



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 12, 2017 – 10:42 pm GMT

PDB ID : 1IB0
Title : CRYSTAL STRUCTURE OF RAT B5R IN COMPLEX WITH FAD AND NAD
Authors : Bewley, M.C.; Marohnic, C.C.; Barber, M.J.
Deposited on : 2001-03-26
Resolution : 2.30 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)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Percentile statistics	:	20161228.v01 (using entries in the PDB archive December 28th 2016)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	recalc28949

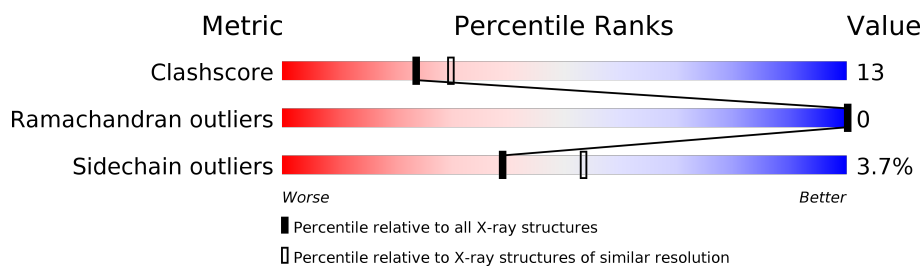
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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	112137	4751 (2.30-2.30)
Ramachandran outliers	110173	4705 (2.30-2.30)
Sidechain outliers	110143	4704 (2.30-2.30)

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	274	

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
3	NAD	A	1994	X	-	-	-

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 2424 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.

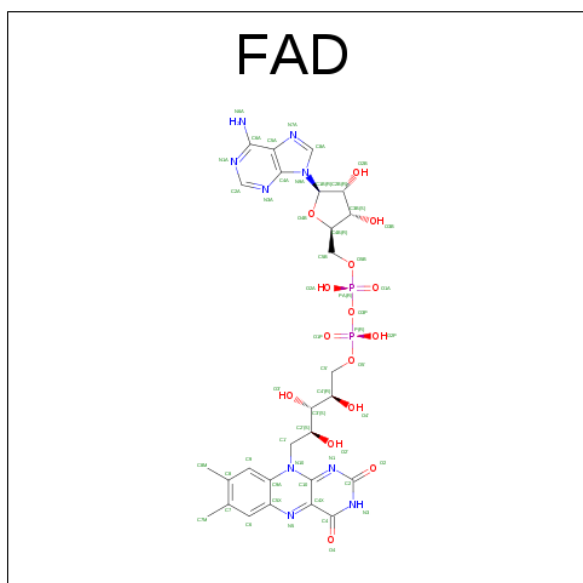
- Molecule 1 is a protein called NADH-CYTOCHROME B5 REDUCTASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	272	2184	1400	380	392	12	0	0	0

There are 8 discrepancies between the modelled and reference sequences:

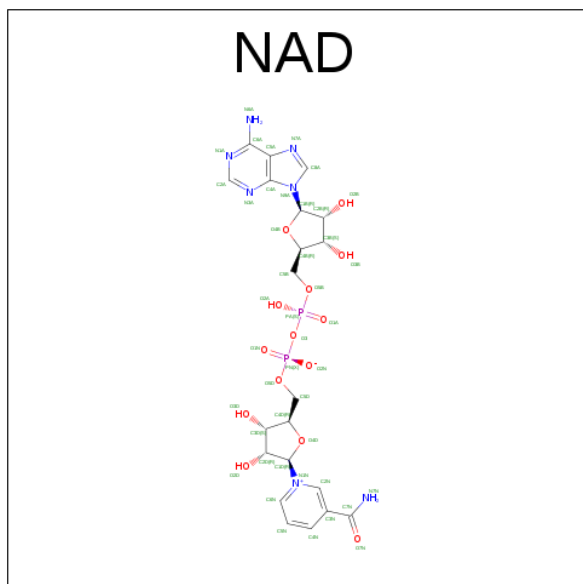
Chain	Residue	Modelled	Actual	Comment	Reference
A	27	MET	-	EXPRESSION TAG	UNP P20070
A	28	HIS	-	EXPRESSION TAG	UNP P20070
A	29	HIS	-	EXPRESSION TAG	UNP P20070
A	30	HIS	-	EXPRESSION TAG	UNP P20070
A	31	HIS	-	EXPRESSION TAG	UNP P20070
A	32	MET	-	EXPRESSION TAG	UNP P20070
A	52	LEU	ILE	CONFLICT	UNP P20070
A	115	GLU	ASP	CONFLICT	UNP P20070

- Molecule 2 is FLAVIN-ADENINE DINUCLEOTIDE (three-letter code: FAD) (formula: $C_{27}H_{33}N_9O_{15}P_2$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			53	27	9	15	2		

- Molecule 3 is NICOTINAMIDE-ADENINE-DINUCLEOTIDE (three-letter code: NAD) (formula: $C_{21}H_{27}N_7O_{14}P_2$).



