



Full wwPDB X-ray Structure Validation Report ⓘ

May 21, 2020 – 11:56 pm BST

PDB ID : 1ENP
Title : BRASSICA NAPUS ENOYL ACP REDUCTASE/NADH BINARY COMPLEX AT PH 8.0 AND ROOM TEMPERATURE
Authors : Rafferty, J.B.; Rice, D.W.
Deposited on : 1995-10-18
Resolution : 2.60 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

<https://www.wwpdb.org/validation/2017/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 : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : **NOT EXECUTED**
EDS : **NOT EXECUTED**
buster-report : 1.1.7 (2018)
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.11

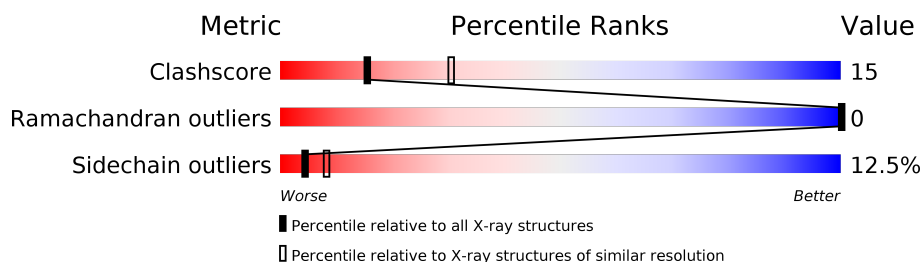
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.60 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	141614	3518 (2.60-2.60)
Ramachandran outliers	138981	3455 (2.60-2.60)
Sidechain outliers	138945	3455 (2.60-2.60)

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 respectively. 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 $\leq 5\%$.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	312	

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 2279 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.

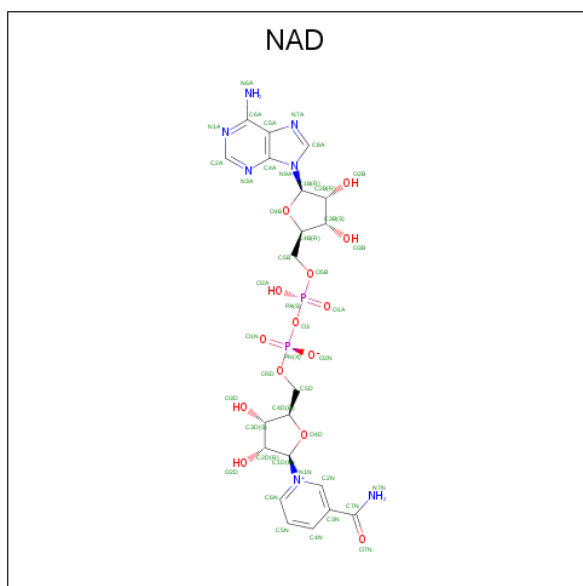
- Molecule 1 is a protein called ENOYL ACYL CARRIER PROTEIN REDUCTASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	297	2181	1386	369	420	6	0	0	0

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	311	ASP	ASN	CONFLICT	UNP P80030

- Molecule 2 is NICOTINAMIDE-ADENINE-DINUCLEOTIDE (three-letter code: NAD) (formula: $C_{21}H_{27}N_7O_{14}P_2$).



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

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	54	Total	O	0	0
			54	54		

4 Data and refinement statistics

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

Property	Value	Source
Space group	P 4 ₂ 2 ₁ 2	Depositor
Cell constants a, b, c, α , β , γ	70.50 Å 70.50 Å 117.80 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	10.00 – 2.60	Depositor
% Data completeness (in resolution range)	(Not available) (10.00-2.60)	Depositor
R_{merge}	0.04	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	TNT	Depositor
R, R_{free}	(Not available) , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	2279	wwPDB-VP
Average B, all atoms (Å ²)	29.0	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: 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	0.98	7/2224 (0.3%)	1.27	28/3028 (0.9%)

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	A	0	2

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	52	TRP	C-N	10.12	1.57	1.34
1	A	77	GLY	C-N	9.49	1.55	1.34
1	A	284	ILE	C-N	7.59	1.51	1.34
1	A	286	GLY	C-O	6.88	1.34	1.23
1	A	285	THR	C-O	6.78	1.36	1.23
1	A	182	ALA	C-N	-5.97	1.20	1.34
1	A	298	MET	C-N	5.54	1.43	1.33

