



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 13, 2017 – 12:41 pm GMT

PDB ID : 2ATW
Title : Structure of a Mycobacterium tuberculosis NusA-RNA complex
Authors : Beuth, B.; Pennell, S.; Arnvig, K.B.; Martin, S.R.; Taylor, I.A.
Deposited on : 2005-08-26
Resolution : 2.25 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

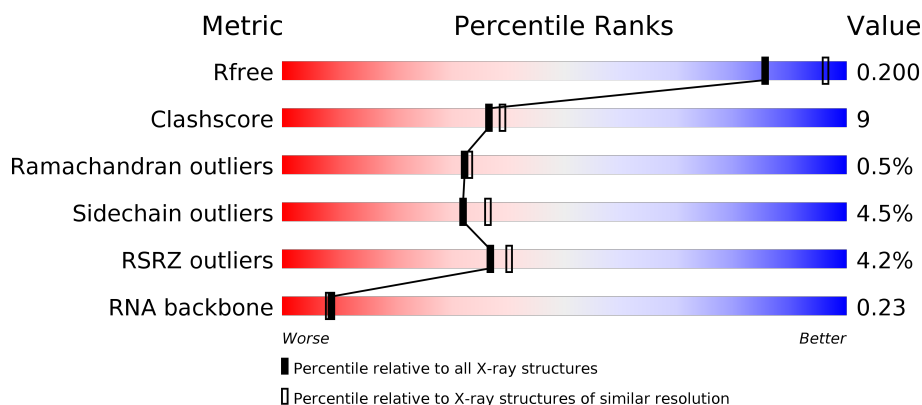
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.25 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	1062 (2.26-2.26)
Clashscore	112137	1178 (2.26-2.26)
Ramachandran outliers	110173	1145 (2.26-2.26)
Sidechain outliers	110143	1146 (2.26-2.26)
RSRZ outliers	101464	1066 (2.26-2.26)
RNA backbone	2435	1122 (2.80-1.72)

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	12	
1	D	12	
2	A	251	
2	C	251	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 4083 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 RNA chain called ribosomal RNA (5'- AGAACUCAAUAG -3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	B	12	Total	C	N	O	P	0	0	0
			255	116	50	78	11			
1	D	12	Total	C	N	O	P	0	0	0
			255	116	50	78	11			

- Molecule 2 is a protein called Transcription elongation protein nusA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	A	223	Total	C	N	O	S	0	0	0
			1676	1037	320	315	4			
2	C	225	Total	C	N	O	S	0	0	0
			1689	1044	322	319	4			

There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	348	LEU	-	CLONING ARTIFACT	UNP P0A5M2
A	349	GLU	-	CLONING ARTIFACT	UNP P0A5M2
A	350	HIS	-	EXPRESSION TAG	UNP P0A5M2
A	351	HIS	-	EXPRESSION TAG	UNP P0A5M2
A	352	HIS	-	EXPRESSION TAG	UNP P0A5M2
A	353	HIS	-	EXPRESSION TAG	UNP P0A5M2
A	354	HIS	-	EXPRESSION TAG	UNP P0A5M2
A	355	HIS	-	EXPRESSION TAG	UNP P0A5M2
C	348	LEU	-	CLONING ARTIFACT	UNP P0A5M2
C	349	GLU	-	CLONING ARTIFACT	UNP P0A5M2
C	350	HIS	-	EXPRESSION TAG	UNP P0A5M2
C	351	HIS	-	EXPRESSION TAG	UNP P0A5M2
C	352	HIS	-	EXPRESSION TAG	UNP P0A5M2
C	353	HIS	-	EXPRESSION TAG	UNP P0A5M2
C	354	HIS	-	EXPRESSION TAG	UNP P0A5M2
C	355	HIS	-	EXPRESSION TAG	UNP P0A5M2

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	102	Total 102	O 102	0	0
3	B	11	Total 11	O 11	0	0
3	C	90	Total 90	O 90	0	0
3	D	5	Total 5	O 5	0	0

