



Full wwPDB X-ray Structure Validation Report

Feb 27, 2014 – 07:47 PM GMT

PDB ID : 2R2T
Title : d(ATTTAGTTAACTAAAT) complexed with MMLV RT catalytic fragment
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Deposited on : 2007-08-27
Resolution : 2.00 Å(reported)

This is a full wwPDB 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/ValidationPDFNotes.html>

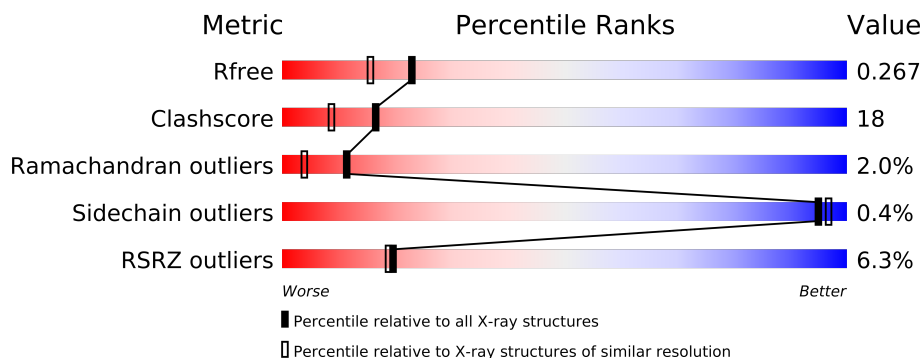
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : dev-1323
EDS : stable22639
Percentile statistics : 21963
Refmac : 5.8.0049
CCP4 : 6.3.0 (Settle)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable22683

1 Overall quality at a glance

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}	66092	4888 (2.00-2.00)
Clashscore	79885	6188 (2.00-2.00)
Ramachandran outliers	78287	6102 (2.00-2.00)
Sidechain outliers	78261	6100 (2.00-2.00)
RSRZ outliers	66119	4890 (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. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	B	8	<div><div></div><div></div></div>
2	G	8	<div><div></div><div></div></div>
3	A	255	<div><div></div><div></div></div>

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 2542 atoms, of which 0 are hydrogen and 0 are deuterium.

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	A	171	Total	O	0	0
			171	171		
4	B	2	Total	O	0	0
			2	2		
4	G	3	Total	O	0	0
			3	3		

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

Chain B: 

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: 

A9
A10
C11
T12
A13
A14
A15
T16

- Molecule 3: Reverse transcriptase

Chain A: 

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
V101
K102
K103
P104
G105
T106
N107
D108
E117
H126
P127
N131
P132
Y133
N134
L135
L136
S137
G138
L139
P140
W145
H161
P162

T163
R173
D174
P175
E176
M177
G178
I179
S180
G181
Q182
L183
D206
Q213
T218
V223
A229
A230
E233
L234
D235
C236
Q237
Q248
N249
L250
Q263
V266
F267
K274
E275
Q276
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 0.233 , 0.267	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)	0.35 , 40.2	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	1 of 25116 reflections (0.004%)	Xtriage
F_o, F_c correlation	0.93	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.

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	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 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

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

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 18.

