



Full wwPDB X-ray Structure Validation Report

Nov 7, 2014 – 07:27 PM EST

PDB ID : 2Q1W
Title : Crystal structure of the Bordetella bronchiseptica enzyme WbmH in complex with NAD+
Authors : King, J.D.; Harmer, N.J.; Maskell, D.J.; Blundell, T.L.
Deposited on : 2007-05-25
Resolution : 2.19 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>

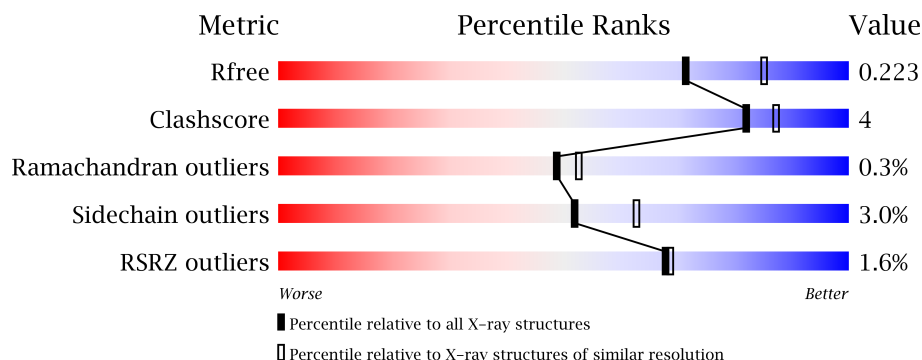
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.16 November 2013
Xtriage (Phenix) : dev-1439
EDS : trunk24195
Percentile statistics : 23426
Refmac : 5.8.0049
CCP4 : 6.1.3
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : trunk24195

1 Overall quality at a glance

The reported resolution of this entry is 2.19 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}	77520	3336 (2.20-2.20)
Clashscore	88313	4037 (2.20-2.20)
Ramachandran outliers	86584	3966 (2.20-2.20)
C α geometry	86677	3967 (2.20-2.20)
Sidechain outliers	86556	3967 (2.20-2.20)
RSRZ outliers	77580	3339 (2.20-2.20)

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	A	333	
1	B	333	
1	C	333	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 7367 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 protein called Putative nucleotide sugar epimerase/ dehydratase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	300	Total	C	N	O	S	8	1	0
			2286	1453	396	431	6			
1	B	299	Total	C	N	O	S	0	0	0
			2302	1464	399	433	6			
1	C	296	Total	C	N	O	S	0	0	0
			2232	1421	384	421	6			

There are 60 discrepancies between the modelled and reference sequences:

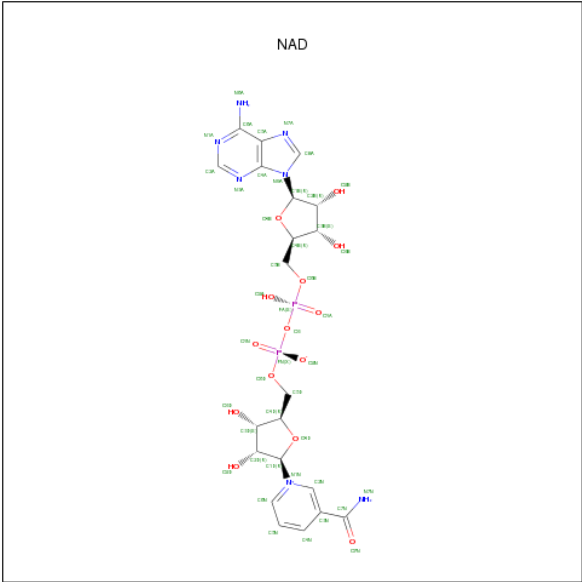
Chain	Residue	Modelled	Actual	Comment	Reference
A	-19	MET	-	EXPRESSION TAG	UNP O87987
A	-18	GLY	-	EXPRESSION TAG	UNP O87987
A	-17	SER	-	EXPRESSION TAG	UNP O87987
A	-16	SER	-	EXPRESSION TAG	UNP O87987
A	-15	HIS	-	INSERTION	UNP O87987
A	-14	HIS	-	INSERTION	UNP O87987
A	-13	HIS	-	INSERTION	UNP O87987
A	-12	HIS	-	INSERTION	UNP O87987
A	-11	HIS	-	INSERTION	UNP O87987
A	-10	HIS	-	INSERTION	UNP O87987
A	-9	SER	-	EXPRESSION TAG	UNP O87987
A	-8	SER	-	EXPRESSION TAG	UNP O87987
A	-7	GLY	-	EXPRESSION TAG	UNP O87987
A	-6	LEU	-	EXPRESSION TAG	UNP O87987
A	-5	VAL	-	EXPRESSION TAG	UNP O87987
A	-4	PRO	-	EXPRESSION TAG	UNP O87987
A	-3	ARG	-	EXPRESSION TAG	UNP O87987
A	-2	GLY	-	EXPRESSION TAG	UNP O87987
A	-1	SER	-	EXPRESSION TAG	UNP O87987
A	0	HIS	-	EXPRESSION TAG	UNP O87987
B	-19	MET	-	EXPRESSION TAG	UNP O87987
B	-18	GLY	-	EXPRESSION TAG	UNP O87987
B	-17	SER	-	EXPRESSION TAG	UNP O87987

