachandran outliers	138981	1891 (1.66-1.66)
Sidechain outliers	138945	1891 (1.66-1.66)
RSRZ outliers	127900	1791 (1.66-1.66)

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% 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	Qua	ality of chain	
1	А	198	5%	91%	6% • •
1	М	198	4%	91%	5% • •
2	J	9	44%	44%	11%



2 Entry composition (i)

There are 3 unique types of molecules in this entry. The entry contains 3640 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.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	Δ	193	Total	С	Ν	0	S	0	0	0
1	A	195	1515	960	278	273	4	0	0	0
1	М	192	Total	С	Ν	0	S	0	0	0
	1 M	192	1500	951	273	272	4	0	0	U

• Molecule 1 is a protein called WYL domain-containing protein.

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
А	130	GLY	-	expression tag	UNP Q9A999
А	131	SER	-	expression tag	UNP Q9A999
А	132	HIS	-	expression tag	UNP Q9A999
А	133	MET	-	expression tag	UNP Q9A999
М	130	GLY	-	expression tag	UNP Q9A999
М	131	SER	-	expression tag	UNP Q9A999
М	132	HIS	-	expression tag	UNP Q9A999
М	133	MET	-	expression tag	UNP Q9A999

• Molecule 2 is a DNA chain called DNA (5'-D(*TP*AP*GP*TP*CP*TP*AP*CP*T)-3').

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace		
0	т	0	Total	С	Ν	Ο	Р	0	0	0
	J	9	179	88	29	54	8	0	0	0

• Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	А	191	Total O 191 191	0	0
3	М	237	Total O 237 237	0	0
3	J	18	Total O 18 18	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 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: WYL domain-containing protein



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants	58.02Å 85.34Å 93.33Å	Depositor
a, b, c, α , β , γ	90.00° 90.00° 90.00°	Depositor
Resolution (Å)	31.49 - 1.65	Depositor
Resolution (A)	31.49 - 1.65	EDS
% Data completeness	98.1 (31.49-1.65)	Depositor
(in resolution range)	98.1 (31.49-1.65)	EDS
R _{merge}	(Not available)	Depositor
R _{sym}	0.05	Depositor
$< I/\sigma(I) > 1$	1.56 (at 1.65 Å)	Xtriage
Refinement program	PHENIX 1.17.1_3660	Depositor
D D.	0.192 , 0.223	Depositor
R, R_{free}	0.192 , 0.224	DCC
R_{free} test set	1998 reflections (3.61%)	wwPDB-VP
Wilson B-factor $(Å^2)$	26.9	Xtriage
Anisotropy	0.350	Xtriage
Bulk solvent $k_{sol}(e/Å^3), B_{sol}(Å^2)$	0.32 , 43.7	EDS
L-test for twinning ²	$ \langle L \rangle = 0.50, \langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	3640	wwPDB-VP
Average B, all atoms $(Å^2)$	37.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 5.31% of the height of the origin peak. No significant pseudotranslation is detected.

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



¹Intensities estimated from amplitudes.

5 Model quality (i)

5.1 Standard geometry (i)

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	Boi	nd lengths	Bond angles		
	Mol Chain		# Z > 5	RMSZ	# Z > 5	
1	А	0.67	2/1553~(0.1%)	0.73	0/2103	
1	М	0.74	1/1538~(0.1%)	0.84	2/2085~(0.1%)	
2	J	1.96	5/199~(2.5%)	1.69	8/305~(2.6%)	
All	All	0.84	8/3290~(0.2%)	0.88	10/4493~(0.2%)	

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Ζ	$\operatorname{Observed}(\operatorname{\AA})$	Ideal(Å)
2	J	23	DC	C3'-O3'	-9.21	1.31	1.44
1	М	202	LYS	C-N	8.40	1.50	1.34
1	А	146	GLU	CB-CG	5.82	1.63	1.52
2	J	21	DT	C3'-O3'	-5.77	1.36	1.44
1	А	146	GLU	CG-CD	5.75	1.60	1.51
2	J	21	DT	C5-C6	5.67	1.38	1.34
2	J	20	DC	C3'-O3'	-5.33	1.37	1.44
2	J	24	DT	P-OP1	-5.09	1.40	1.49

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$Observed(^{o})$	$Ideal(^{o})$
2	J	23	DC	O5'-P-OP1	-9.29	97.33	105.70
2	J	20	DC	O5'-P-OP2	-7.70	98.77	105.70
2	J	23	DC	OP1-P-OP2	7.13	130.29	119.60
2	J	24	DT	OP1-P-OP2	6.96	130.04	119.60
1	М	191	ASN	CB-CA-C	6.12	122.64	110.40
1	М	275	ASP	CB-CG-OD2	5.83	123.55	118.30
2	J	24	DT	O5'-P-OP2	-5.74	100.54	105.70
2	J	22	DA	N1-C2-N3	-5.71	126.44	129.30
2	J	24	DT	N3-C4-O4	5.66	123.29	119.90
2	J	24	DT	C5-C4-O4	-5.43	121.10	124.90

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	А	1515	0	1492	15	0
1	М	1500	0	1468	16	0
2	J	179	0	105	1	0
3	А	191	0	0	3	1
3	J	18	0	0	2	0
3	М	237	0	0	2	1
All	All	3640	0	3065	25	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 4.

