



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 13, 2017 – 11:50 pm GMT

PDB ID : 3QBD
Title : 3-Dehydroquinate Synthase (aroB) from Mycobacterium tuberculosis in complex with NAD
Authors : Cheng, W.C.; Chen, T.J.; Wang, W.C.
Deposited on : 2011-01-13
Resolution : 2.47 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
Mogul	:	1.7.2 (RC1), CSD as538be (2017)
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

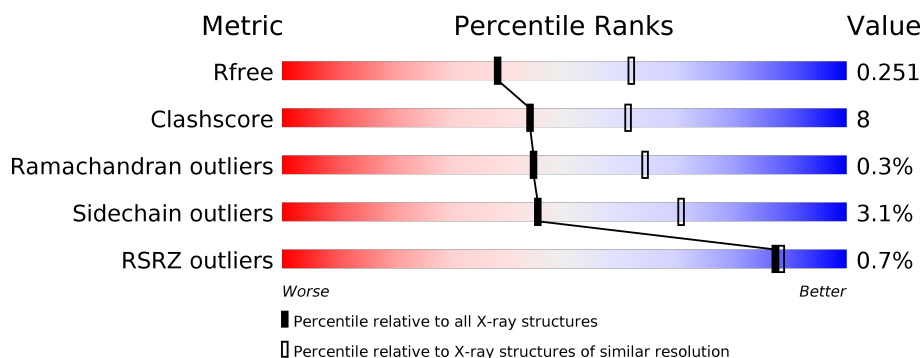
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.47 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	4719 (2.50-2.46)
Clashscore	112137	5483 (2.50-2.46)
Ramachandran outliers	110173	5388 (2.50-2.46)
Sidechain outliers	110143	5390 (2.50-2.46)
RSRZ outliers	101464	4754 (2.50-2.46)

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	A	368	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 1%, orange 1%, yellow 11%, green 80%, grey 7%);"></div> <div style="display: flex; justify-content: space-between; font-size: small;"> % 80% 11% •• 7% </div> </div>
1	B	368	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, green 76%, yellow 16%, orange 7%);"></div> <div style="display: flex; justify-content: space-between; font-size: small;"> 76% 16% • 7% </div> </div>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 5260 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 protein called 3-dehydroquinate synthase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	344	Total	C	N	O	S	0	0	0
			2546	1621	451	467	7			
1	B	342	Total	C	N	O	S	0	0	0
			2537	1616	449	465	7			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-5	HIS	-	EXPRESSION TAG	UNP A5U5N7
A	-4	HIS	-	EXPRESSION TAG	UNP A5U5N7
A	-3	HIS	-	EXPRESSION TAG	UNP A5U5N7
A	-2	HIS	-	EXPRESSION TAG	UNP A5U5N7
A	-1	HIS	-	EXPRESSION TAG	UNP A5U5N7
A	0	HIS	-	EXPRESSION TAG	UNP A5U5N7
B	-5	HIS	-	EXPRESSION TAG	UNP A5U5N7
B	-4	HIS	-	EXPRESSION TAG	UNP A5U5N7
B	-3	HIS	-	EXPRESSION TAG	UNP A5U5N7
B	-2	HIS	-	EXPRESSION TAG	UNP A5U5N7
B	-1	HIS	-	EXPRESSION TAG	UNP A5U5N7
B	0	HIS	-	EXPRESSION TAG	UNP A5U5N7

- 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		

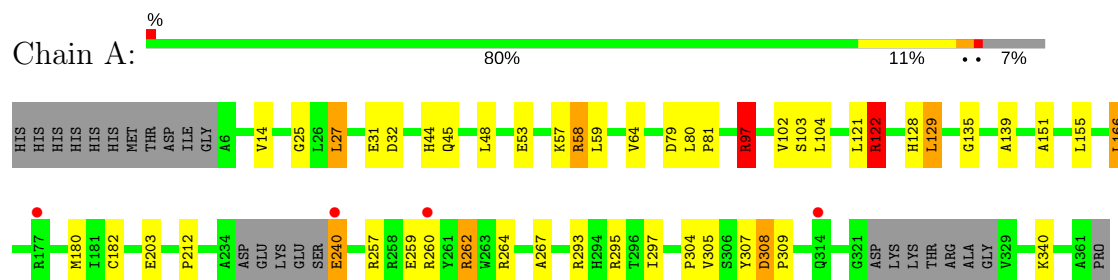
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	53	Total	O	0	0
			53	53		
3	B	36	Total	O	0	0
			36	36		

