



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 12, 2017 – 07:59 pm GMT

PDB ID : 2NL8
Title : The origin binding domain of the SV40 large T antigen bound non specifically to a 17 bp palindrome DNA (sites 1 and 3)
Authors : Martynowski, D.; Bochkareva, E.; Bochkarev, A.
Deposited on : 2006-10-19
Resolution : 2.30 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

<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
Xtriage (Phenix) : 1.9-1692
EDS : trunk28620
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)
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) : recalc28949

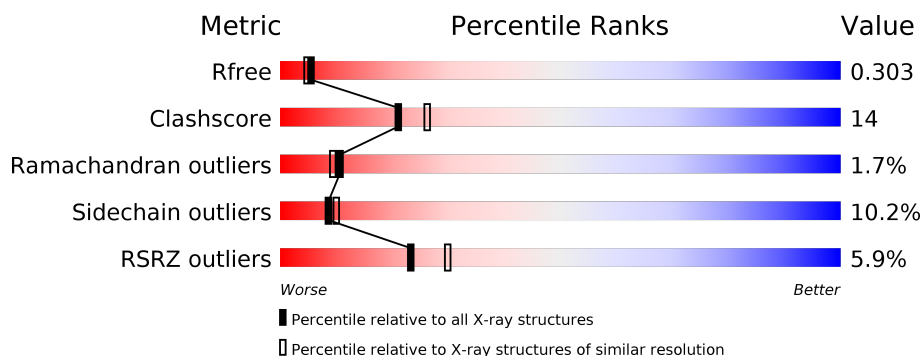
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION



The reported resolution of this entry is 2.30 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}	100719	4130 (2.30-2.30)
Clashscore	112137	4751 (2.30-2.30)
Ramachandran outliers	110173	4705 (2.30-2.30)
Sidechain outliers	110143	4704 (2.30-2.30)
RSRZ outliers	101464	4156 (2.30-2.30)

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	W	18	
2	A	133	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 1406 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 18-nt PEN element of the SV40 DNA origin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	W	18	Total	C	N	O	P	0	0	0
			363	173	67	106	17			

- Molecule 2 is a protein called large T antigen.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	A	117	Total	C	N	O	S	0	0	0
			961	626	159	172	4			

There are 5 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	127	GLY	-	CLONING ARTIFACT	UNP P03070
A	128	SER	-	CLONING ARTIFACT	UNP P03070
A	129	HIS	-	CLONING ARTIFACT	UNP P03070
A	130	MET	-	CLONING ARTIFACT	UNP P03070
A	216	SER	CYS	ENGINEERED	UNP P03070

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	56	Total	O	0	0
			56	56		
3	W	26	Total	O	0	0
			26	26		

4 Data and refinement statistics

Property	Value	Source
Space group	P 43 21 2	Depositor
Cell constants a, b, c, α , β , γ	43.87Å 43.87Å 238.31Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.92 – 2.30 19.62 – 2.30	Depositor EDS
% Data completeness (in resolution range)	94.6 (19.92-2.30) 94.6 (19.62-2.30)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.04	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.14 (at 2.30Å)	Xtriage
Refinement program	REFMAC 5.2	Depositor
R, R_{free}	0.237 , 0.299 0.233 , 0.303	Depositor DCC
R_{free} test set	518 reflections (5.15%)	DCC
Wilson B-factor (Å ²)	49.8	Xtriage
Anisotropy	0.019	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 40.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	1406	wwPDB-VP
Average B, all atoms (Å ²)	43.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.28% 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, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

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	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	W	1.71	3/406 (0.7%)	2.68	36/624 (5.8%)
2	A	0.88	0/987	0.91	2/1334 (0.1%)
All	All	1.18	3/1393 (0.2%)	1.69	38/1958 (1.9%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	W	21	DT	C1'-N1	7.24	1.58	1.49
1	W	14	DG	C3'-O3'	-5.71	1.36	1.44
1	W	20	DA	C5-C4	-5.35	1.35	1.38

All (38) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	W	21	DT	C2-N3-C4	-11.96	120.03	127.20
1	W	25	DC	O4'-C1'-N1	11.24	115.87	108.00
1	W	21	DT	N3-C2-O2	-11.07	115.66	122.30
1	W	21	DT	O4'-C1'-C2'	-10.70	97.34	105.90
1	W	21	DT	C4-C5-C6	10.43	124.26	118.00
1	W	21	DT	N1-C2-N3	10.12	120.67	114.60
1	W	24	DC	O4'-C1'-N1	10.01	115.00	108.00
2	A	156	LEU	CA-CB-CG	9.98	138.26	115.30
1	W	14	DG	P-O5'-C5'	9.31	135.79	120.90
1	W	22	DG	O4'-C1'-N9	9.01	114.31	108.00
1	W	20	DA	N1-C2-N3	-8.88	124.86	129.30
1	W	21	DT	C5-C6-N1	-8.46	118.62	123.70
1	W	19	DC	O4'-C1'-C2'	-7.45	99.94	105.90
1	W	25	DC	N1-C2-O2	7.37	123.32	118.90
1	W	17	DA	O4'-C4'-C3'	-7.18	101.63	104.50
1	W	23	DG	O4'-C1'-N9	6.99	112.89	108.00
1	W	15	DC	N3-C4-N4	-6.90	113.17	118.00
1	W	11	DG	C3'-C2'-C1'	-6.79	94.35	102.50
1	W	21	DT	O4'-C4'-C3'	-6.74	101.81	104.50

