



# Full wwPDB X-ray Structure Validation Report

Feb 27, 2014 – 01:35 PM GMT

PDB ID : 1DFH  
Title : X-RAY STRUCTURE OF ESCHERICHIA COLI ENOYL REDUCTASE  
WITH BOUND NAD AND THIENO-DIAZABORINE  
Authors : Baldock, C.; Rafferty, J.B.; Rice, D.W.  
Deposited on : 1997-01-16  
Resolution : 2.20 Å(reported)

This is a full wwPDB validation report for a publicly released PDB entry.  
We welcome your comments at [validation@mail.wwpdb.org](mailto:validation@mail.wwpdb.org)  
A user guide is available at <http://wwpdb.org/ValidationPDFNotes.html>

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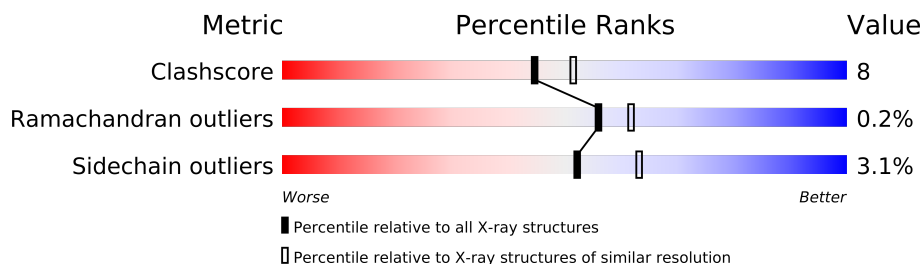
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.15 2013  
Xtriage (Phenix) : **NOT EXECUTED**  
EDS : **NOT EXECUTED**  
Percentile statistics : 21963  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)  
Validation Pipeline (wwPDB-VP) : stable22683

# 1 Overall quality at a glance

The reported resolution of this entry is 2.20 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	79885	3751 (2.20-2.20)
Ramachandran outliers	78287	3681 (2.20-2.20)
Sidechain outliers	78261	3682 (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  $\geq 3$ , 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.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	261	
1	B	261	

The following table lists non-polymeric compounds that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Geometry	Electron density
3	TDB	A	0	X	-

## 2 Entry composition i

There are 3 unique types of molecules in this entry. The entry contains 3936 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 ENOYL ACYL CARRIER PROTEIN REDUCTASE.

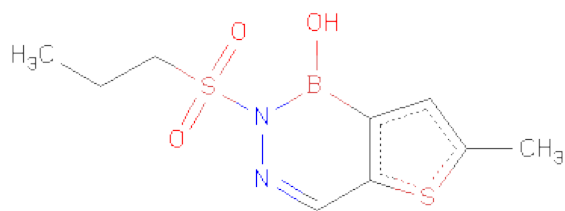
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	257	Total	C	N	O	S	0	0	0
			1904	1200	327	364	13			
1	B	257	Total	C	N	O	S	0	0	0
			1910	1203	330	364	13			

- 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
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 6-METHYL-2(PROPANE-1-SULFONYL)-2H-THIENO[3,2-D][1,2,3]DIAZABORININ-1-OL (three-letter code: TDB) (formula:  $C_9H_{13}BN_2O_3S_2$ ).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
			Total	B	C	N	O	S		
3	A	1	17	1	9	2	3	2	0	0
3	B	1	17	1	9	2	3	2	0	0

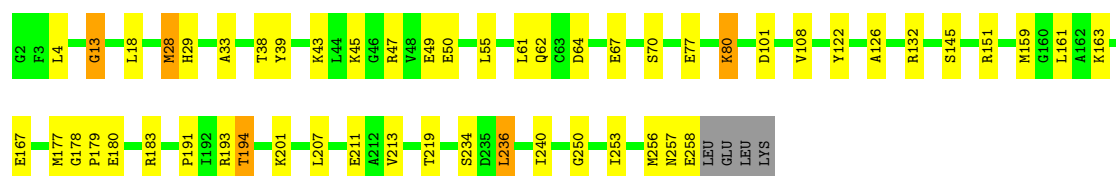
### 3 Residue-property plots

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.

