



wwPDB X-ray Structure Validation Summary Report ⓘ

Jan 31, 2016 – 07:44 PM GMT

PDB ID : 1GZU
Title : CRYSTAL STRUCTURE OF HUMAN NICOTINAMIDE MONONUCLEOTIDE ADENYLYLTRANSFERASE IN COMPLEX WITH NMN
Authors : Werner, E.; Ziegler, M.; Lerner, F.; Schweiger, M.; Heinemann, U.
Deposited on : 2002-06-06
Resolution : 2.90 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
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : **NOT EXECUTED**
EDS : **NOT EXECUTED**
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

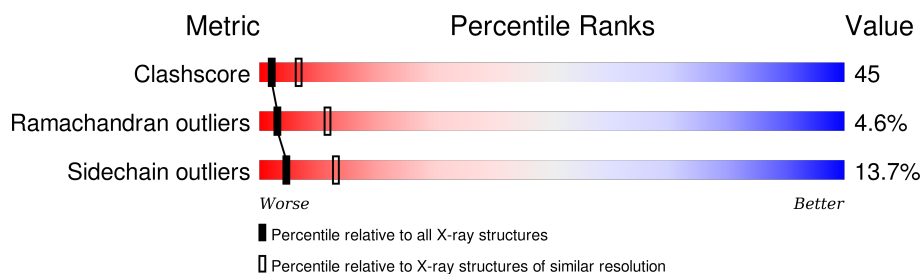
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.90 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))
Clashscore	102246	1668 (2.90-2.90)
Ramachandran outliers	100387	1630 (2.90-2.90)
Sidechain outliers	100360	1632 (2.90-2.90)

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.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	290	
1	B	290	
1	C	290	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	NMN	A	501	X	-	-	-

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Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	NMN	B	501	X	-	-	-
2	NMN	C	501	X	-	-	-

2 Entry composition [i](#)

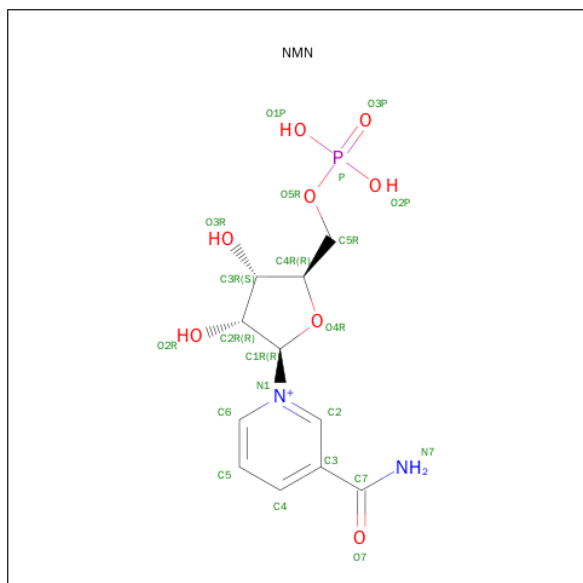
There are 2 unique types of molecules in this entry. The entry contains 5544 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 NICOTINAMIDE MONONUCLEOTIDE ADENYLYL TRANSFERASE.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	228	Total	C	N	O	S	Se	4	0	1
			1826	1168	321	331	3	3			
1	B	228	Total	C	N	O	S	Se	3	0	1
			1826	1168	321	331	3	3			
1	C	228	Total	C	N	O	S	Se	19	0	1
			1826	1168	321	331	3	3			

- Molecule 2 is BETA-NICOTINAMIDE RIBOSE MONOPHOSPHATE (three-letter code: NMN) (formula: $C_{11}H_{16}N_2O_8P$).

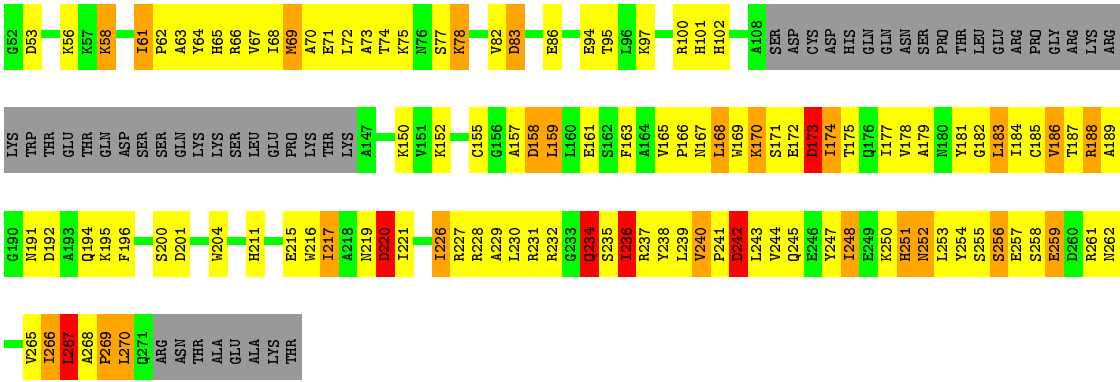


Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			22	11	2	8	1		
2	B	1	Total	C	N	O	P	0	0
			22	11	2	8	1		

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	C	1	Total	C	N	O	P	0	0
			22	11	2	8	1		



4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section will therefore be incomplete.