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

- Molecule 4 is water.

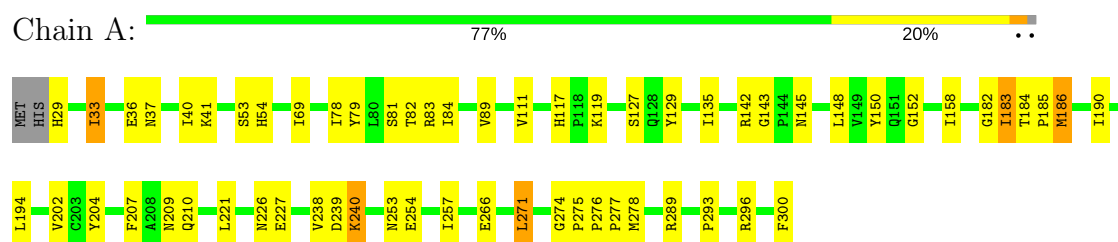
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	143	Total	O	0	0
			143	143		

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.

Note EDS was not executed.

• Molecule 1: NADH-CYTOCHROME B5 REDUCTASE



4 Data and refinement statistics

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

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	67.61Å 69.46Å 78.79Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.00 – 2.30	Depositor
% Data completeness (in resolution range)	(Not available) (30.00-2.30)	Depositor
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.218 , 0.244	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	2424	wwPDB-VP
Average B, all atoms (Å ²)	19.0	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: FAD, 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	0.52	0/2243	0.70	0/3040

There are no bond length outliers.

There are no bond angle outliers.

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	2184	0	2187	54	0
2	A	53	0	31	2	0
3	A	44	0	26	6	0
4	A	143	0	0	1	0
All	All	2424	0	2244	57	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 13.