Note EDS was not executed.

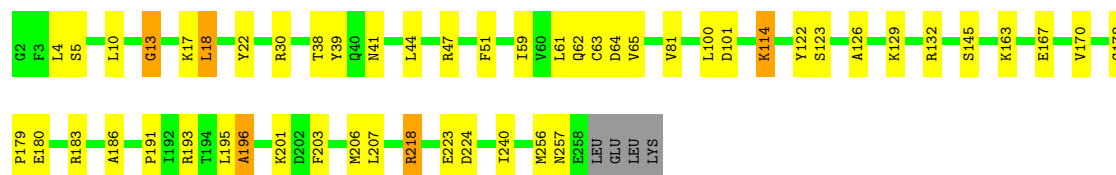
#### • Molecule 1: ENOYL ACYL CARRIER PROTEIN REDUCTASE

Chain A: 



#### • Molecule 1: ENOYL ACYL CARRIER PROTEIN REDUCTASE

Chain B: 



## 4 Data and refinement statistics

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

Property	Value	Source
Space group	P 61 2 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	80.90Å 80.90Å 328.30Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	10.00 – 2.20	Depositor
% Data completeness (in resolution range)	93.0 (10.00-2.20)	Depositor
$R_{merge}$	0.06	Depositor
$R_{sym}$	(Not available)	Depositor
Refinement program	TNT	Depositor
R, $R_{free}$	0.194 , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	3936	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	27.0	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: TDB, 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	0.70	0/1936	0.78	5/2616 (0.2%)
1	B	0.69	0/1942	0.78	2/2623 (0.1%)
All	All	0.70	0/3878	0.78	7/5239 (0.1%)

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	1
1	B	0	2
All	All	0	3

There are no bond length outliers.

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	13	GLY	O-C-N	-11.46	104.36	122.70
1	B	13	GLY	O-C-N	-10.60	105.75	122.70
1	B	256	MET	CG-SD-CE	5.91	109.66	100.20
1	A	159	MET	CG-SD-CE	5.81	109.50	100.20
1	A	177	MET	CG-SD-CE	5.67	109.26	100.20
1	A	151	ARG	NE-CZ-NH2	-5.62	117.49	120.30
1	A	28	MET	CG-SD-CE	5.61	109.18	100.20

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	13	GLY	Mainchain
1	B	13	GLY	Mainchain
1	B	257	ASN	Mainchain

## 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	1904	0	1887	36	0
1	B	1910	0	1898	29	0
2	A	44	0	24	2	0
2	B	44	0	24	2	0
3	A	17	0	13	1	0
3	B	17	0	13	1	0
All	All	3936	0	3859	66	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 8.

All (66) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:218:ARG:NH2	1:B:224:ASP:OD2	2.04	0.90
1:B:59:ILE:HD13	1:B:81:VAL:HG21	1.57	0.86
1:A:258:GLU:O	1:A:258:GLU:HG3	1.82	0.78
1:A:132:ARG:NH2	1:A:180:GLU:OE2	2.14	0.76
1:B:38:THR:HA	1:B:61:LEU:O	1.88	0.72
1:A:193:ARG:HG3	1:A:207:LEU:HD21	1.72	0.72
1:A:108:VAL:O	1:B:129:LYS:HE3	1.89	0.72
1:B:193:ARG:HA	1:B:207:LEU:HD11	1.81	0.62
1:A:193:ARG:HA	1:A:207:LEU:HD21	1.82	0.62
1:A:193:ARG:HG3	1:A:207:LEU:CD2	2.28	0.61
1:A:39:TYR:CZ	1:A:62:GLN:HB2	2.39	0.57
2:A:501:NAD:C2D	3:A:0:TDB:N1	2.67	0.56
1:A:38:THR:HA	1:A:61:LEU:O	2.06	0.56
1:B:10:LEU:HD11	1:B:38:THR:HG23	1.88	0.55
1:B:17:LYS:HG3	1:B:22:TYR:CE2	2.40	0.55

