



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 27, 2014 – 01:08 AM GMT

PDB ID : 3GD4
Title : Crystal structure of the reduced, NAD-bound form of murine apoptosis inducing factor
Authors : Sevrioukova, I.F.
Deposited on : 2009-02-23
Resolution : 2.24 Å(reported)

This is a full wwPDB 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/ValidationPDFNotes.html>

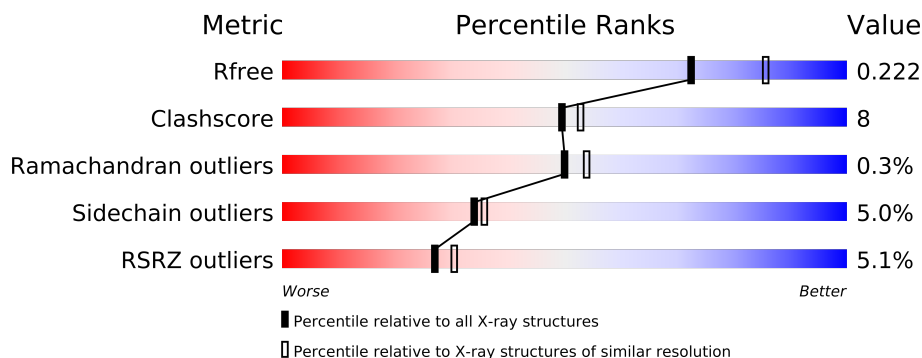
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : dev-1323
EDS : stable22639
Percentile statistics : 21963
Refmac : 5.8.0049
CCP4 : 6.3.0 (Settle)
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 2.24 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}	66092	1112 (2.26-2.22)
Clashscore	79885	1317 (2.26-2.22)
Ramachandran outliers	78287	1282 (2.26-2.22)
Sidechain outliers	78261	1282 (2.26-2.22)
RSRZ outliers	66119	1112 (2.26-2.22)

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. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	511	
1	B	511	

2 Entry composition i

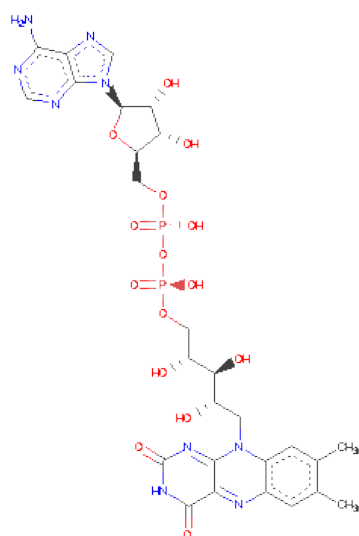
There are 4 unique types of molecules in this entry. The entry contains 7518 atoms, of which 0 are hydrogen and 0 are deuterium.

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 Apoptosis-inducing factor 1, mitochondrial.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	445	Total	C	N	O	S	0	0	0
			3437	2187	609	630	11			
1	B	435	Total	C	N	O	S	0	0	0
			3366	2142	599	614	11			

- 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		
2	B	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		
3	B	1	Total	C	N	O	P	0	0
			44	21	7	14	2		