All (25) 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:261:ARG:HH22	1:M:190:SER:HB3	1.34	0.91
1:A:261:ARG:NH2	1:M:190:SER:HB3	1.91	0.85
1:M:190:SER:O	1:M:192:TYR:CE2	2.46	0.68
1:A:261:ARG:NH1	1:M:190:SER:HB2	2.10	0.67
1:M:192:TYR:CE1	1:M:207:ARG:HG3	2.36	0.60
1:M:202:LYS:NZ	3:M:401:HOH:O	1.93	0.60
1:M:191:ASN:HB2	1:M:208:LEU:HD12	1.84	0.58
1:A:196:LEU:HD22	1:A:201:GLY:O	2.04	0.57
1:A:261:ARG:NH1	1:M:190:SER:CB	2.70	0.55
1:M:136:ILE:HG22	3:M:536:HOH:O	2.10	0.50
1:M:190:SER:OG	1:M:207:ARG:HG2	2.12	0.49
1:A:146:GLU:HG2	3:A:500:HOH:O	2.12	0.48
1:A:261:ARG:HH12	1:M:190:SER:CB	2.26	0.48
1:A:255:ARG:NH2	3:A:408:HOH:O	2.47	0.48
1:A:236:SER:HB3	3:J:107:HOH:O	2.14	0.47
1:A:261:ARG:NH2	1:M:190:SER:CB	2.72	0.45
1:M:260:LEU:HD23	1:M:270:VAL:HG11	1.99	0.45
1:A:165:SER:HA	1:A:178:ARG:O	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:322:ARG:HG3	3:A:553:HOH:O	2.16	0.44
1:A:261:ARG:CZ	1:M:190:SER:CB	2.96	0.44
1:A:303:ILE:O	1:A:310:LYS:HE2	2.17	0.43
1:A:245:GLN:HB3	1:A:306:PRO:HG3	2.01	0.42
1:M:144:PRO:HG3	1:M:190:SER:HA	2.02	0.42
1:M:260:LEU:CD2	1:M:270:VAL:HG11	2.50	0.42
2:J:21:DT:H72	3:J:116:HOH:O	2.20	0.42

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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)
3:A:418:HOH:O	3:M:560:HOH:O[4_545]	2.13	0.07

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	Favoured	Allowed	Outliers	Perce	\mathbf{ntiles}
1	А	191/198~(96%)	189 (99%)	2(1%)	0	100	100
1	М	190/198~(96%)	187 (98%)	3~(2%)	0	100	100
All	All	381/396~(96%)	376~(99%)	5 (1%)	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



Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
1	А	153/158~(97%)	153 (100%)	0	100 100
1	М	151/158~(96%)	150 (99%)	1 (1%)	84 73
All	All	304/316~(96%)	303 (100%)	1 (0%)	92 88

analysed, and the total number of residues.

All (1) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	М	165	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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 monosaccharides in this entry.

5.6 Ligand geometry (i)

There are no ligands in this entry.

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 $>$	#RSRZ>2	$OWAB(Å^2)$	Q < 0.9
1	А	193/198~(97%)	0.17	10 (5%) 27 26	22, 37, 60, 75	0
1	М	192/198~(96%)	0.04	7 (3%) 42 43	19, 30, 54, 64	0
2	J	9/9~(100%)	-0.20	0 100 100	32, 37, 44, 47	0
All	All	394/405~(97%)	0.10	17 (4%) 35 34	19, 34, 57, 75	0

All (17) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	А	225	GLN	4.6
1	А	327	TRP	4.6
1	М	190	SER	4.1
1	А	226	ASP	4.0
1	А	235	GLU	3.9
1	М	175	GLY	3.8
1	А	199	LYS	3.7
1	М	327	TRP	2.8
1	А	274	ALA	2.8
1	М	136	ILE	2.7
1	А	198	GLY	2.7
1	А	176	ARG	2.3
1	М	173	THR	2.3
1	М	226	ASP	2.1
1	А	236	SER	2.1
1	А	200	GLY	2.1
1	М	174	PRO	2.1

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 monosaccharides in this entry.

6.4 Ligands (i)

There are no ligands in this entry.

6.5 Other polymers (i)

There are no such residues in this entry.