All (28) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	65	ARG	NE-CZ-NH1	9.38	124.99	120.30
1	A	91	VAL	CA-CB-CG2	8.77	124.06	110.90
1	A	52	TRP	C-N-CA	-7.69	102.48	121.70
1	A	77	GLY	C-N-CA	-7.48	103.01	121.70
1	A	53	VAL	O-C-N	-7.46	106.94	121.10
1	A	29	ASP	CB-CG-OD1	7.20	124.78	118.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	29	ASP	CB-CG-OD2	-6.94	112.05	118.30
1	A	212	ASP	CB-CG-OD1	-6.94	112.06	118.30
1	A	76	ASP	CB-CG-OD1	-6.76	112.22	118.30
1	A	267	ASP	CB-CG-OD2	-6.69	112.28	118.30
1	A	76	ASP	CB-CG-OD2	6.47	124.13	118.30
1	A	93	ASP	CB-CG-OD1	-6.30	112.63	118.30
1	A	287	ALA	O-C-N	6.13	132.51	122.70
1	A	131	ASP	CB-CG-OD1	-6.02	112.88	118.30
1	A	80	MET	CG-SD-CE	6.00	109.80	100.20
1	A	126	ASP	CB-CG-OD1	-5.83	113.06	118.30
1	A	65	ARG	NE-CZ-NH2	-5.77	117.42	120.30
1	A	267	ASP	CB-CG-OD1	5.77	123.49	118.30
1	A	28	ASP	CB-CG-OD1	-5.68	113.19	118.30
1	A	93	ASP	CB-CG-OD2	5.58	123.32	118.30
1	A	91	VAL	CG1-CB-CG2	-5.54	102.03	110.90
1	A	227	ARG	NE-CZ-NH2	-5.40	117.60	120.30
1	A	101	ASP	CB-CG-OD1	-5.37	113.47	118.30
1	A	131	ASP	CB-CG-OD2	5.32	123.09	118.30
1	A	222	ARG	NE-CZ-NH1	5.19	122.90	120.30
1	A	287	ALA	CA-C-N	-5.19	105.79	117.20
1	A	28	ASP	CB-CG-OD2	5.08	122.87	118.30
1	A	64	ARG	NE-CZ-NH1	5.07	122.83	120.30

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	47	ILE	Mainchain
1	A	53	VAL	Mainchain

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	2181	0	2157	65	0
2	A	44	0	25	1	0
3	A	54	0	0	3	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	2279	0	2182	65	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 15.