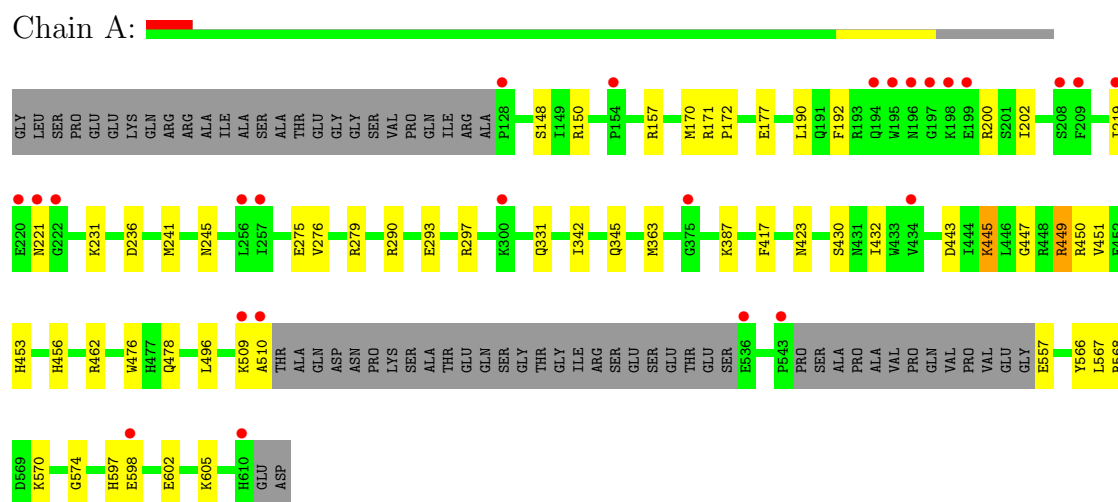
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	257	Total	O	0	0
			257	257		
4	B	264	Total	O	0	0
			264	264		

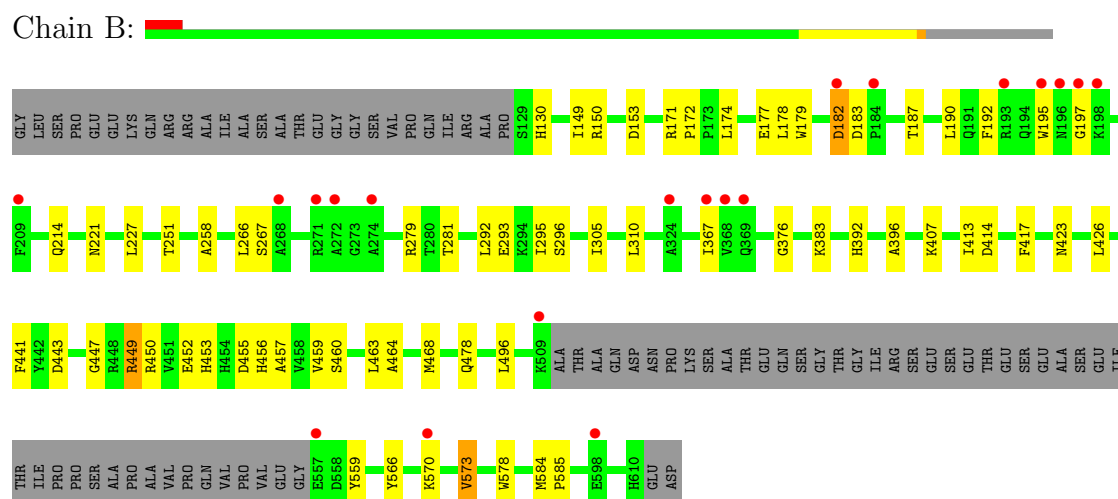
3 Residue-property plots

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.

- Molecule 1: Apoptosis-inducing factor 1, mitochondrial



- Molecule 1: Apoptosis-inducing factor 1, mitochondrial



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	60.26Å 120.39Å 178.19Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.67 – 2.24 30.67 – 2.08	Depositor EDS
% Data completeness (in resolution range)	95.1 (30.67-2.24) 80.4 (30.67-2.08)	Depositor EDS
R_{merge}	0.04	Depositor
R_{sym}	0.04	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.83 (at 2.08Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.187 , 0.222 0.188 , 0.222	Depositor DCC
R_{free} test set	3182 reflections (5.32%)	DCC
Wilson B-factor (Å ²)	41.4	Xtriage
Anisotropy	0.245	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 39.8	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 63021 reflections	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	7518	wwPDB-VP
Average B, all atoms (Å ²)	48.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.16% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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	0.59	0/3506	0.69	0/4736
1	B	0.62	0/3434	0.74	1/4638 (0.0%)
All	All	0.61	0/6940	0.71	1/9374 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	449	ARG	NE-CZ-NH1	5.23	122.92	120.30

There are no chirality outliers.