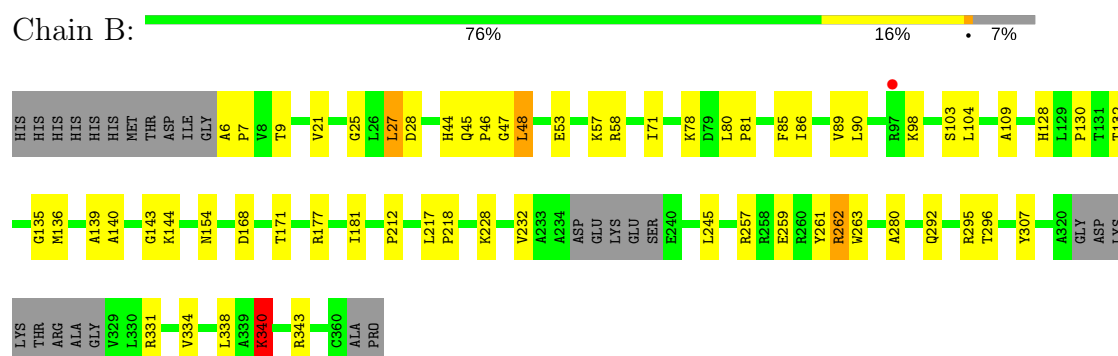
3 Residue-property plots [i](#)

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: 3-dehydroquinase synthase



- Molecule 1: 3-dehydroquinase synthase



4 Data and refinement statistics

Property	Value	Source
Space group	P 43	Depositor
Cell constants a, b, c, α , β , γ	142.83Å 142.83Å 37.17Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.00 – 2.47 28.70 – 2.47	Depositor EDS
% Data completeness (in resolution range)	99.8 (30.00-2.47) 99.8 (28.70-2.47)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	0.09	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.56 (at 2.48Å)	Xtriage
Refinement program	REFMAC 5.5.0109	Depositor
R, R_{free}	0.196 , 0.258 0.196 , 0.251	Depositor DCC
R_{free} test set	1389 reflections (5.30%)	DCC
Wilson B-factor (Å ²)	34.8	Xtriage
Anisotropy	0.166	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 37.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.031 for h,-k,-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	5260	wwPDB-VP
Average B, all atoms (Å ²)	32.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 72.50 % of the origin peak, indicating pseudo translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo translational symmetry is equal to 2.1765e-06. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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

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	1.11	3/2589 (0.1%)	0.92	7/3528 (0.2%)
1	B	1.08	1/2580 (0.0%)	0.89	2/3516 (0.1%)
All	All	1.09	4/5169 (0.1%)	0.91	9/7044 (0.1%)

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	307	TYR	CD2-CE2	-5.48	1.31	1.39
1	A	307	TYR	CD1-CE1	-5.36	1.31	1.39
1	A	14	VAL	CB-CG2	-5.31	1.41	1.52
1	B	307	TYR	CD2-CE2	-5.20	1.31	1.39

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	297	ILE	CG1-CB-CG2	-5.90	98.42	111.40
1	B	340	LYS	CD-CE-NZ	-5.76	98.46	111.70
1	A	308	ASP	CB-CG-OD1	5.69	123.42	118.30
1	A	308	ASP	CB-CG-OD2	-5.68	113.19	118.30
1	A	122	ARG	NE-CZ-NH2	-5.48	117.56	120.30
1	A	295	ARG	NE-CZ-NH1	5.34	122.97	120.30
1	B	143	GLY	N-CA-C	5.34	126.44	113.10
1	A	97	ARG	NE-CZ-NH2	-5.11	117.75	120.30
1	A	32	ASP	CB-CG-OD2	-5.04	113.77	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	A	2546	0	2623	43	0
1	B	2537	0	2615	39	0
2	A	44	0	26	0	0
2	B	44	0	26	0	0
3	A	53	0	0	2	0
3	B	36	0	0	0	0
All	All	5260	0	5290	79	0

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 8.