All (83) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(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
1:B:1:DA:H2''	1:B:2:DT:H5''	1.67	0.76
3:A:206:ASP:HB3	3:A:250:LEU:HD13	1.69	0.75
3:A:274:LYS:HG2	3:A:275:GLU:HG3	1.71	0.71
1:B:2:DT:H2''	1:B:3:DT:C5'	2.22	0.69
2:G:14:DA:C2'	2:G:15:DA:H5'	2.21	0.69
3:A:104:PRO:HG2	3:A:107:ASN:HB2	1.74	0.68
1:B:5:DA:C2'	1:B:6:DG:H5''	2.22	0.68
1:B:6:DG:H2''	1:B:7:DT:H5'	1.75	0.68
1:B:1:DA:H2''	1:B:2:DT:C5'	2.24	0.67
1:B:6:DG:H2''	1:B:7:DT:C5'	2.25	0.67
3:A:102:LYS:HD2	3:A:106:THR:HA	1.76	0.66
2:G:11:DC:H2''	2:G:12:DT:C5'	2.27	0.65
1:B:5:DA:H2''	1:B:6:DG:C5'	2.27	0.64
2:G:13:DA:H1'	2:G:14:DA:H5''	1.79	0.64
3:A:174:ASP:HB3	3:A:178:GLY:HA2	1.80	0.64
1:B:1:DA:C2'	1:B:2:DT:H5''	2.27	0.64
1:B:2:DT:H2''	1:B:3:DT:H5'	1.81	0.63
3:A:70:ALA:HB1	3:A:100:PRO:HB3	1.82	0.62
3:A:161:HIS:HD2	3:A:163:THR:H	1.49	0.61
3:A:173:ARG:HD2	4:A:434:HOH:O	2.02	0.59
3:A:68:GLN:O	3:A:72:LEU:HD13	2.02	0.59
3:A:103:LYS:HB2	3:A:104:PRO:HD3	1.84	0.59
1:B:3:DT:H2''	1:B:4:DT:C5'	2.29	0.59
3:A:161:HIS:CD2	3:A:163:THR:H	2.23	0.57
3:A:233:GLU:O	3:A:237:GLN:HG3	2.06	0.56
2:G:10:DA:H1'	2:G:11:DC:H5''	1.89	0.55
2:G:11:DC:H2''	2:G:12:DT:H5''	1.89	0.55
3:A:132:PRO:O	3:A:136:LEU:HD23	2.08	0.54
3:A:61:ILE:HD11	3:A:117:GLU:HG3	1.89	0.53
2:G:13:DA:C2'	2:G:14:DA:H5''	2.39	0.53
1:B:2:DT:H2''	1:B:3:DT:H5''	1.90	0.53
3:A:274:LYS:HE2	3:A:275:GLU:OE1	2.08	0.53
2:G:13:DA:H2''	2:G:14:DA:C5'	2.39	0.53
2:G:11:DC:H2''	2:G:12:DT:H5'	1.92	0.53
3:A:179:ILE:HG22	3:A:180:SER:N	2.24	0.52
3:A:233:GLU:HG2	3:A:237:GLN:NE2	2.13	0.52
3:A:173:ARG:HH11	3:A:173:ARG:HG2	1.74	0.52
3:A:26:LEU:O	3:A:26:LEU:HD23	2.08	0.52

