



wwPDB X-ray Structure Validation Summary Report ⓘ

Oct 15, 2017 – 12:12 PM EDT

PDB ID : 2R2T
Title : d(ATTTAGTTAACTAAAT) complexed with MMLV RT catalytic fragment
Authors : Goodwin, K.D.; Lewis, M.A.; Long, E.C.; Georgiadis, M.M.
Deposited on : unknown
Resolution : 2.00 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
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20030345
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)	:	rb-20030345

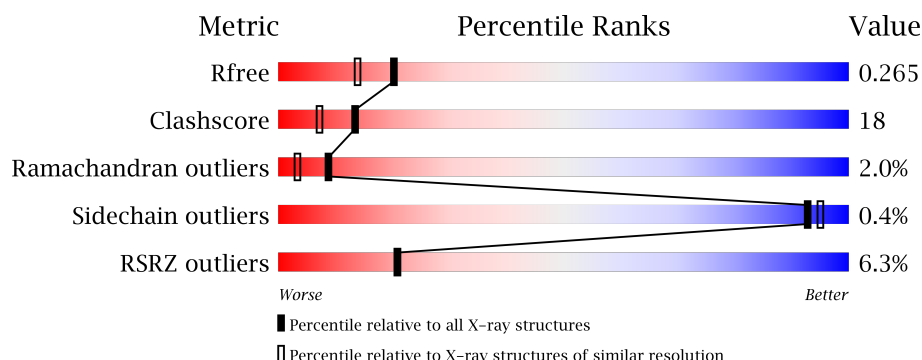
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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	6609 (2.00-2.00)
Clashscore	112137	7775 (2.00-2.00)
Ramachandran outliers	110173	7679 (2.00-2.00)
Sidechain outliers	110143	7678 (2.00-2.00)
RSRZ outliers	101464	6696 (2.00-2.00)

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	B	8	<div style="width: 100%; height: 10px; background-color: yellow;"></div> 100%
2	G	8	<div style="display: flex; justify-content: space-between; align-items: center;"> <div style="width: 13%; height: 10px; background-color: green;"></div> <div style="width: 88%; height: 10px; background-color: yellow;"></div> </div> 13% 88%
3	A	255	<div style="display: flex; justify-content: space-between; align-items: center;"> <div style="width: 7%; height: 10px; background-color: red;"></div> <div style="width: 75%; height: 10px; background-color: green;"></div> <div style="width: 24%; height: 10px; background-color: yellow;"></div> </div> 7% 75% 24% .

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 2542 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(*DAP*DTP*DTP*DTP*DAP*DGP*DTP*D T)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	B	8	Total	C	N	O	P	0	0	0
			161	80	25	49	7			

- Molecule 2 is a DNA chain called DNA (5'-D(P*DAP*DAP*DCP*DTP*DAP*DAP*DAP*DT)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	G	8	Total	C	N	O	P	0	0	0
			164	79	32	45	8			

- Molecule 3 is a protein called Reverse transcriptase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	A	255	Total	C	N	O	S	0	0	0
			2041	1311	356	367	7			

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	2	Total	O	0	0
			2	2		
4	G	3	Total	O	0	0
			3	3		
4	A	171	Total	O	0	0
			171	171		

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 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: DNA (5'-D(*DAP*DTP*DTP*DTP*DAP*DGP*DTP*DT)-3')

Chain B:  100%


A1
T2
T3
T4
A5
G6
T7
T8

- Molecule 2: DNA (5'-D(P*DAP*DAP*DCP*DTP*DAP*DAP*DAP*DT)-3')

Chain G:  13% 88%

A9
A10
C11
T12
A13
A14
A15
T16

- Molecule 3: Reverse transcriptase

Chain A:  7% 75% 24%

T24
W25
L26
S27
D28
F29
P30
Q31
I49
I50
P51
T57
P58
I61
M66
S67
Q68
E69
A70
R71
L72
G73
I74
K75
P76
P100
Y101
K102
K103
P104
G105
T106
M107
D108
E117
H126
P127
M131
P132
Y133
M134
L135
L136
S137
G138
L139
P140
W145
H161
P162

T163
R173
D174
P175
E176
M177
G178
I179
S180
G181
Q182
L183
D206
Q213
I218
V223
A229
A230
E233
L234
D235
C236
Q237
G248
N249
L250
Q263
V266
K274
E275
G276
Q277
R278

4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	54.64Å 145.80Å 46.88Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 2.00 43.72 – 2.00	Depositor EDS
% Data completeness (in resolution range)	92.3 (50.00-2.00) 92.4 (43.72-2.00)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.42 (at 2.00Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.224 , 0.264 0.225 , 0.265	Depositor DCC
R_{free} test set	1166 reflections (4.84%)	DCC
Wilson B-factor (Å ²)	29.1	Xtriage
Anisotropy	0.276	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 49.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	2542	wwPDB-VP
Average B, all atoms (Å ²)	40.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.18% 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	B	0.23	0/179	0.64	0/275
2	G	0.23	0/184	0.68	0/281
3	A	0.32	0/2097	0.59	0/2858
All	All	0.30	0/2460	0.60	0/3414

There are no bond length outliers.

There are no bond angle outliers.

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	B	161	0	95	17	1
2	G	164	0	91	16	1
3	A	2041	0	2056	50	0
4	A	171	0	0	3	0
4	B	2	0	0	0	0
4	G	3	0	0	0	0
All	All	2542	0	2242	83	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 18.

The worst 5 of 83 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:G:14:DA:H2''	2:G:15:DA:H5'	1.40	1.03
3:A:233:GLU:HG2	3:A:237:GLN:HE21	1.32	0.93
1:B:3:DT:H2''	1:B:4:DT:H5'	1.51	0.91
1:B:5:DA:H2''	1:B:6:DG:H5''	1.60	0.83
2:G:12:DT:H2''	2:G:13:DA:H5'	1.67	0.77

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:B:8:DT:O3'	2:G:9:DA:P[2_765]	1.61	0.59

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	Percentiles
3	A	253/255 (99%)	241 (95%)	7 (3%)	5 (2%)	9 3

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	A	106	THR
3	A	223	VAL
3	A	104	PRO
3	A	177	MET
3	A	181	GLY

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 analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
3	A	224/224 (100%)	223 (100%)	1 (0%)	93	95

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

Mol	Chain	Res	Type
3	A	177	MET

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 11 such sidechains are listed below:

Mol	Chain	Res	Type
3	A	161	HIS
3	A	213	GLN
3	A	238	GLN
3	A	144	GLN
3	A	237	GLN

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 ⓘ

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	B	8/8 (100%)	0.46	0	100	100	46, 61, 75, 79	0
2	G	8/8 (100%)	0.76	0	100	100	33, 74, 85, 85	0
3	A	255/255 (100%)	0.49	17 (6%)	19	18	18, 32, 67, 85	0
All	All	271/271 (100%)	0.49	17 (6%)	21	21	18, 33, 74, 85	0

The worst 5 of 17 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	A	106	THR	10.1
3	A	105	GLY	9.3
3	A	179	ILE	7.8
3	A	178	GLY	7.0
3	A	180	SER	6.8

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.