All (65) 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:115:THR:OG1	1:A:118:GLU:HG3	1.89	0.72
1:A:159:ILE:HD13	1:A:205:ALA:HB2	1.70	0.72
1:A:35:ALA:HB2	1:A:266:ALA:HB1	1.70	0.71
1:A:202:MET:HE3	1:A:202:MET:HA	1.72	0.71
1:A:32:TYR:CZ	1:A:236:LEU:HB2	2.29	0.66
1:A:20:ARG:HB3	1:A:46:GLU:HB3	1.78	0.65
1:A:84:LYS:HD3	1:A:86:TYR:CZ	2.34	0.62
1:A:53:VAL:N	1:A:54:PRO:HD2	2.16	0.60
1:A:62:SER:HB3	1:A:67:LYS:HE3	1.84	0.59
1:A:179:PRO:HB3	1:A:224:GLN:O	2.03	0.58
1:A:142:GLU:HB3	1:A:145:LYS:HG3	1.87	0.56
1:A:202:MET:HE3	1:A:202:MET:CA	2.36	0.56
1:A:20:ARG:CB	1:A:46:GLU:HB3	2.36	0.55
1:A:94:ASN:O	1:A:97:ASP:HB2	2.07	0.55
1:A:287:ALA:HA	3:A:550:HOH:O	2.07	0.54
1:A:29:ASP:HA	1:A:34:TRP:CG	2.43	0.54
1:A:86:TYR:HH	1:A:127:PHE:HE2	1.56	0.54
1:A:13:PRO:HD2	3:A:502:HOH:O	2.07	0.53
1:A:174:LEU:HD11	1:A:226:ILE:HG13	1.89	0.53
1:A:56:LEU:HD22	1:A:87:PRO:HB3	1.90	0.52
1:A:184:ILE:HD11	1:A:231:ILE:HG13	1.93	0.51
1:A:189:ILE:CG1	1:A:189:ILE:O	2.59	0.51
1:A:86:TYR:OH	1:A:127:PHE:HE2	1.95	0.50
1:A:248:ASP:HB2	3:A:540:HOH:O	2.10	0.50
1:A:260:ILE:CG2	1:A:262:LYS:HD3	2.42	0.50
1:A:243:ALA:HA	1:A:248:ASP:HB3	1.94	0.50
1:A:94:ASN:HB3	1:A:117:GLN:NE2	2.26	0.49
1:A:20:ARG:HD2	1:A:129:SER:O	2.11	0.49
1:A:215:VAL:O	1:A:218:PHE:HB3	2.13	0.49
1:A:265:THR:O	1:A:268:GLU:HB2	2.13	0.49
1:A:189:ILE:O	1:A:189:ILE:HG13	2.14	0.48
1:A:76:ASP:CG	1:A:78:SER:H	2.17	0.47
1:A:242:LYS:O	1:A:244:ILE:HG12	2.14	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:35:ALA:CB	1:A:266:ALA:HB1	2.40	0.47
1:A:184:ILE:HA	1:A:229:ASN:O	2.15	0.47
1:A:140:GLY:O	1:A:143:VAL:HB	2.15	0.46
1:A:26:ILE:CD1	1:A:34:TRP:HA	2.46	0.46
1:A:295:LEU:HA	1:A:295:LEU:HD23	1.77	0.46
1:A:189:ILE:C	1:A:191:SER:H	2.20	0.45
1:A:202:MET:CE	1:A:202:MET:HA	2.44	0.45
1:A:26:ILE:HD11	1:A:33:GLY:C	2.37	0.45
1:A:133:LEU:O	1:A:183:SER:HA	2.17	0.45
1:A:174:LEU:N	1:A:175:PRO:HD2	2.33	0.44
1:A:99:PRO:HG2	1:A:102:VAL:CG2	2.47	0.44
1:A:140:GLY:HA2	1:A:141:PRO:HD2	1.77	0.44
1:A:53:VAL:N	1:A:54:PRO:CD	2.80	0.44
1:A:12:LEU:HD23	1:A:12:LEU:HA	1.82	0.44
1:A:240:ALA:HB2	1:A:248:ASP:OD2	2.17	0.44
1:A:306:VAL:HG23	1:A:307:PHE:CD1	2.53	0.44
1:A:202:MET:HB3	1:A:202:MET:HE2	1.87	0.43
1:A:32:TYR:OH	1:A:236:LEU:HB2	2.18	0.43
1:A:115:THR:HG1	1:A:118:GLU:HG3	1.84	0.43
1:A:255:TYR:O	1:A:261:GLN:HG3	2.18	0.42
1:A:105:ASN:O	1:A:106:LYS:C	2.58	0.42
1:A:50:GLY:HA2	1:A:86:TYR:O	2.19	0.42
1:A:229:ASN:OD1	1:A:287:ALA:HB3	2.20	0.42
1:A:235:PRO:HG3	1:A:250:MET:HB3	2.00	0.42
1:A:29:ASP:HA	1:A:34:TRP:CD2	2.55	0.42
1:A:159:ILE:HD13	1:A:205:ALA:CB	2.47	0.41
1:A:210:GLU:O	1:A:213:THR:HB	2.21	0.41
1:A:295:LEU:O	1:A:296:ASN:C	2.57	0.41
1:A:108:TYR:HB3	1:A:114:TRP:CZ2	2.55	0.41
1:A:303:ASP:O	1:A:305:PRO:HD3	2.20	0.41
1:A:15:ASP:OD1	1:A:17:ARG:HB2	2.21	0.40
1:A:32:TYR:HE2	2:A:501:NAD:C3N	2.34	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	295/312 (95%)	271 (92%)	24 (8%)	0	100	100

There are no Ramachandran outliers to report.