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-16	SER	-	EXPRESSION TAG	UNP O87987
B	-15	HIS	-	INSERTION	UNP O87987
B	-14	HIS	-	INSERTION	UNP O87987
B	-13	HIS	-	INSERTION	UNP O87987
B	-12	HIS	-	INSERTION	UNP O87987
B	-11	HIS	-	INSERTION	UNP O87987
B	-10	HIS	-	INSERTION	UNP O87987
B	-9	SER	-	EXPRESSION TAG	UNP O87987
B	-8	SER	-	EXPRESSION TAG	UNP O87987
B	-7	GLY	-	EXPRESSION TAG	UNP O87987
B	-6	LEU	-	EXPRESSION TAG	UNP O87987
B	-5	VAL	-	EXPRESSION TAG	UNP O87987
B	-4	PRO	-	EXPRESSION TAG	UNP O87987
B	-3	ARG	-	EXPRESSION TAG	UNP O87987
B	-2	GLY	-	EXPRESSION TAG	UNP O87987
B	-1	SER	-	EXPRESSION TAG	UNP O87987
B	0	HIS	-	EXPRESSION TAG	UNP O87987
C	-19	MET	-	EXPRESSION TAG	UNP O87987
C	-18	GLY	-	EXPRESSION TAG	UNP O87987
C	-17	SER	-	EXPRESSION TAG	UNP O87987
C	-16	SER	-	EXPRESSION TAG	UNP O87987
C	-15	HIS	-	INSERTION	UNP O87987
C	-14	HIS	-	INSERTION	UNP O87987
C	-13	HIS	-	INSERTION	UNP O87987
C	-12	HIS	-	INSERTION	UNP O87987
C	-11	HIS	-	INSERTION	UNP O87987
C	-10	HIS	-	INSERTION	UNP O87987
C	-9	SER	-	EXPRESSION TAG	UNP O87987
C	-8	SER	-	EXPRESSION TAG	UNP O87987
C	-7	GLY	-	EXPRESSION TAG	UNP O87987
C	-6	LEU	-	EXPRESSION TAG	UNP O87987
C	-5	VAL	-	EXPRESSION TAG	UNP O87987
C	-4	PRO	-	EXPRESSION TAG	UNP O87987
C	-3	ARG	-	EXPRESSION TAG	UNP O87987
C	-2	GLY	-	EXPRESSION TAG	UNP O87987
C	-1	SER	-	EXPRESSION TAG	UNP O87987
C	0	HIS	-	EXPRESSION TAG	UNP O87987

- Molecule 2 is NICOTINAMIDE-ADENINE-DINUCLEOTIDE (three-letter code: NAD) (formula: C₂₁H₂₇N₇O₁₄P₂).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			44	21	7	14	2		
2	B	1	Total	C	N	O	P	0	0
			44	21	7	14	2		
2	C	1	Total	C	N	O	P	0	0
			44	21	7	14	2		

- Molecule 3 is water.