All (57) 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:186:MET:HE2	1:A:186:MET:HA	1.54	0.89
1:A:186:MET:HA	1:A:186:MET:CE	2.07	0.84
3:A:1994:NAD:H8A	3:A:1994:NAD:H51A	1.74	0.69
1:A:143:GLY:HA2	1:A:145:ASN:ND2	2.11	0.65
1:A:266:GLU:HA	1:A:296:ARG:HE	1.63	0.64
1:A:274:GLY:HA3	1:A:278:MET:CE	2.27	0.63
1:A:275:PRO:HD3	3:A:1994:NAD:H51N	1.81	0.63
1:A:276:PRO:HB2	1:A:277:PRO:HD3	1.81	0.62
1:A:186:MET:CE	1:A:186:MET:CA	2.78	0.60
1:A:274:GLY:HA3	1:A:278:MET:HE1	1.85	0.59
1:A:184:THR:OG1	1:A:185:PRO:HD3	2.02	0.58
1:A:78:ILE:HD12	1:A:78:ILE:C	2.24	0.58
1:A:82:THR:HG23	1:A:84:ILE:HD11	1.85	0.58
1:A:79:TYR:HB2	1:A:142:ARG:HG2	1.85	0.58
2:A:301:FAD:H2'	2:A:301:FAD:N1	2.21	0.56
1:A:84:ILE:HD12	1:A:129:TYR:CD1	2.40	0.56
1:A:202:VAL:HG12	1:A:204:TYR:CE1	2.41	0.56
1:A:210:GLN:O	1:A:240:LYS:HG2	2.08	0.54
1:A:150:TYR:CE2	1:A:152:GLY:HA2	2.44	0.53
1:A:182:GLY:O	1:A:185:PRO:HD2	2.11	0.51
1:A:274:GLY:HA3	1:A:278:MET:HE2	1.93	0.51
1:A:33:ILE:HG22	1:A:69:ILE:HG13	1.92	0.51
1:A:186:MET:HE1	1:A:300:PHE:HE2	1.75	0.50
1:A:210:GLN:HA	1:A:239:ASP:OD2	2.12	0.50
1:A:81:SER:HA	1:A:89:VAL:O	2.12	0.50
1:A:148:LEU:HD23	1:A:148:LEU:C	2.33	0.49
1:A:82:THR:CG2	1:A:84:ILE:HD11	2.43	0.47
1:A:209:ASN:O	1:A:238:VAL:HA	2.15	0.47
1:A:183:ILE:HD11	1:A:221:LEU:HD11	1.96	0.46
1:A:276:PRO:HG3	3:A:1994:NAD:C7N	2.46	0.45
1:A:33:ILE:HD12	1:A:33:ILE:O	2.16	0.45
1:A:186:MET:HE1	1:A:271:LEU:HD23	1.98	0.45
1:A:84:ILE:CD1	1:A:129:TYR:CD1	3.00	0.45
2:A:301:FAD:H1'1	2:A:301:FAD:H9	1.66	0.45
1:A:37:ASN:HB3	1:A:40:ILE:HG12	1.99	0.45
1:A:254:GLU:OE1	1:A:289:ARG:HD2	2.18	0.44
1:A:29:HIS:HE1	1:A:36:GLU:O	2.01	0.43
1:A:117:HIS:CD2	1:A:119:LYS:H	2.37	0.43
1:A:183:ILE:HG13	1:A:207:PHE:CZ	2.53	0.43
1:A:293:PRO:HG3	1:A:296:ARG:HH11	1.84	0.43
1:A:111:VAL:HA	1:A:127:SER:HB2	2.01	0.42
1:A:186:MET:HE3	1:A:186:MET:HA	1.93	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:227:GLU:HB3	4:A:436:HOH:O	2.18	0.42
1:A:83:ARG:C	1:A:84:ILE:HD13	2.40	0.42
1:A:83:ARG:O	1:A:84:ILE:HD13	2.20	0.42
1:A:117:HIS:CD2	1:A:119:LYS:HB2	2.55	0.41
1:A:274:GLY:HA2	3:A:1994:NAD:H4D	2.02	0.41
1:A:253:ASN:O	1:A:257:ILE:HG12	2.21	0.41
1:A:275:PRO:HG3	3:A:1994:NAD:H52A	2.01	0.41
1:A:33:ILE:HG22	1:A:69:ILE:CG1	2.50	0.41
1:A:183:ILE:HB	1:A:207:PHE:CE1	2.55	0.41
1:A:33:ILE:C	1:A:33:ILE:HD12	2.40	0.41
1:A:275:PRO:HG3	3:A:1994:NAD:C5B	2.50	0.41
1:A:53:SER:O	1:A:54:HIS:C	2.58	0.41
1:A:82:THR:HG23	1:A:84:ILE:CD1	2.49	0.40
1:A:190:ILE:O	1:A:194:LEU:HG	2.21	0.40
1:A:276:PRO:CB	1:A:277:PRO:HD3	2.49	0.40

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	270/274 (98%)	261 (97%)	9 (3%)	0	100	100

There are no Ramachandran outliers to report.

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	243/245 (99%)	234 (96%)	9 (4%)	39 53

All (9) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	33	ILE
1	A	41	LYS
1	A	135	ILE
1	A	158	ILE
1	A	183	ILE
1	A	186	MET
1	A	226	ASN
1	A	240	LYS
1	A	271	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	29	HIS
1	A	76	GLN
1	A	117	HIS
1	A	145	ASN
1	A	292	HIS

5.3.3 RNA ⓘ

There are no RNA molecules 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

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	$\# Z > 2$
3	NAD	A	1994	-	41,48,48	1.81	8 (19%)	43,73,73	1.99	11 (25%)
2	FAD	A	301	-	51,58,58	2.24	21 (41%)	54,89,89	3.42	26 (48%)

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	NAD	A	1994	-	3/3/11/11	0/22/62/62	0/5/5/5
2	FAD	A	301	-	-	0/28/50/50	0/6/6/6

All (29) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	301	FAD	PA-O2A	-4.09	1.34	1.55
2	A	301	FAD	P-O5'	-3.00	1.46	1.59
3	A	1994	NAD	O4B-C1B	-2.99	1.37	1.41
2	A	301	FAD	C5A-N7A	-2.93	1.29	1.39
2	A	301	FAD	P-O2P	-2.77	1.41	1.55
2	A	301	FAD	PA-O5B	-2.42	1.48	1.59
3	A	1994	NAD	C2D-C1D	-2.12	1.50	1.53
3	A	1994	NAD	C6N-C5N	-2.07	1.34	1.38
2	A	301	FAD	C8-C7	2.17	1.46	1.41
3	A	1994	NAD	C2A-N1A	2.22	1.38	1.33
2	A	301	FAD	O4B-C4B	2.25	1.50	1.45
2	A	301	FAD	O5'-C5'	2.48	1.54	1.44
2	A	301	FAD	C5B-C4B	2.53	1.59	1.51
2	A	301	FAD	C2A-N1A	2.65	1.38	1.33
2	A	301	FAD	C2-N3	2.70	1.43	1.38
2	A	301	FAD	C4-N3	2.71	1.38	1.33