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:G:13:DA:H2''	2:G:14:DA:H5''	1.93	0.51
3:A:24:THR:HB	3:A:27:SER:OG	2.11	0.51
2:G:13:DA:C1'	2:G:14:DA:H5''	2.41	0.50
1:B:2:DT:C2'	1:B:3:DT:H5''	2.42	0.50
2:G:9:DA:H2''	2:G:10:DA:C8	2.48	0.49
3:A:104:PRO:CG	3:A:107:ASN:HB2	2.42	0.48
3:A:145:TRP:CH2	3:A:233:GLU:HB2	2.48	0.48
3:A:179:ILE:HD13	3:A:183:LEU:HD21	1.95	0.48
3:A:57:THR:HG23	3:A:58:PRO:HD2	1.96	0.47
1:B:2:DT:H1'	1:B:3:DT:H5''	1.97	0.47
3:A:71:ARG:HD3	4:A:428:HOH:O	2.15	0.47
3:A:75:LYS:HB3	3:A:76:PRO:HD3	1.96	0.47
3:A:179:ILE:O	3:A:180:SER:C	2.52	0.46
2:G:11:DC:C2'	2:G:12:DT:H5''	2.46	0.46
3:A:138:GLY:O	3:A:140:PRO:HD3	2.14	0.46
3:A:277:GLN:OE1	3:A:277:GLN:HA	2.15	0.46
3:A:102:LYS:CD	3:A:106:THR:HA	2.43	0.46
3:A:105:GLY:O	3:A:106:THR:HB	2.17	0.45
3:A:31:GLN:HE22	3:A:249:ASN:HD22	1.64	0.45
1:B:4:DT:H2''	1:B:5:DA:C8	2.51	0.45
3:A:277:GLN:O	3:A:278:ARG:HB2	2.16	0.45
3:A:248:GLY:HA3	4:A:444:HOH:O	2.16	0.45
3:A:173:ARG:NH1	3:A:181:GLY:HA2	2.32	0.45
3:A:29:PHE:N	3:A:30:PRO:HD3	2.32	0.45
3:A:263:GLN:HB2	3:A:266:VAL:CG1	2.48	0.44
3:A:173:ARG:NH1	3:A:173:ARG:HG2	2.33	0.44
3:A:131:ASN:HD21	3:A:134:ASN:ND2	2.16	0.43
3:A:49:ILE:O	3:A:51:PRO:HD3	2.18	0.43
3:A:213:GLN:HA	3:A:213:GLN:NE2	2.34	0.43
3:A:218:ILE:HB	3:A:229:ALA:HB3	2.00	0.43
2:G:12:DT:C2'	2:G:13:DA:H5'	2.44	0.42
3:A:26:LEU:C	3:A:26:LEU:HD23	2.40	0.41
2:G:11:DC:H1'	2:G:12:DT:H5''	2.03	0.41
3:A:66:MET:HG2	3:A:70:ALA:HB3	2.02	0.41
1:B:6:DG:H2''	1:B:7:DT:H5''	2.00	0.41
3:A:126:HIS:HA	3:A:127:PRO:HD3	1.95	0.41
3:A:145:TRP:CZ2	3:A:233:GLU:HB2	2.56	0.41
3:A:74:ILE:HD11	3:A:100:PRO:HA	2.02	0.41
3:A:71:ARG:CZ	3:A:175:PRO:HG2	2.51	0.40
3:A:230:ALA:HB3	3:A:236:CYS:HB2	2.04	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	Distance(Å)	Clash(Å)
1:B:8:DT:O3'	2:G:9:DA:P[2_765]	1.61	0.59

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
3	A	253/255 (99%)	241 (95%)	7 (3%)	5 (2%)	11 4

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 ⓘ

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%)	95 97

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. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
3	A	31	GLN
3	A	84	GLN
3	A	134	ASN
3	A	144	GLN
3	A	161	HIS
3	A	213	GLN
3	A	214	HIS
3	A	237	GLN
3	A	238	GLN
3	A	245	GLN
3	A	249	ASN

5.3.3 RNA ⓘ

There are no RNA chains 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 carbohydrates 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 ⓘ

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(Å²)	Q<0.9
1	B	8/8 (100%)	0.36	0	100 100	46, 61, 75, 79	0
2	G	8/8 (100%)	0.65	0	100 100	33, 74, 85, 85	0
3	A	255/255 (100%)	0.36	17 (6%)	17 17	18, 32, 67, 85	0
All	All	271/271 (100%)	0.37	17 (6%)	19 19	18, 33, 74, 85	0

All (17) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	A	106	THR	9.4
3	A	105	GLY	8.4
3	A	179	ILE	7.6
3	A	180	SER	7.1
3	A	178	GLY	4.9
3	A	177	MET	4.1
3	A	103	LYS	3.5
3	A	104	PRO	3.5
3	A	176	GLU	3.4
3	A	173	ARG	3.2
3	A	175	PRO	2.5
3	A	108	ASP	2.5
3	A	107	ASN	2.3
3	A	234	LEU	2.1
3	A	102	LYS	2.0
3	A	267	LYS	2.0
3	A	237	GLN	2.0

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

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

6.3 Carbohydrates

There are no carbohydrates in this entry.

6.4 Ligands

There are no ligands in this entry.

6.5 Other polymers

There are no such residues in this entry.