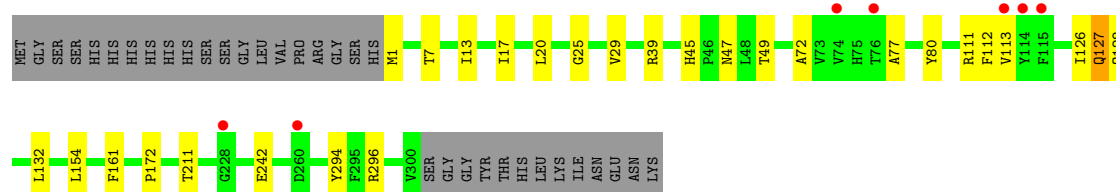
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	124	Total	O	0	0
			124	124		
3	B	180	Total	O	0	0
			180	180		
3	C	111	Total	O	0	0
			111	111		

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.

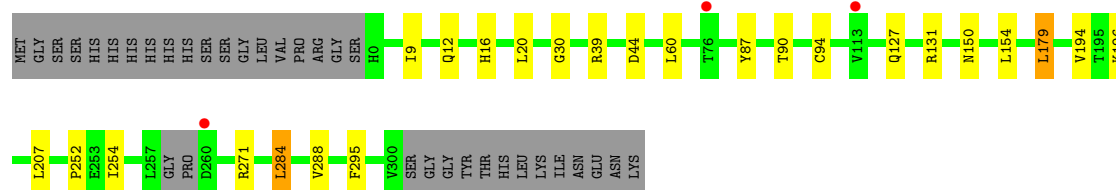
- Molecule 1: Putative nucleotide sugar epimerase/ dehydratase

Chain A: 



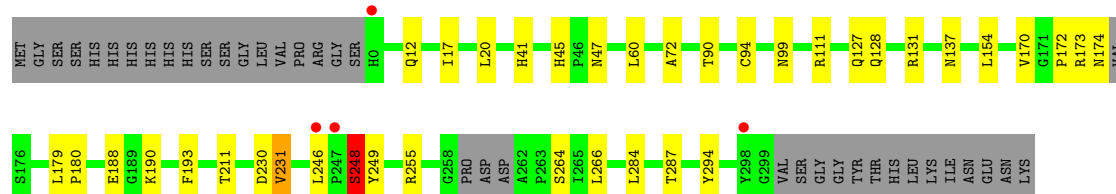
- Molecule 1: Putative nucleotide sugar epimerase/ dehydratase

Chain B: 



- Molecule 1: Putative nucleotide sugar epimerase/ dehydratase

Chain C: 



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	93.67Å 158.99Å 68.22Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.00 – 2.19 29.73 – 2.19	Depositor EDS
% Data completeness (in resolution range)	93.2 (30.00-2.19) 93.2 (29.73-2.19)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.05 (at 2.18Å)	Xtriage
Refinement program	REFMAC 5.0	Depositor
R, R_{free}	0.176 , 0.227 0.175 , 0.223	Depositor DCC
R_{free} test set	2559 reflections (5.40%)	DCC
Wilson B-factor (Å ²)	30.0	Xtriage
Anisotropy	0.093	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 28.8	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	1 of 49944 reflections (0.002%)	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	7367	wwPDB-VP
Average B, all atoms (Å ²)	29.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.95% 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, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: NAD

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	A	0.75	0/2343	0.72	0/3193
1	B	0.80	0/2355	0.77	2/3204 (0.1%)
1	C	0.73	0/2284	0.71	0/3114
All	All	0.76	0/6982	0.73	2/9511 (0.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	C	0	1

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	271	ARG	NE-CZ-NH1	9.50	125.05	120.30
1	B	271	ARG	NE-CZ-NH2	-7.66	116.47	120.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	C	248	SER	Peptide

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	A	2286	0	2195	14	0
1	B	2302	0	2230	12	0
1	C	2232	0	2112	23	0
2	A	44	0	26	1	0
2	B	44	0	26	1	0
2	C	44	0	26	0	0
3	A	124	0	0	0	0
3	B	180	0	0	0	0
3	C	111	0	0	0	0
All	All	7367	0	6615	49	0

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 4.