4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	64.56Å 89.54Å 100.89Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	15.00 – 2.25 14.99 – 2.25	Depositor EDS
% Data completeness (in resolution range)	99.4 (15.00-2.25) 99.4 (14.99-2.25)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	0.08	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.28 (at 2.24Å)	Xtriage
Refinement program	REFMAC 5	Depositor
R, R_{free}	0.206 , 0.273 0.210 , 0.200	Depositor DCC
R_{free} test set	1402 reflections (5.25%)	DCC
Wilson B-factor (Å ²)	32.7	Xtriage
Anisotropy	0.014	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 35.7	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	4083	wwPDB-VP
Average B, all atoms (Å ²)	33.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.41% 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	1.42	2/286 (0.7%)	2.08	11/444 (2.5%)
1	D	1.31	2/286 (0.7%)	2.01	9/444 (2.0%)
2	A	0.77	0/1698	0.91	3/2300 (0.1%)
2	C	0.70	0/1711	0.94	5/2317 (0.2%)
All	All	0.86	4/3981 (0.1%)	1.18	28/5505 (0.5%)

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	3	A	N7-C5	-6.42	1.35	1.39
1	D	10	A	C5-C6	-5.63	1.35	1.41
1	B	1	G	C2-N3	5.54	1.37	1.32
1	D	0	A	N7-C5	-5.46	1.35	1.39

All (28) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	10	A	O4'-C1'-N9	-10.65	99.68	108.20
1	B	2	A	O4'-C1'-N9	9.98	116.19	108.20
1	D	2	A	O4'-C1'-N9	7.68	114.34	108.20
1	B	3	A	O4'-C1'-N9	-7.34	102.32	108.20
2	C	263	ASP	CB-CG-OD2	7.09	124.68	118.30
1	D	10	A	C4-C5-N7	7.07	114.23	110.70
1	D	10	A	N1-C6-N6	6.56	122.53	118.60
1	D	9	U	P-O3'-C3'	-6.48	111.93	119.70
2	C	259	ASP	CB-CG-OD2	6.45	124.11	118.30
2	A	284	ASP	CB-CG-OD2	6.44	124.10	118.30
1	B	1	G	C6-N1-C2	-6.36	121.28	125.10
1	B	7	A	C4'-C3'-C2'	-6.34	96.26	102.60
1	B	1	G	O4'-C1'-N9	-6.24	103.21	108.20
1	B	9	U	C5-C4-O4	-5.93	122.34	125.90
2	C	296	ASP	CB-CG-OD2	5.88	123.59	118.30
2	C	122	ASP	CB-CG-OD2	5.84	123.56	118.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	2	A	N1-C2-N3	-5.82	126.39	129.30
1	D	9	U	O4'-C1'-N1	5.77	112.82	108.20
2	A	122	ASP	CB-CG-OD2	5.77	123.50	118.30
2	A	326	ASP	CB-CG-OD2	5.47	123.22	118.30
1	D	10	A	N9-C4-C5	-5.42	103.63	105.80
1	B	7	A	C8-N9-C4	-5.35	103.66	105.80
1	D	1	G	C6-C5-N7	-5.27	127.24	130.40
1	B	4	C	O4'-C1'-N1	-5.25	104.00	108.20
1	B	10	A	P-O3'-C3'	5.25	126.00	119.70
1	D	6	C	P-O3'-C3'	5.22	125.96	119.70
1	B	11	G	O4'-C1'-N9	-5.13	104.09	108.20
2	C	261	ASP	CB-CG-OD2	5.11	122.90	118.30

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	B	255	0	132	4	0
1	D	255	0	132	2	0
2	A	1676	0	1717	30	1
2	C	1689	0	1726	32	0
3	A	102	0	0	14	0
3	B	11	0	0	1	0
3	C	90	0	0	12	1
3	D	5	0	0	0	0
All	All	4083	0	3707	68	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 9.