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	224/245 (91%)	196 (88%)	28 (12%)	4	8

All (28) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	12	LEU
1	A	38	LYS
1	A	53	VAL
1	A	62	SER
1	A	67	LYS
1	A	80	MET
1	A	83	LYS
1	A	94	ASN
1	A	96	GLU
1	A	101	ASP
1	A	111	SER
1	A	124	ARG
1	A	137	LEU
1	A	144	SER
1	A	145	LYS
1	A	151	SER
1	A	153	LYS
1	A	162	SER
1	A	170	LEU

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Mol	Chain	Res	Type
1	A	204	SER
1	A	239	ARG
1	A	242	LYS
1	A	244	ILE
1	A	247	ILE
1	A	260	ILE
1	A	263	THR
1	A	278	SER
1	A	295	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	117	GLN
1	A	256	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](#)

1 ligand is 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	-	42,48,48	2.38	15 (35%)	50,73,73	2.59	16 (32%)

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	-	-	7/26/62/62	0/5/5/5

All (15) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	501	NAD	C3N-C7N	6.86	1.60	1.50
2	A	501	NAD	C2A-N3A	6.49	1.42	1.32
2	A	501	NAD	C4A-N3A	5.33	1.43	1.35
2	A	501	NAD	C7N-N7N	3.96	1.40	1.33
2	A	501	NAD	C2N-N1N	3.49	1.39	1.35
2	A	501	NAD	O4B-C1B	-2.79	1.37	1.41
2	A	501	NAD	O4D-C4D	2.75	1.51	1.45
2	A	501	NAD	O3D-C3D	-2.71	1.36	1.43
2	A	501	NAD	C3D-C4D	2.67	1.59	1.53
2	A	501	NAD	C2B-C1B	2.38	1.57	1.53
2	A	501	NAD	C6A-C5A	2.35	1.52	1.43
2	A	501	NAD	C5N-C4N	2.32	1.43	1.38
2	A	501	NAD	C2N-C3N	-2.28	1.35	1.39
2	A	501	NAD	O2B-C2B	-2.17	1.37	1.43
2	A	501	NAD	PN-O1N	-2.05	1.43	1.50

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	501	NAD	C2N-C3N-C4N	8.14	127.49	118.26
2	A	501	NAD	C3N-C2N-N1N	-7.08	113.51	120.43
2	A	501	NAD	C5N-C4N-C3N	-6.33	112.86	120.34
2	A	501	NAD	C6N-N1N-C2N	4.60	126.17	121.97
2	A	501	NAD	C2A-N1A-C6A	4.35	126.20	118.75
2	A	501	NAD	C5A-C6A-N6A	3.88	126.24	120.35
2	A	501	NAD	N3A-C2A-N1A	-3.87	122.63	128.68
2	A	501	NAD	C2N-C3N-C7N	-3.64	108.91	119.46
2	A	501	NAD	C3D-C2D-C1D	-3.56	95.62	100.98
2	A	501	NAD	C3B-C2B-C1B	-3.53	95.66	100.98

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	501	NAD	C5A-C6A-N1A	-3.45	112.53	120.35
2	A	501	NAD	O3D-C3D-C4D	-2.77	103.04	111.05
2	A	501	NAD	O4D-C4D-C3D	-2.72	99.72	105.11
2	A	501	NAD	C6N-C5N-C4N	2.29	122.77	119.44
2	A	501	NAD	C5N-C6N-N1N	-2.24	117.18	120.40
2	A	501	NAD	O2N-PN-O1N	2.02	122.22	112.24

There are no chirality outliers.