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	W	20	DA	C2-N3-C4	6.64	113.92	110.60
1	W	17	DA	C5-N7-C8	-6.62	100.59	103.90
1	W	21	DT	C6-C5-C7	-6.56	118.97	122.90
1	W	15	DC	O4'-C4'-C3'	-6.54	101.89	104.50
1	W	13	DG	O4'-C1'-C2'	-6.51	100.69	105.90
1	W	24	DC	O4'-C1'-C2'	-6.42	100.77	105.90
1	W	19	DC	N1-C2-O2	6.36	122.71	118.90
1	W	19	DC	N3-C2-O2	-6.30	117.49	121.90
1	W	14	DG	O4'-C4'-C3'	-6.15	102.04	104.50
1	W	12	DA	O4'-C1'-N9	-6.01	103.79	108.00
1	W	15	DC	C5-C4-N4	5.75	124.22	120.20
1	W	12	DA	N1-C2-N3	-5.61	126.50	129.30
1	W	14	DG	C5-C6-N1	5.50	114.25	111.50
1	W	17	DA	N7-C8-N9	5.32	116.46	113.80
1	W	22	DG	C2-N3-C4	5.24	114.52	111.90
1	W	22	DG	N1-C2-N3	-5.23	120.76	123.90
1	W	26	DT	O4'-C1'-N1	-5.14	104.40	108.00
2	A	195	LEU	CA-CB-CG	5.09	127.00	115.30
1	W	14	DG	N3-C4-C5	-5.06	126.07	128.60

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	W	363	0	203	2	1
2	A	961	0	958	32	1
3	A	56	0	0	7	0
3	W	26	0	0	0	0
All	All	1406	0	1161	34	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 14.

All (34) 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:A:239:ASP:HA	2:A:241:PHE:H	1.26	0.97
2:A:144:SER:HB3	3:A:45:HOH:O	1.77	0.84
2:A:249:PRO:HB3	3:A:82:HOH:O	1.84	0.78
2:A:246:GLU:HG3	2:A:248:LEU:O	1.84	0.78
2:A:238:ARG:O	2:A:239:ASP:HB2	1.87	0.74
2:A:214:LYS:HD2	3:A:62:HOH:O	1.86	0.74
2:A:210:ASN:O	2:A:214:LYS:HD3	1.89	0.72
2:A:239:ASP:HA	2:A:241:PHE:N	2.03	0.72
2:A:246:GLU:OE2	2:A:249:PRO:HA	1.93	0.68
2:A:238:ARG:HG3	3:A:68:HOH:O	1.94	0.67
2:A:145:PHE:HA	2:A:228:LYS:HG3	1.81	0.62
2:A:239:ASP:CA	2:A:241:PHE:H	2.09	0.60
2:A:166:GLU:HB2	3:A:60:HOH:O	2.02	0.59
2:A:191:ASN:HD22	2:A:192:HIS:CD2	2.22	0.58
2:A:197:PHE:HZ	2:A:208:ILE:CD1	2.17	0.57
2:A:208:ILE:HG12	2:A:223:CYS:SG	2.45	0.56
1:W:13:DG:H2'	1:W:14:DG:OP2	2.06	0.56
2:A:235:ALA:O	2:A:241:PHE:HD1	1.92	0.53
2:A:191:ASN:H	2:A:191:ASN:ND2	2.06	0.52
2:A:139:PRO:HD2	2:A:142:LEU:HD12	1.93	0.51
2:A:188:ASN:HB2	3:A:80:HOH:O	2.10	0.51
2:A:160:ALA:HA	2:A:195:LEU:O	2.12	0.50
2:A:187:HIS:O	2:A:193:ASN:HA	2.11	0.50
2:A:141:GLU:H	2:A:141:GLU:CD	2.15	0.50
2:A:197:PHE:CZ	2:A:208:ILE:CD1	2.95	0.49
2:A:181:VAL:HG13	2:A:183:PHE:O	2.11	0.49
2:A:249:PRO:HD3	3:A:61:HOH:O	2.14	0.47
2:A:155:THR:HB	2:A:202:ARG:HB3	1.97	0.47
2:A:192:HIS:C	2:A:193:ASN:HD22	2.19	0.45
2:A:134:ASP:O	2:A:136:LYS:NZ	2.50	0.45
2:A:238:ARG:O	2:A:239:ASP:CB	2.57	0.44
1:W:16:DC:H2'	1:W:17:DA:C8	2.53	0.44
2:A:197:PHE:CZ	2:A:208:ILE:HD12	2.56	0.41
2:A:181:VAL:HG11	2:A:184:ILE:HB	2.02	0.40

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)
1:W:21:DT:C7	2:A:152:SER:OG[8_665]	2.01	0.19

5.3 Torsion angles

5.3.1 Protein backbone

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	Percentiles
2	A	115/133 (86%)	101 (88%)	12 (10%)	2 (2%)	11 9

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	A	239	ASP
2	A	240	PRO

5.3.2 Protein sidechains

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	Percentiles
2	A	108/121 (89%)	97 (90%)	11 (10%)	8 10

All (11) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	A	141	GLU
2	A	156	LEU
2	A	162	TYR
2	A	171	LEU
2	A	191	ASN
2	A	193	ASN
2	A	195	LEU
2	A	208	ILE
2	A	214	LYS

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Mol	Chain	Res	Type
2	A	228	LYS
2	A	242	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
2	A	191	ASN
2	A	192	HIS

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](#)

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, 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	W	18/18 (100%)	-0.81	0	100	100	30, 42, 49, 49	0
2	A	117/133 (87%)	0.30	8 (6%)	18	24	26, 43, 64, 74	0
All	All	135/151 (89%)	0.15	8 (5%)	23	30	26, 43, 62, 74	0

All (8) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	A	230	TYR	4.7
2	A	151	PHE	3.7
2	A	248	LEU	3.6
2	A	238	ARG	3.5
2	A	194	ILE	3.4
2	A	195	LEU	2.5
2	A	160	ALA	2.4
2	A	161	ILE	2.2

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](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.