All (68) 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:329:PRO:HD2	3:A:2021:HOH:O	1.61	0.99
2:A:185:HIS:HD2	2:A:187:ASN:H	1.11	0.94
2:A:138:GLU:HG3	3:A:2039:HOH:O	1.68	0.93
2:C:185:HIS:HD2	2:C:187:ASN:H	1.15	0.90
2:C:113:GLU:HG3	3:C:1035:HOH:O	1.73	0.88
2:C:176:GLU:O	2:C:176:GLU:HG2	1.73	0.87
2:C:320:ARG:HD3	3:C:2038:HOH:O	1.76	0.86
3:B:2012:HOH:O	2:A:315:ARG:HD2	1.78	0.84
2:A:284:ASP:CG	3:A:1081:HOH:O	2.16	0.82
2:C:121:ARG:CG	2:C:121:ARG:HH11	1.94	0.80
2:C:106:GLU:OE2	2:C:172:ARG:NH1	2.15	0.80
2:A:288:ARG:HD3	3:A:1072:HOH:O	1.80	0.79
2:A:185:HIS:CD2	2:A:187:ASN:H	1.98	0.79
2:A:320:ARG:HD3	3:A:1133:HOH:O	1.84	0.77
2:C:134:ARG:NH1	3:C:1109:HOH:O	2.04	0.75
2:C:185:HIS:CD2	2:C:187:ASN:H	2.02	0.74
2:C:296:ASP:HB2	3:C:2004:HOH:O	1.88	0.74
2:C:110:ARG:NE	3:C:1030:HOH:O	2.28	0.67
2:A:134:ARG:NE	3:A:1028:HOH:O	2.22	0.67
1:B:9:U:OP1	1:B:11:G:O3'	2.12	0.66
2:C:247:MET:HG2	2:C:255:ILE:HD12	1.76	0.66
2:A:126:ASN:O	2:A:129:GLY:N	2.30	0.64
2:A:185:HIS:HD2	2:A:187:ASN:N	1.89	0.64
2:A:288:ARG:CD	3:A:1072:HOH:O	2.41	0.64
2:A:284:ASP:OD1	2:A:287:ALA:HB3	1.97	0.63
2:A:134:ARG:NH2	3:A:1028:HOH:O	2.24	0.62
2:C:121:ARG:HG3	2:C:121:ARG:HH11	1.64	0.62
2:C:185:HIS:CD2	2:C:186:PRO:HD2	2.35	0.61
2:C:121:ARG:HG2	2:C:121:ARG:HH11	1.67	0.59
2:C:229:LEU:HD12	2:C:230:ASN:H	1.70	0.57
1:D:9:U:OP1	1:D:11:G:O3'	2.19	0.56
2:C:229:LEU:HD12	2:C:230:ASN:N	2.21	0.56
2:C:107:PHE:HZ	2:C:172:ARG:HB3	1.73	0.54
2:A:134:ARG:CZ	3:A:1028:HOH:O	2.56	0.53
2:A:130:LEU:HD13	2:A:145:VAL:HG21	1.91	0.53
2:A:284:ASP:OD1	2:A:287:ALA:CB	2.56	0.53
2:C:276:LYS:HE3	3:C:2051:HOH:O	2.09	0.52
2:C:176:GLU:HB2	3:C:1087:HOH:O	2.08	0.52
2:C:327:ALA:N	3:C:2004:HOH:O	2.25	0.51
1:D:10:A:H2'	1:D:10:A:N3	2.24	0.51
2:A:138:GLU:CG	3:A:2039:HOH:O	2.38	0.51
2:C:176:GLU:N	3:C:2003:HOH:O	2.45	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:219:LYS:HG2	2:A:256:ASP:HB2	1.94	0.49
2:C:303:ILE:HG22	2:C:307:GLY:HA2	1.95	0.48
2:A:122:ASP:C	2:A:122:ASP:OD1	2.52	0.48
2:C:159:HIS:HB3	3:C:1054:HOH:O	2.13	0.48
2:A:285:GLN:HG2	3:A:1157:HOH:O	2.16	0.46
2:C:145:VAL:HG22	2:C:146:ILE:N	2.32	0.44
1:B:7:A:H2'	1:B:7:A:O5'	2.17	0.44
2:A:320:ARG:NH1	3:A:1143:HOH:O	2.49	0.44
2:C:270:ASN:O	2:C:273:SER:HB2	2.17	0.44
2:C:121:ARG:CG	2:C:121:ARG:NH1	2.65	0.44
2:C:139:THR:OG1	2:C:140:LYS:N	2.50	0.43
2:A:244:ARG:HG3	2:A:247:MET:HE3	1.99	0.43
2:A:251:SER:HA	3:A:1193:HOH:O	2.18	0.43
2:C:130:LEU:HB3	2:C:145:VAL:CG2	2.49	0.43
2:C:138:GLU:OE1	3:C:1185:HOH:O	2.21	0.43
2:C:106:GLU:HA	3:C:2006:HOH:O	2.17	0.42
2:C:175:ARG:O	2:C:176:GLU:HB3	2.19	0.42
2:A:329:PRO:CD	3:A:2021:HOH:O	2.40	0.42
2:A:146:ILE:HG22	2:A:151:GLN:HG3	2.02	0.42
2:A:195:LEU:HD23	2:A:195:LEU:HA	1.83	0.41
2:A:112:GLY:HA2	2:A:166:TYR:CZ	2.54	0.41
2:A:273:SER:OG	2:A:274:PRO:HA	2.20	0.41
2:A:172:ARG:H	2:A:172:ARG:HG2	1.67	0.41
1:B:9:U:H4'	1:B:10:A:C4	2.57	0.40
2:C:280:VAL:HG22	2:C:292:VAL:HG22	2.04	0.40
1:B:10:A:H4'	1:B:11:G:O5'	2.22	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)
2:A:140:LYS:NZ	3:C:1128:HOH:O[1_554]	2.18	0.02