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:122:TYR:CE2	1:A:126:ALA:HB2	2.41	0.55
1:B:39:TYR:CZ	1:B:62:GLN:HB2	2.42	0.54
1:A:193:ARG:CG	1:A:207:LEU:HD21	2.38	0.54
1:B:203:PHE:HA	1:B:206:MET:HE3	1.90	0.54
1:A:77:GLU:O	1:A:80:LYS:HB2	2.08	0.53
1:B:122:TYR:CE2	1:B:126:ALA:HB2	2.44	0.53
2:B:501:NAD:H2D	3:B:0:TDB:N1	2.23	0.53
1:A:45:LYS:HE2	1:A:49:GLU:OE2	2.09	0.52
1:A:101:ASP:OD1	1:A:101:ASP:C	2.48	0.52
1:A:183:ARG:HD2	1:A:240:ILE:O	2.10	0.51
1:A:234:SER:OG	1:A:236:LEU:HB2	2.11	0.51
1:A:18:LEU:HD13	1:A:194:THR:HG22	1.91	0.51
1:A:258:GLU:O	1:A:258:GLU:CG	2.52	0.49
1:B:101:ASP:OD1	1:B:201:LYS:HD3	2.12	0.49
1:B:62:GLN:HG2	1:B:63:CYS:N	2.26	0.49
1:A:47:ARG:O	1:A:50:GLU:HB2	2.13	0.49
1:B:132:ARG:NH2	1:B:180:GLU:OE2	2.43	0.49
1:A:29:HIS:CG	1:A:55:LEU:HD12	2.47	0.49
1:A:145:SER:O	2:A:501:NAD:H6N	2.14	0.48
1:B:59:ILE:HD13	1:B:81:VAL:CG2	2.37	0.47
1:B:145:SER:O	2:B:501:NAD:H6N	2.14	0.47
1:A:64:ASP:OD1	1:A:64:ASP:C	2.52	0.47
1:A:29:HIS:CD2	1:A:55:LEU:O	2.68	0.47
1:B:170:VAL:HG21	1:B:186:ALA:HB2	1.97	0.47
1:B:163:LYS:O	1:B:167:GLU:HG3	2.14	0.47
1:A:250:GLY:O	1:A:253:ILE:HG13	2.15	0.47
1:B:191:PRO:HB2	1:B:207:LEU:HD23	1.98	0.46
1:A:4:LEU:HB3	1:A:33:ALA:HB2	1.97	0.45
1:A:191:PRO:HB2	1:A:207:LEU:HD12	1.97	0.45
1:B:114:LYS:HB2	1:B:114:LYS:HE3	1.57	0.45
1:B:47:ARG:HG3	1:B:51:PHE:CE2	2.53	0.44
1:A:67:GLU:HB2	1:A:70:SER:CB	2.48	0.43
1:B:30:ARG:O	1:B:30:ARG:HG2	2.17	0.43
1:B:17:LYS:HB2	1:B:17:LYS:NZ	2.34	0.43
1:B:178:GLY:N	1:B:179:PRO:CD	2.82	0.43
1:A:43:LYS:O	1:A:43:LYS:HG2	2.18	0.43
1:B:183:ARG:HD2	1:B:240:ILE:O	2.19	0.43
1:B:64:ASP:OD1	1:B:64:ASP:C	2.58	0.42
1:A:256:MET:C	1:A:258:GLU:H	2.22	0.42
1:A:193:ARG:CB	1:A:207:LEU:HD21	2.49	0.42
1:A:178:GLY:N	1:A:179:PRO:CD	2.83	0.42
1:A:4:LEU:HD12	1:A:28:MET:HA	2.02	0.42