All (49) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:296[B]:ARG:HH11	1:A:296[B]:ARG:HG2	1.51	0.75
1:C:127:GLN:NE2	1:C:131:ARG:H	1.86	0.74
1:C:45:HIS:HD2	1:C:47:ASN:H	1.37	0.72
1:B:127:GLN:NE2	1:B:131:ARG:H	1.92	0.67
1:C:231:VAL:HG21	1:C:284:LEU:HD22	1.79	0.65
1:A:45:HIS:HD2	1:A:47:ASN:H	1.49	0.61
1:C:193:PHE:HB3	1:C:255:ARG:HH11	1.67	0.59
1:A:126:ILE:O	1:A:127:GLN:HB2	2.04	0.58
1:C:179:LEU:HB2	1:C:180:PRO:HD3	1.85	0.58
1:C:193:PHE:HB3	1:C:255:ARG:NH1	2.19	0.57
1:B:12:GLN:O	1:B:16:HIS:HD2	1.87	0.57
1:C:127:GLN:HE21	1:C:131:ARG:H	1.52	0.57
1:A:296[B]:ARG:NH1	1:A:296[B]:ARG:HG2	2.18	0.56
1:B:90:THR:O	1:B:94:CYS:HB2	2.07	0.55
1:C:17:ILE:HD13	1:C:211:THR:HG21	1.89	0.55
1:C:45:HIS:CD2	1:C:47:ASN:H	2.22	0.55
1:C:246:LEU:HB3	1:C:248:SER:HB2	1.90	0.54
1:B:179:LEU:HD22	1:B:295:PHE:HE2	1.72	0.54

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:45:HIS:CD2	1:A:47:ASN:H	2.26	0.53
1:B:194:VAL:CG2	1:B:254:ILE:HD13	2.40	0.52
1:C:72:ALA:HA	1:C:111:ARG:O	2.12	0.50
1:C:41:HIS:CD2	1:C:173:ARG:HG2	2.48	0.49
1:A:7:THR:HG1	1:A:77:ALA:H	1.62	0.48
1:C:128:GLN:HE22	1:C:266:LEU:HG	1.79	0.47
1:B:194:VAL:HG23	1:B:254:ILE:HD13	1.98	0.46
1:A:29:VAL:HG22	1:A:49:THR:HB	1.98	0.45
1:A:13:ILE:HD12	2:A:501:NAD:H51N	1.99	0.45
1:A:113:VAL:HG11	1:A:211:THR:HG23	1.99	0.45
1:B:87:TYR:HD1	1:C:99:ASN:HD21	1.65	0.45
1:C:128:GLN:NE2	1:C:266:LEU:HG	2.33	0.44
1:B:194:VAL:HG22	1:B:252:PRO:HB2	1.99	0.43
1:B:196:LYS:HD2	1:B:254:ILE:HG22	2.00	0.43
1:C:230:ASP:OD2	1:C:264:SER:OG	2.32	0.43
1:C:172:PRO:HB3	1:C:294:TYR:HB2	2.00	0.43
1:B:9:ILE:HD12	1:B:30:GLY:HA3	2.01	0.43
1:C:179:LEU:CB	1:C:180:PRO:HD3	2.49	0.43
1:A:172:PRO:HB3	1:A:294:TYR:HB2	2.01	0.42
1:B:284:LEU:HD22	1:B:288:VAL:HG23	2.01	0.42
1:C:188:GLU:OE1	1:C:190:LYS:HE2	2.19	0.42
1:C:179:LEU:HA	1:C:179:LEU:HD23	1.78	0.41
1:A:17:ILE:HD13	1:A:211:THR:HG21	2.02	0.41
1:C:12:GLN:NE2	1:C:174:ASN:HB2	2.35	0.41
1:C:90:THR:O	1:C:94:CYS:HB2	2.20	0.41
1:A:112:PHE:O	1:A:161:PHE:HA	2.20	0.41
1:C:170:VAL:HG23	1:C:287:THR:CG2	2.51	0.41
1:B:127:GLN:HE21	1:B:131:ARG:H	1.67	0.41
1:A:72:ALA:HA	1:A:111:ARG:O	2.21	0.41
1:A:1:MET:HG3	1:A:25:GLY:O	2.21	0.41
2:B:501:NAD:O1N	2:B:501:NAD:N7N	2.55	0.40

There are no symmetry-related clashes.