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	140.80 Å 235.70 Å 89.00 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 2.90	Depositor
% Data completeness (in resolution range)	86.0 (20.00-2.90)	Depositor
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	REFMAC 5.0	Depositor
R, R_{free}	0.246 , 0.286	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	5544	wwPDB-VP
Average B, all atoms (Å ²)	53.0	wwPDB-VP

5 Model quality

5.1 Standard geometry

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

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	0.85	1/1860 (0.1%)	0.84	7/2513 (0.3%)
1	B	0.79	0/1860	0.85	13/2513 (0.5%)
1	C	1.17	5/1860 (0.3%)	1.05	13/2513 (0.5%)
All	All	0.95	6/5580 (0.1%)	0.92	33/7539 (0.4%)

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

The worst 5 of 6 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	78	LYS	CG-CD	25.39	2.38	1.52
1	C	267	LEU	CB-CG	20.50	2.12	1.52
1	C	220	ASP	CB-CG	-18.46	1.12	1.51
1	A	260	ASP	CA-CB	8.63	1.73	1.53
1	C	150	LYS	CG-CD	-7.64	1.26	1.52

The worst 5 of 33 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	78	LYS	CB-CG-CD	-20.62	57.98	111.60
1	C	78	LYS	CG-CD-CE	11.98	147.85	111.90
1	C	220	ASP	CA-CB-CG	10.68	136.90	113.40
1	C	267	LEU	CB-CG-CD2	10.09	128.16	111.00
1	C	267	LEU	CA-CB-CG	-9.59	93.23	115.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	C	220	ASP	Sidechain

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	1826	0	1858	131	0
1	B	1826	0	1858	145	0
1	C	1826	0	1858	234	0
2	A	22	0	9	2	0
2	B	22	0	8	3	0
2	C	22	0	8	4	0
All	All	5544	0	5599	501	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 45.

The worst 5 of 501 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:C:250:LYS:O	1:C:251:HIS:ND1	1.60	1.33
1:C:18:ASN:HB3	1:C:19:PRO:CD	1.57	1.32
1:C:27:LEU:HD22	1:C:186:VAL:CG2	1.69	1.21
1:C:168:LEU:HD23	1:C:169:TRP:CD1	1.75	1.20
1:B:58:LYS:HE2	1:B:58:LYS:N	1.59	1.17

There are no symmetry-related clashes.

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	
1	A	224/290 (77%)	189 (84%)	25 (11%)	10 (4%)	3	12
1	B	224/290 (77%)	188 (84%)	26 (12%)	10 (4%)	3	12
1	C	224/290 (77%)	184 (82%)	29 (13%)	11 (5%)	3	10
All	All	672/870 (77%)	561 (84%)	80 (12%)	31 (5%)	3	11

5 of 31 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	18	ASN
1	A	251	HIS
1	A	265	VAL
1	B	18	ASN
1	B	251	HIS

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	
1	A	199/253 (79%)	178 (89%)	21 (11%)	8	25
1	B	199/253 (79%)	168 (84%)	31 (16%)	3	10
1	C	199/253 (79%)	169 (85%)	30 (15%)	3	11
All	All	597/759 (79%)	515 (86%)	82 (14%)	4	13

5 of 82 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	187	THR
1	B	205	LYS
1	C	240	VAL
1	B	188	ARG
1	B	197	ILE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 18 such sidechains are listed below:

Mol	Chain	Res	Type
1	B	103	GLN
1	B	176	GLN
1	B	245	GLN
1	B	76	ASN
1	B	102	HIS

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](#)

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	NMN	A	501	-	20,23,23	3.75	5 (25%)	26,34,34	3.51	11 (42%)
2	NMN	B	501	-	20,23,23	3.87	5 (25%)	26,34,34	4.54	10 (38%)
2	NMN	C	501	-	20,23,23	3.66	5 (25%)	26,34,34	3.29	8 (30%)

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	NMN	A	501	-	2/2/6/6	0/10/30/30	0/2/2/2
2	NMN	B	501	-	1/1/6/6	0/10/30/30	0/2/2/2
2	NMN	C	501	-	1/1/6/6	0/10/30/30	0/2/2/2

The worst 5 of 15 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	501	NMN	O3R-C3R	-9.30	1.20	1.43
2	A	501	NMN	O3R-C3R	-9.10	1.21	1.43
2	C	501	NMN	O2R-C2R	-8.66	1.22	1.43
2	C	501	NMN	O3R-C3R	-8.37	1.22	1.43
2	B	501	NMN	O2R-C2R	-8.35	1.23	1.43

The worst 5 of 29 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	501	NMN	C4R-O4R-C1R	-13.76	94.59	109.72
2	C	501	NMN	C4R-O4R-C1R	-11.47	97.11	109.72
2	A	501	NMN	C4R-O4R-C1R	-9.05	99.78	109.72
2	A	501	NMN	C5-C4-C3	-2.52	117.17	120.33
2	B	501	NMN	C3-C7-N7	-2.44	115.15	117.82

All (4) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
2	A	501	NMN	C2R
2	A	501	NMN	C4R
2	B	501	NMN	C2R
2	C	501	NMN	C2R

There are no torsion outliers.

There are no ring outliers.

3 monomers are involved in 9 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	501	NMN	2	0
2	B	501	NMN	3	0
2	C	501	NMN	4	0

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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

EDS was not executed - this section will therefore be empty.

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

EDS was not executed - this section will therefore be empty.

6.3 Carbohydrates [i](#)

EDS was not executed - this section will therefore be empty.

6.4 Ligands [i](#)

EDS was not executed - this section will therefore be empty.

6.5 Other polymers [i](#)

EDS was not executed - this section will therefore be empty.