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:65:VAL:HB	1:B:123:SER:HB2	2.02	0.41
1:A:257:ASN:HD22	1:A:257:ASN:HA	1.56	0.41
1:B:195:LEU:C	1:B:196:ALA:O	2.56	0.41
1:B:18:LEU:HD12	1:B:18:LEU:HA	1.81	0.41
1:B:41:ASN:OD1	1:B:41:ASN:C	2.59	0.41
1:A:211:GLU:HG2	1:A:219:THR:HG23	2.03	0.40
1:A:163:LYS:O	1:A:167:GLU:HG3	2.21	0.40
1:A:213:VAL:HB	1:A:257:ASN:HB2	2.03	0.40
1:A:161:LEU:N	1:A:161:LEU:HD12	2.37	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	255/261 (98%)	243 (95%)	12 (5%)	0	100	100
1	B	255/261 (98%)	247 (97%)	7 (3%)	1 (0%)	43	45
All	All	510/522 (98%)	490 (96%)	19 (4%)	1 (0%)	56	62

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	196	ALA

### 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.

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
1	A	195/200 (98%)	191 (98%)	4 (2%)	66 78
1	B	196/200 (98%)	188 (96%)	8 (4%)	41 49
All	All	391/400 (98%)	379 (97%)	12 (3%)	52 63

All (12) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	80	LYS
1	A	194	THR
1	A	201	LYS
1	A	236	LEU
1	B	4	LEU
1	B	5	SER
1	B	18	LEU
1	B	44	LEU
1	B	100	LEU
1	B	114	LYS
1	B	218	ARG
1	B	223	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	257	ASN
1	B	40	GLN
1	B	54	GLN

### 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

4 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
3	TDB	A	0	2	18,18,18	5.91	14 (77%)	20,27,27	7.35	16 (80%)
2	NAD	A	501	3	48,48,48	1.65	11 (22%)	73,73,73	2.35	20 (27%)
3	TDB	B	0	2	18,18,18	5.43	11 (61%)	20,27,27	5.22	15 (75%)
2	NAD	B	501	3	48,48,48	1.81	14 (29%)	73,73,73	2.47	22 (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
3	TDB	A	0	2	-	0/4/23/23	0/0/2/2
2	NAD	A	501	3	-	0/30/62/62	0/3/5/5
3	TDB	B	0	2	-	0/4/23/23	0/0/2/2
2	NAD	B	501	3	-	0/30/62/62	0/3/5/5

All (50) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	0	TDB	B1-O1	13.46	1.56	1.36
3	A	0	TDB	B1-O1	13.04	1.55	1.36
3	A	0	TDB	C14-S1	-11.74	1.60	1.72
3	B	0	TDB	C14-S1	-11.18	1.61	1.72
3	A	0	TDB	O15-S15	8.88	1.53	1.43
3	B	0	TDB	N1-N2	-8.10	1.32	1.41