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	
1	A	299/333 (90%)	291 (97%)	6 (2%)	2 (1%)	29	27
1	B	295/333 (89%)	288 (98%)	7 (2%)	0	100	100
1	C	290/333 (87%)	275 (95%)	14 (5%)	1 (0%)	49	53
All	All	884/999 (88%)	854 (97%)	27 (3%)	3 (0%)	49	53

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	127	GLN
1	A	80	TYR
1	C	249	TYR

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	
1	A	236/277 (85%)	230 (98%)	6 (2%)	58	70
1	B	241/277 (87%)	232 (96%)	9 (4%)	44	53
1	C	226/277 (82%)	220 (97%)	6 (3%)	56	67
All	All	703/831 (85%)	682 (97%)	21 (3%)	52	63

All (21) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	20	LEU
1	A	39	ARG
1	A	128	GLN
1	A	132	LEU
1	A	154	LEU
1	A	242	GLU
1	B	20	LEU
1	B	39	ARG
1	B	44	ASP

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Mol	Chain	Res	Type
1	B	60	LEU
1	B	150	ASN
1	B	154	LEU
1	B	179	LEU
1	B	207	LEU
1	B	284	LEU
1	C	20	LEU
1	C	60	LEU
1	C	137	ASN
1	C	154	LEU
1	C	231	VAL
1	C	248	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (19) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	12	GLN
1	A	45	HIS
1	A	69	GLN
1	A	99	ASN
1	A	140	ASN
1	A	150	ASN
1	B	12	GLN
1	B	16	HIS
1	B	99	ASN
1	B	102	GLN
1	B	127	GLN
1	C	12	GLN
1	C	45	HIS
1	C	69	GLN
1	C	99	ASN
1	C	102	GLN
1	C	127	GLN
1	C	137	ASN
1	C	140	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 ⓘ

3 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. 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 $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# $ Z > 2$	Counts	RMSZ	# $ Z > 2$
2	NAD	A	501	-	48,48,48	1.58	5 (10%)	73,73,73	2.02	10 (13%)
2	NAD	B	501	-	48,48,48	1.53	4 (8%)	73,73,73	2.06	12 (16%)
2	NAD	C	501	-	48,48,48	1.69	5 (10%)	73,73,73	2.01	12 (16%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the chemical component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	NAD	A	501	-	-	0/30/62/62	0/5/5/5
2	NAD	B	501	-	-	0/30/62/62	0/5/5/5
2	NAD	C	501	-	-	0/30/62/62	0/5/5/5

All (14) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	501	NAD	O7N-C7N	8.29	1.42	1.24
2	B	501	NAD	O7N-C7N	7.28	1.40	1.24
2	A	501	NAD	O7N-C7N	7.24	1.40	1.24
2	C	501	NAD	C2A-N3A	4.18	1.39	1.32

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	501	NAD	C2A-N3A	3.78	1.38	1.32
2	B	501	NAD	C2A-N3A	3.48	1.38	1.32
2	A	501	NAD	C2N-N1N	3.13	1.39	1.35
2	B	501	NAD	C2A-N1A	3.11	1.40	1.33
2	C	501	NAD	C2N-N1N	2.79	1.38	1.35
2	A	501	NAD	C2A-N1A	2.71	1.39	1.33
2	C	501	NAD	O4D-C1D	2.70	1.44	1.41
2	C	501	NAD	C2A-N1A	2.46	1.38	1.33
2	B	501	NAD	C2N-N1N	2.41	1.38	1.35
2	A	501	NAD	C4A-N3A	2.02	1.38	1.35