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	
2	A	221/251 (88%)	210 (95%)	10 (4%)	1 (0%)	32	33
2	C	223/251 (89%)	216 (97%)	6 (3%)	1 (0%)	38	42
All	All	444/502 (88%)	426 (96%)	16 (4%)	2 (0%)	32	33

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	A	174	ALA
2	C	176	GLU

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	176/199 (88%)	170 (97%)	6 (3%)	42	52
2	C	177/199 (89%)	167 (94%)	10 (6%)	25	25
All	All	353/398 (89%)	337 (96%)	16 (4%)	32	36

All (16) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	A	156	SER
2	A	172	ARG
2	A	175	ARG
2	A	277	VAL
2	A	279	SER
2	A	281	SER
2	C	110	ARG
2	C	121	ARG
2	C	138	GLU
2	C	140	LYS
2	C	143	GLU

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Mol	Chain	Res	Type
2	C	176	GLU
2	C	239	MET
2	C	248	SER
2	C	276	LYS
2	C	277	VAL

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
2	A	126	ASN
2	A	159	HIS
2	A	185	HIS
2	A	187	ASN
2	C	161	ASN
2	C	185	HIS
2	C	187	ASN
2	C	285	GLN

5.3.3 RNA ⓘ

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
1	B	11/12 (91%)	3 (27%)	0
1	D	11/12 (91%)	3 (27%)	0
All	All	22/24 (91%)	6 (27%)	0

All (6) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
1	B	7	A
1	B	10	A
1	B	11	G
1	D	7	A
1	D	10	A
1	D	11	G

There are no RNA pucker outliers to report.

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

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(Å ²)	Q<0.9
1	B	12/12 (100%)	-0.44	0 100 100	30, 32, 37, 38	0
1	D	12/12 (100%)	-0.35	0 100 100	28, 36, 42, 50	0
2	A	223/251 (88%)	0.12	11 (4%) 30 34	19, 28, 51, 64	0
2	C	225/251 (89%)	0.16	9 (4%) 39 42	21, 31, 50, 70	0
All	All	472/526 (89%)	0.12	20 (4%) 37 40	19, 30, 51, 70	0

All (20) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	C	329	PRO	9.8
2	C	175	ARG	7.3
2	C	328	PRO	7.1
2	A	329	PRO	5.5
2	C	327	ALA	5.2
2	A	172	ARG	4.4
2	C	174	ALA	4.0
2	A	107	PHE	3.9
2	A	124	ARG	3.1
2	A	175	ARG	3.1
2	C	320	ARG	2.7
2	A	176	GLU	2.7
2	A	127	ALA	2.6
2	A	121	ARG	2.5
2	C	107	PHE	2.4
2	C	127	ALA	2.4
2	A	158	GLU	2.3
2	A	174	ALA	2.3
2	C	153	PRO	2.3
2	A	159	HIS	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 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.