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	0	TDB	N1-N2	-7.84	1.32	1.41
3	A	0	TDB	O16-S15	7.52	1.52	1.43
3	A	0	TDB	C7-S1	-6.25	1.62	1.74
3	B	0	TDB	O16-S15	5.40	1.49	1.43
3	B	0	TDB	C15-S15	5.08	1.83	1.78
3	B	0	TDB	B1-N1	-4.99	1.43	1.47
3	B	0	TDB	C2-N2	4.98	1.34	1.29
3	A	0	TDB	B1-N1	-4.82	1.43	1.47
2	A	501	NAD	C3N-C7N	4.48	1.58	1.50
2	B	501	NAD	O2B-C2B	-4.44	1.32	1.43
3	B	0	TDB	C7-S1	-4.18	1.66	1.74
2	B	501	NAD	C4A-N9A	4.09	1.43	1.37
2	B	501	NAD	C2N-N1N	3.82	1.40	1.35
3	A	0	TDB	B1-C13	-3.64	1.47	1.54
3	B	0	TDB	B1-C13	-3.47	1.48	1.54
3	A	0	TDB	C8-C7	-3.46	1.39	1.50
2	A	501	NAD	O3B-C3B	3.22	1.50	1.43
2	B	501	NAD	O4D-C4D	3.18	1.52	1.45
2	B	501	NAD	C3N-C7N	3.05	1.55	1.50
2	A	501	NAD	O4D-C4D	3.03	1.52	1.45
2	A	501	NAD	PN-O1N	-2.90	1.41	1.48
2	B	501	NAD	PN-O1N	-2.88	1.41	1.48
2	B	501	NAD	O4B-C4B	2.85	1.51	1.45
2	A	501	NAD	C8A-N7A	-2.85	1.29	1.34
2	B	501	NAD	C2A-N3A	2.79	1.37	1.32
3	A	0	TDB	C2-N2	2.76	1.32	1.29
2	A	501	NAD	O2B-C2B	-2.73	1.36	1.43
3	A	0	TDB	C16-C15	-2.56	1.38	1.52
3	A	0	TDB	C14-C2	-2.49	1.39	1.44
2	B	501	NAD	O3B-C3B	2.39	1.48	1.43
3	B	0	TDB	O15-S15	2.39	1.46	1.43
3	A	0	TDB	C15-S15	2.36	1.80	1.78
2	B	501	NAD	C5A-N7A	-2.32	1.31	1.40
2	B	501	NAD	C2A-N1A	-2.31	1.29	1.33
2	A	501	NAD	C2A-N1A	-2.25	1.29	1.33
2	A	501	NAD	O3D-C3D	-2.21	1.37	1.43
3	B	0	TDB	C8-C7	-2.21	1.43	1.50
2	A	501	NAD	C5A-N7A	-2.20	1.32	1.40
2	A	501	NAD	C2A-N3A	2.20	1.36	1.32
2	B	501	NAD	O4B-C1B	-2.16	1.38	1.41
3	A	0	TDB	C12-C7	2.12	1.43	1.37
2	B	501	NAD	O2D-C2D	-2.04	1.38	1.43

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	501	NAD	C1B-N9A	-2.03	1.42	1.48
2	B	501	NAD	C6A-N6A	-2.03	1.28	1.35

All (73) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	0	TDB	O16-S15-O15	26.94	146.31	118.95
3	B	0	TDB	O16-S15-O15	10.66	129.78	118.95
3	B	0	TDB	O16-S15-C15	-10.42	89.55	107.68
2	B	501	NAD	C8A-N9A-C4A	-9.07	99.97	106.90
2	B	501	NAD	C2N-C3N-C4N	8.90	128.39	118.31
2	A	501	NAD	C2N-C3N-C4N	8.75	128.22	118.31
3	B	0	TDB	O1-B1-C13	-7.56	100.52	123.77
2	A	501	NAD	N3A-C2A-N1A	-7.37	122.55	128.71
3	A	0	TDB	O16-S15-C15	-7.15	95.23	107.68
3	B	0	TDB	C16-C15-S15	6.46	134.89	113.72
3	A	0	TDB	O1-B1-C13	-6.39	104.11	123.77
3	B	0	TDB	O15-S15-C15	6.34	118.71	107.68
2	A	501	NAD	C8A-N9A-C4A	-6.27	102.11	106.90
3	B	0	TDB	C8-C7-S1	6.23	126.74	119.88
2	B	501	NAD	C5N-C4N-C3N	-6.12	112.38	120.32
3	A	0	TDB	C13-B1-N1	6.11	117.48	115.51
3	A	0	TDB	O15-S15-C15	-5.86	97.48	107.68
3	A	0	TDB	C16-C15-S15	5.77	132.63	113.72
3	A	0	TDB	C12-C13-C14	5.64	117.52	112.15
2	A	501	NAD	C5N-C4N-C3N	-5.58	113.07	120.32
2	B	501	NAD	N3A-C2A-N1A	-5.57	124.05	128.71
2	B	501	NAD	C3N-C2N-N1N	-5.36	111.32	120.36
3	A	0	TDB	C7-C12-C13	-5.21	103.35	112.06
3	B	0	TDB	C12-C13-C14	5.02	116.93	112.15
3	B	0	TDB	O1-B1-N1	-4.83	117.84	123.97
3	A	0	TDB	O1-B1-N1	-4.79	117.89	123.97
2	A	501	NAD	C4B-O4B-C1B	-4.75	104.59	109.75
2	A	501	NAD	C3N-C2N-N1N	-4.68	112.47	120.36
2	A	501	NAD	C2A-N1A-C6A	4.51	126.91	118.77
3	A	0	TDB	C7-S1-C14	4.29	97.73	92.95
3	A	0	TDB	B1-N1-N2	4.11	123.71	122.33
3	B	0	TDB	C7-S1-C14	4.11	97.52	92.95
3	B	0	TDB	O15-S15-N1	-4.05	101.81	106.60
3	B	0	TDB	C7-C12-C13	-4.02	105.35	112.06
2	B	501	NAD	C6N-N1N-C2N	3.92	126.46	122.04
2	B	501	NAD	C2A-N1A-C6A	3.90	125.81	118.77