All (79) 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)
1:A:58:ARG:HH11	1:A:58:ARG:HG2	1.29	0.97
1:A:97:ARG:HH12	1:B:144:LYS:HE2	1.32	0.93
1:A:97:ARG:NH2	1:A:122:ARG:HH11	1.65	0.93
1:A:97:ARG:HH22	1:A:122:ARG:HH11	1.09	0.90
1:A:58:ARG:CG	1:A:58:ARG:HH11	1.85	0.88
1:A:27:LEU:HD13	1:A:58:ARG:NE	1.92	0.85
1:A:97:ARG:HH22	1:A:122:ARG:NH1	1.77	0.83
1:B:48:LEU:HD23	1:B:48:LEU:N	1.94	0.82
1:B:338:LEU:O	1:B:340:LYS:HD2	1.80	0.80
1:A:97:ARG:NH2	1:A:122:ARG:NH1	2.32	0.77
1:A:97:ARG:NH1	1:B:144:LYS:HE2	2.03	0.74
1:B:212:PRO:HA	1:B:217:LEU:HD23	1.72	0.71
1:A:264:ARG:HD2	3:A:415:HOH:O	1.92	0.69
1:A:58:ARG:NH1	1:A:58:ARG:HG2	2.09	0.67
1:B:9:THR:HG22	1:B:21:VAL:HG13	1.78	0.66
1:A:103:SER:O	1:A:129:LEU:HB2	1.95	0.66
1:B:132:THR:O	1:B:136:MET:HG3	1.96	0.65
1:B:28:ASP:OD1	1:B:58:ARG:NH2	2.27	0.65
1:A:58:ARG:CG	1:A:58:ARG:NH1	2.57	0.62
1:B:257:ARG:HD2	1:B:261:TYR:OH	1.99	0.62
1:B:47:GLY:C	1:B:48:LEU:HD23	2.20	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:27:LEU:CD1	1:A:58:ARG:NE	2.62	0.61
1:A:240:GLU:HA	1:A:240:GLU:OE1	1.99	0.61
1:B:48:LEU:CD2	1:B:48:LEU:N	2.65	0.59
1:A:121:LEU:O	1:A:122:ARG:HB2	2.03	0.58
1:B:44:HIS:HD2	1:B:45:GLN:O	1.84	0.58
1:A:129:LEU:HD12	1:A:166:LEU:HB2	1.85	0.57
1:A:259:GLU:O	1:A:262:ARG:HG2	2.05	0.56
1:A:80:LEU:HB3	1:A:81:PRO:HD3	1.88	0.56
1:B:48:LEU:HD12	1:B:104:LEU:HG	1.89	0.54
1:A:53:GLU:O	1:A:57:LYS:HB2	2.08	0.53
1:B:331:ARG:NH2	1:B:343:ARG:HD3	2.24	0.53
1:B:86:ILE:O	1:B:90:LEU:HG	2.08	0.53
1:B:262:ARG:HH11	1:B:262:ARG:HG3	1.73	0.53
1:A:27:LEU:HD13	1:A:58:ARG:HE	1.71	0.52
1:B:280:ALA:HB2	1:B:334:VAL:HB	1.92	0.52
1:A:48:LEU:HD13	1:A:104:LEU:HG	1.93	0.51
1:A:58:ARG:NH1	1:A:58:ARG:HB3	2.25	0.51
1:A:44:HIS:HD2	1:A:45:GLN:O	1.94	0.50
1:A:151:ALA:O	1:A:155:LEU:HD12	2.11	0.50
1:A:58:ARG:CB	1:A:58:ARG:HH11	2.24	0.49
1:B:177:ARG:HG2	1:B:181:ILE:HD12	1.94	0.49
1:A:122:ARG:NH2	1:B:154:ASN:O	2.25	0.48
1:B:85:PHE:O	1:B:89:VAL:HG23	2.13	0.48
1:B:53:GLU:O	1:B:57:LYS:HG2	2.13	0.48
1:B:331:ARG:HH22	1:B:343:ARG:NH1	2.10	0.47
1:A:102:VAL:CG1	1:A:129:LEU:HD22	2.44	0.47
1:B:78:LYS:NZ	1:B:154:ASN:HD21	2.12	0.47
1:B:259:GLU:HG3	1:B:263:TRP:CD1	2.49	0.47
1:A:31:GLU:OE1	1:A:58:ARG:NH1	2.44	0.47
1:B:135:GLY:HA2	1:B:139:ALA:HB3	1.96	0.47
1:A:267:ALA:HB1	1:A:304:PRO:HG3	1.97	0.46
1:A:103:SER:OG	1:A:128:HIS:HA	2.16	0.46
1:B:46:PRO:HD3	1:B:71:ILE:O	2.17	0.44
1:A:203:GLU:OE1	1:A:293:ARG:NH2	2.45	0.44
1:A:259:GLU:O	1:A:260:ARG:C	2.54	0.44
1:B:292:GLN:HG3	1:B:295:ARG:HH21	1.83	0.44
1:A:135:GLY:HA2	1:A:139:ALA:HB3	2.00	0.43
1:A:58:ARG:CB	1:A:58:ARG:NH1	2.82	0.43
1:B:6:ALA:N	1:B:7:PRO:CD	2.81	0.43
1:B:27:LEU:HD13	1:B:58:ARG:HD3	2.01	0.42
1:B:130:PRO:HG3	1:B:140:ALA:CB	2.49	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:180:MET:CE	1:A:212:PRO:HG3	2.49	0.42
1:A:27:LEU:CD1	1:A:58:ARG:HE	2.30	0.42
1:A:267:ALA:HB1	1:A:304:PRO:CG	2.50	0.42
1:A:308:ASP:HA	1:A:309:PRO:HD3	1.91	0.42
1:A:262:ARG:HE	1:A:262:ARG:HB3	1.45	0.42
1:B:45:GLN:HG3	1:B:109:ALA:HB2	2.03	0.41
1:B:80:LEU:N	1:B:81:PRO:HD2	2.35	0.41
1:B:217:LEU:HB3	1:B:218:PRO:HD3	2.01	0.41
1:A:59:LEU:O	1:A:64:VAL:HB	2.21	0.41
1:B:98:LYS:HE2	1:B:98:LYS:HB3	1.94	0.41
1:B:103:SER:OG	1:B:128:HIS:HA	2.21	0.41
1:B:228:LYS:HE3	1:B:245:LEU:HD13	2.02	0.41
1:A:182:CYS:HB2	3:A:415:HOH:O	2.21	0.41
1:A:79:ASP:O	1:A:80:LEU:C	2.59	0.40
1:B:331:ARG:HH22	1:B:343:ARG:HH11	1.68	0.40
1:B:168:ASP:O	1:B:171:THR:OG1	2.32	0.40
1:B:228:LYS:O	1:B:232:VAL:HG13	2.21	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	338/368 (92%)	326 (96%)	11 (3%)	1 (0%)	44	64
1	B	336/368 (91%)	317 (94%)	18 (5%)	1 (0%)	44	64
All	All	674/736 (92%)	643 (95%)	29 (4%)	2 (0%)	44	64