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	301	FAD	C5X-N5	2.79	1.39	1.35
2	A	301	FAD	C4A-N3A	2.84	1.39	1.35
3	A	1994	NAD	C6N-N1N	2.93	1.43	1.35
2	A	301	FAD	C2A-N3A	3.14	1.37	1.32
2	A	301	FAD	C4X-N5	3.55	1.38	1.33
3	A	1994	NAD	O4D-C1D	3.59	1.46	1.41
2	A	301	FAD	C4-C4X	3.67	1.48	1.41
2	A	301	FAD	O4B-C1B	3.68	1.46	1.41
2	A	301	FAD	C10-N1	3.71	1.38	1.33
2	A	301	FAD	C4X-C10	4.18	1.48	1.41
3	A	1994	NAD	C2N-C3N	5.02	1.46	1.39
2	A	301	FAD	C9A-N10	5.43	1.45	1.38
3	A	1994	NAD	C4N-C3N	6.68	1.50	1.39

All (37) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	301	FAD	C1'-N10-C9A	-8.00	111.01	118.35
2	A	301	FAD	C4-C4X-C10	-7.21	114.13	119.96
2	A	301	FAD	C4'-C3'-C2'	-6.53	99.35	113.41
3	A	1994	NAD	O4D-C4D-C5D	-6.39	87.82	109.40
2	A	301	FAD	C4X-C10-N10	-5.89	116.43	120.52
2	A	301	FAD	O5B-PA-O1A	-4.71	90.27	109.25
2	A	301	FAD	C5X-C9A-N10	-4.68	114.19	117.66
3	A	1994	NAD	O7N-C7N-C3N	-4.55	114.30	119.62
2	A	301	FAD	C4A-C5A-N7A	-3.81	105.73	109.41
2	A	301	FAD	O4B-C4B-C3B	-3.64	97.92	105.17
2	A	301	FAD	O5'-P-O1P	-2.70	98.34	109.25
2	A	301	FAD	C8M-C8-C9	-2.58	113.86	120.34
3	A	1994	NAD	C5N-C4N-C3N	-2.44	117.48	120.35
2	A	301	FAD	N3A-C2A-N1A	-2.21	126.93	128.86
2	A	301	FAD	C4X-C4-N3	-2.17	120.39	123.48
2	A	301	FAD	O2'-C2'-C3'	-2.11	103.86	109.09
3	A	1994	NAD	O4D-C4D-C3D	-2.06	101.08	105.17
2	A	301	FAD	C6-C5X-N5	2.03	121.35	118.97
3	A	1994	NAD	C6N-C5N-C4N	2.12	122.64	119.44
3	A	1994	NAD	C4A-C5A-N7A	2.20	111.54	109.41
3	A	1994	NAD	C5A-C6A-N6A	2.21	124.98	120.47
3	A	1994	NAD	C5N-C6N-N1N	2.45	124.18	120.40
2	A	301	FAD	C1'-C2'-C3'	2.57	117.17	109.82
3	A	1994	NAD	O3D-C3D-C4D	2.69	118.94	111.09
2	A	301	FAD	C5A-C6A-N6A	2.69	125.95	120.47

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	301	FAD	O3'-C3'-C2'	3.03	116.32	108.82
2	A	301	FAD	C8M-C8-C7	3.31	127.67	120.72
2	A	301	FAD	O2A-PA-O5B	3.36	123.99	108.14
3	A	1994	NAD	C4B-O4B-C1B	3.65	113.65	109.77
2	A	301	FAD	C4B-O4B-C1B	3.91	113.93	109.77
2	A	301	FAD	C1'-N10-C10	4.56	123.18	118.50
2	A	301	FAD	O2A-PA-O1A	4.72	136.71	112.28
2	A	301	FAD	C4X-N5-C5X	4.98	122.02	116.76
3	A	1994	NAD	C3N-C7N-N7N	5.35	123.88	117.77
2	A	301	FAD	C4-C4X-N5	6.53	125.84	118.68
2	A	301	FAD	C4-N3-C2	7.97	122.13	115.16
2	A	301	FAD	O4'-C4'-C3'	8.82	130.98	109.09

All (3) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
3	A	1994	NAD	C3D
3	A	1994	NAD	C2D
3	A	1994	NAD	C4D

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	1994	NAD	6	0
2	A	301	FAD	2	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

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.