*Continued on next page...*

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	501	NAD	O4D-C1D-N1N	-3.74	104.12	107.95
2	B	501	NAD	O3D-C3D-C4D	-3.72	100.11	111.08
3	B	0	TDB	B1-N1-N2	3.70	123.57	122.33
3	B	0	TDB	C14-C2-N2	-3.60	119.76	125.73
3	A	0	TDB	O15-S15-N1	-3.45	102.52	106.60
2	A	501	NAD	O4B-C1B-N9A	-3.28	105.39	108.44
2	A	501	NAD	O3D-C3D-C4D	-3.24	101.54	111.08
2	A	501	NAD	C6N-N1N-C2N	3.20	125.65	122.04
3	A	0	TDB	C14-C2-N2	-3.15	120.50	125.73
2	A	501	NAD	C2N-C3N-C7N	-2.97	110.50	119.35
2	B	501	NAD	C6N-N1N-C1D	-2.95	112.02	119.33
2	B	501	NAD	N3A-C4A-N9A	2.93	130.72	125.43
2	A	501	NAD	N3A-C4A-N9A	2.88	130.64	125.43
3	B	0	TDB	C15-S15-N1	2.67	109.17	105.06
2	B	501	NAD	C5A-C6A-N6A	2.62	126.64	120.72
2	B	501	NAD	C2N-C3N-C7N	-2.56	111.74	119.35
2	A	501	NAD	C3B-C2B-C1B	-2.51	96.98	100.91
2	B	501	NAD	C5A-C4A-N3A	-2.50	120.26	125.70
3	A	0	TDB	O16-S15-N1	-2.46	103.69	106.60
2	B	501	NAD	C8A-N9A-C1B	2.40	131.10	126.38
2	A	501	NAD	C6N-N1N-C1D	-2.35	113.50	119.33
2	A	501	NAD	O4D-C4D-C3D	-2.35	100.41	105.17
2	B	501	NAD	O7N-C7N-C3N	-2.31	116.97	119.58
2	B	501	NAD	C6A-C5A-C4A	2.27	121.41	117.25
2	B	501	NAD	C2D-C1D-N1N	2.27	117.70	113.86
2	B	501	NAD	O2B-C2B-C3B	2.23	119.09	111.83
2	B	501	NAD	C6N-C5N-C4N	2.22	122.97	119.44
2	A	501	NAD	C5A-C6A-N6A	2.22	125.75	120.72
2	A	501	NAD	C2B-C3B-C4B	2.20	107.03	102.65
2	A	501	NAD	C3D-C2D-C1D	-2.19	97.48	100.91
2	B	501	NAD	C3D-C2D-C1D	-2.16	97.53	100.91
2	A	501	NAD	C2D-C3D-C4D	2.15	106.94	102.65
2	B	501	NAD	C4B-O4B-C1B	-2.13	107.44	109.75
3	A	0	TDB	C12-C7-S1	2.10	113.65	111.87
3	A	0	TDB	C8-C7-S1	2.07	122.17	119.88
2	A	501	NAD	C5A-C6A-N1A	-2.05	112.40	119.27
3	B	0	TDB	C8-C7-C12	-2.02	121.36	126.28

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 ⓘ

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

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

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

### 6.3 Carbohydrates ⓘ

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

### 6.4 Ligands ⓘ

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

### 6.5 Other polymers ⓘ

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