All (7) torsion outliers are listed below:

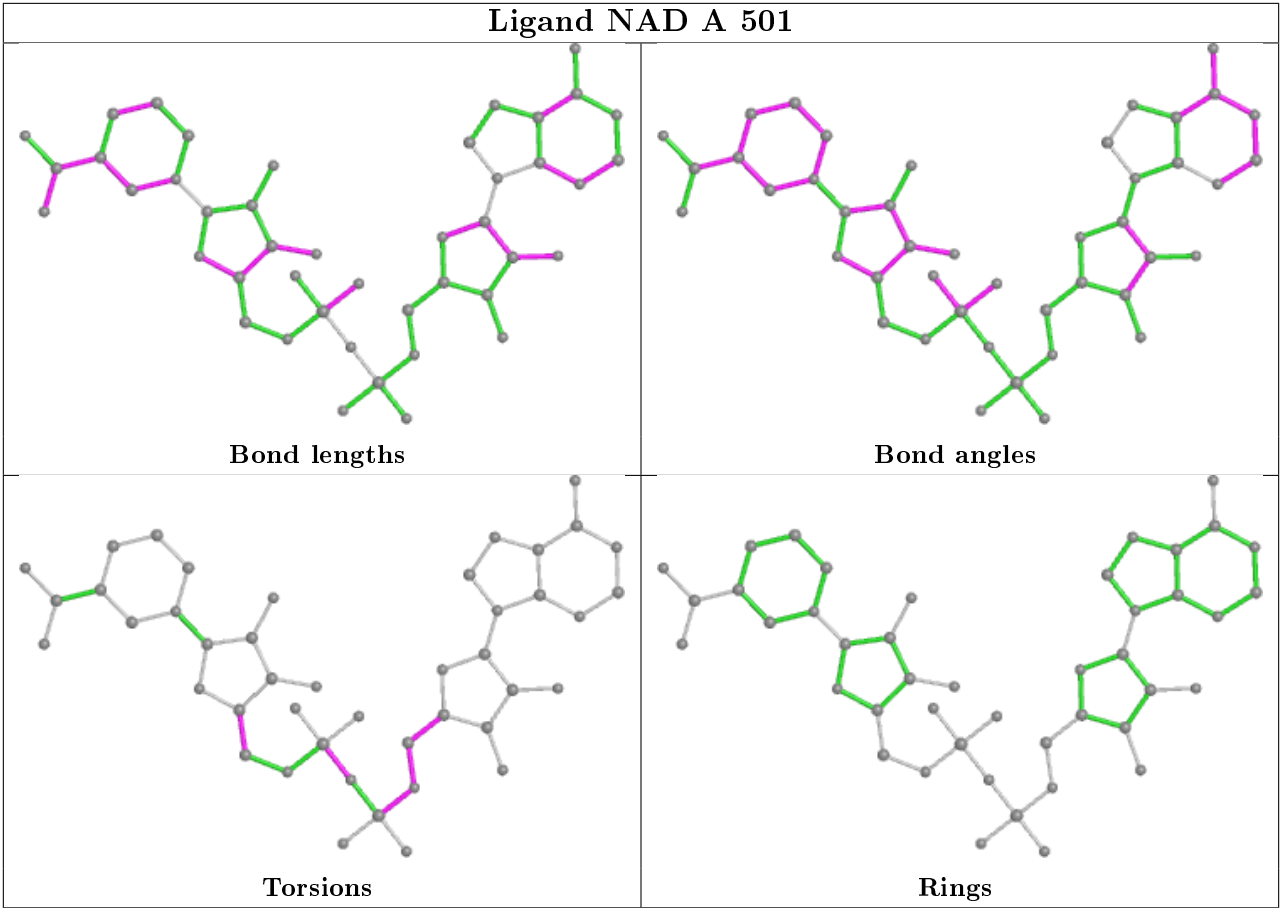
Mol	Chain	Res	Type	Atoms
2	A	501	NAD	O4B-C4B-C5B-O5B
2	A	501	NAD	O4D-C4D-C5D-O5D
2	A	501	NAD	C3D-C4D-C5D-O5D
2	A	501	NAD	C3B-C4B-C5B-O5B
2	A	501	NAD	C4B-C5B-O5B-PA
2	A	501	NAD	PA-O3-PN-O1N
2	A	501	NAD	C5B-O5B-PA-O1A

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	501	NAD	1	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.



5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

The following chains have linkage breaks:

Mol	Chain	Number of breaks
1	A	1

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	A	182:ALA	C	183:SER	N	1.20

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

EDS was not executed - this section is therefore empty.

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

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates ⓘ

EDS was not executed - this section is therefore empty.

6.4 Ligands ⓘ

EDS was not executed - this section is therefore empty.

6.5 Other polymers ⓘ

EDS was not executed - this section is therefore empty.