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	25	GLY

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Mol	Chain	Res	Type
1	A	25	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	
1	A	259/280 (92%)	248 (96%)	11 (4%)	34	58
1	B	259/280 (92%)	254 (98%)	5 (2%)	62	83
All	All	518/560 (92%)	502 (97%)	16 (3%)	45	71

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

Mol	Chain	Res	Type
1	A	27	LEU
1	A	58	ARG
1	A	97	ARG
1	A	122	ARG
1	A	129	LEU
1	A	166	LEU
1	A	240	GLU
1	A	257	ARG
1	A	262	ARG
1	A	305	VAL
1	A	340	LYS
1	B	27	LEU
1	B	48	LEU
1	B	262	ARG
1	B	296	THR
1	B	340	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	44	HIS

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Mol	Chain	Res	Type
1	A	154	ASN
1	B	44	HIS
1	B	154	ASN

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](#)

2 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	400	-	41,48,48	0.96	2 (4%)	43,73,73	1.72	7 (16%)
2	NAD	B	400	-	41,48,48	0.89	2 (4%)	43,73,73	1.83	7 (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	400	-	-	0/22/62/62	0/5/5/5

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	NAD	B	400	-	-	0/22/62/62	0/5/5/5

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	400	NAD	C3N-C7N	2.06	1.53	1.50
2	A	400	NAD	C2A-N3A	2.10	1.35	1.32
2	B	400	NAD	O4B-C1B	2.24	1.44	1.41
2	A	400	NAD	O4B-C1B	2.86	1.45	1.41

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	400	NAD	N3A-C2A-N1A	-7.29	122.51	128.86
2	B	400	NAD	N3A-C2A-N1A	-7.26	122.54	128.86
2	B	400	NAD	O7N-C7N-N7N	-3.82	117.14	122.58
2	A	400	NAD	O7N-C7N-C3N	-3.01	116.10	119.62
2	A	400	NAD	O3B-C3B-C4B	-2.98	102.39	111.09
2	A	400	NAD	C5N-C4N-C3N	-2.75	117.11	120.35
2	A	400	NAD	C4D-O4D-C1D	-2.35	107.27	109.77
2	B	400	NAD	O3B-C3B-C4B	-2.24	104.54	111.09
2	B	400	NAD	C3N-C2N-N1N	2.22	122.66	120.43
2	A	400	NAD	O2N-PN-O1N	2.46	125.00	112.28
2	B	400	NAD	N6A-C6A-N1A	2.46	123.64	118.77
2	A	400	NAD	C3N-C7N-N7N	2.76	120.93	117.77
2	B	400	NAD	O2A-PA-O1A	3.02	127.90	112.28
2	B	400	NAD	C3N-C7N-N7N	4.52	122.93	117.77

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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 [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	A	344/368 (93%)	-0.38	4 (1%) 79 80	13, 27, 45, 61	0
1	B	342/368 (92%)	-0.24	1 (0%) 93 94	17, 34, 55, 64	0
All	All	686/736 (93%)	-0.31	5 (0%) 87 88	13, 30, 51, 64	0

All (5) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	97	ARG	3.2
1	A	260	ARG	3.1
1	A	240	GLU	2.5
1	A	177	ARG	2.3
1	A	314	GLN	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](#)

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	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
2	NAD	A	400	44/44	0.96	0.12	-0.34	18,25,32,33	0
2	NAD	B	400	44/44	0.96	0.12	-0.56	22,31,40,45	0

6.5 Other polymers [i](#)

There are no such residues in this entry.