All (34) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	501	NAD	N3A-C2A-N1A	-12.79	117.64	128.89
2	C	501	NAD	N3A-C2A-N1A	-12.00	118.33	128.89
2	A	501	NAD	N3A-C2A-N1A	-11.65	118.64	128.89
2	C	501	NAD	O4D-C1D-N1N	4.83	113.44	108.13
2	A	501	NAD	C5A-C4A-N3A	-4.83	121.27	125.98
2	C	501	NAD	C5A-C4A-N3A	-4.55	121.55	125.98
2	A	501	NAD	O4D-C1D-N1N	4.50	113.07	108.13
2	B	501	NAD	O4D-C1D-C2D	-4.37	100.33	106.69
2	A	501	NAD	C3N-C7N-N7N	4.20	122.46	117.78
2	A	501	NAD	N3A-C4A-N9A	4.12	132.45	125.39
2	A	501	NAD	O7N-C7N-C3N	-3.75	115.40	119.59
2	B	501	NAD	C2A-N1A-C6A	3.43	124.86	118.76
2	C	501	NAD	N3A-C4A-N9A	3.24	130.94	125.39
2	B	501	NAD	C5A-C4A-N3A	-3.10	122.96	125.98
2	B	501	NAD	C4D-O4D-C1D	3.08	113.11	109.72
2	C	501	NAD	C4A-C5A-N7A	-3.08	106.44	109.41
2	B	501	NAD	N3A-C4A-N9A	3.02	130.57	125.39
2	B	501	NAD	C8A-N9A-C4A	2.98	109.38	106.96
2	A	501	NAD	C2A-N3A-C4A	2.85	121.47	113.27
2	C	501	NAD	C2A-N1A-C6A	2.73	123.62	118.76
2	A	501	NAD	O3D-C3D-C2D	-2.64	103.29	111.83
2	C	501	NAD	O4B-C1B-C2B	-2.59	102.92	106.69
2	B	501	NAD	C1B-N9A-C4A	-2.53	122.26	126.64
2	C	501	NAD	PN-O3-PA	-2.47	125.10	131.93
2	B	501	NAD	O4D-C1D-N1N	2.39	110.76	108.13
2	C	501	NAD	C8A-N7A-C5A	2.33	110.80	103.58
2	A	501	NAD	C1B-N9A-C4A	-2.25	122.74	126.64
2	A	501	NAD	O4B-C1B-C2B	-2.23	103.44	106.69

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	501	NAD	C6A-C5A-C4A	2.22	120.04	117.55
2	C	501	NAD	N7A-C8A-N9A	-2.21	107.18	112.20
2	B	501	NAD	N7A-C8A-N9A	-2.13	107.36	112.20
2	B	501	NAD	O2N-PN-O3	2.10	115.09	105.14
2	C	501	NAD	C2A-N3A-C4A	2.08	119.26	113.27
2	B	501	NAD	C2A-N3A-C4A	2.07	119.23	113.27

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

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(Å ²)	Q<0.9
1	A	300/333 (90%)	-0.15	7 (2%) 57 57	20, 29, 46, 55	1 (0%)
1	B	299/333 (89%)	-0.34	3 (1%) 78 80	15, 23, 35, 49	1 (0%)
1	C	296/333 (88%)	-0.24	4 (1%) 71 72	17, 29, 58, 66	0
All	All	895/999 (89%)	-0.24	14 (1%) 68 69	15, 27, 46, 66	2 (0%)

All (14) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	260	ASP	3.0
1	C	247	PRO	3.0
1	A	76	THR	2.9
1	A	74	VAL	2.8
1	A	115	PHE	2.8
1	A	113	VAL	2.6
1	B	113	VAL	2.5
1	A	114	TYR	2.3
1	C	298	TYR	2.3
1	C	246	LEU	2.3
1	C	0	HIS	2.2
1	B	260	ASP	2.0
1	B	76	THR	2.0
1	A	228	GLY	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

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q < 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSR	LLDF	B-factors(Å ²)	Q<0.9
2	NAD	B	501	44/44	0.10	-0.27	13,18,25,26	0
2	NAD	A	501	44/44	0.11	-0.29	17,23,36,39	0
2	NAD	C	501	44/44	0.09	-0.51	13,22,35,37	0

6.5 Other polymers

There are no such residues in this entry.