There are no planarity outliers.

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	3437	0	7	24	0
1	B	3366	0	0	26	0
2	A	53	0	30	1	0
2	B	53	0	30	3	0
3	A	44	0	26	1	0
3	B	44	0	26	3	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	A	257	0	0	6	0
4	B	264	0	0	1	0
All	All	7518	0	119	54	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 (54) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:150:ARG:NH1	1:B:221:ASN:O	2.23	0.71
1:B:449:ARG:NH2	1:B:478:GLN:NE2	2.44	0.66
1:A:293:GLU:OE2	1:A:297:ARG:NH2	2.32	0.62
1:A:449:ARG:NH2	1:A:478:GLN:OE1	2.32	0.61
1:A:453:HIS:NE2	4:A:643:HOH:O	2.31	0.60
1:B:279:ARG:NH1	1:B:376:GLY:O	2.35	0.59
1:A:157:ARG:NH1	1:A:219:ILE:O	2.38	0.57
1:A:150:ARG:NH2	1:A:221:ASN:O	2.39	0.56
1:B:171:ARG:N	1:B:172:PRO:CD	2.69	0.55
1:B:174:LEU:O	1:B:179:TRP:NE1	2.39	0.55
1:A:236:ASP:N	1:A:241:MET:O	2.39	0.55
1:B:367:ILE:O	1:B:383:LYS:N	2.40	0.55
1:A:445:LYS:NZ	1:A:476:TRP:O	2.40	0.54
1:A:509:LYS:O	1:A:510:ALA:CB	2.56	0.53
2:B:1611:FAD:H1'1	3:B:700:NAD:H1D	1.92	0.50
1:A:170:MET:N	1:A:202:ILE:O	2.45	0.49
1:A:568:ARG:NE	4:A:812:HOH:O	2.44	0.49
1:A:570:LYS:N	4:A:699:HOH:O	2.44	0.49
1:B:443:ASP:O	1:B:447:GLY:N	2.46	0.49
1:B:559:TYR:O	1:B:578:TRP:CD1	2.65	0.48
1:A:171:ARG:N	1:A:172:PRO:CD	2.76	0.48
1:B:130:HIS:ND1	1:B:251:THR:OG1	2.48	0.47
1:B:566:TYR:O	1:B:573:VAL:N	2.47	0.47
1:A:566:TYR:N	1:A:574:GLY:O	2.47	0.47
1:A:279:ARG:NH1	4:A:844:HOH:O	2.48	0.47
1:B:258:ALA:O	2:B:1611:FAD:H52A	2.15	0.47
1:B:460:SER:O	1:B:464:ALA:N	2.48	0.46
1:B:456:HIS:O	1:B:460:SER:N	2.48	0.46
1:A:331:GLN:O	1:A:363:MET:N	2.49	0.45
1:B:281:THR:OG1	1:B:392:HIS:NE2	2.49	0.45
1:A:177:GLU:CG	1:A:192:PHE:CD2	2.99	0.45
1:B:453:HIS:O	1:B:457:ALA:N	2.50	0.44
1:B:177:GLU:CG	1:B:192:PHE:CD2	3.00	0.44

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:292:LEU:O	1:B:296:SER:N	2.51	0.44
1:A:451:VAL:CG2	1:A:456:HIS:CB	2.96	0.43
1:B:464:ALA:O	1:B:468:MET:N	2.51	0.43
1:A:290:ARG:NH1	4:A:1059:HOH:O	2.51	0.43
1:A:602:GLU:OE1	1:A:605:LYS:NZ	2.52	0.43
1:B:149:ILE:O	1:B:153:ASP:N	2.51	0.43
1:B:455:ASP:O	1:B:459:VAL:N	2.51	0.43
1:B:566:TYR:CD1	1:B:566:TYR:N	2.87	0.42
1:A:148:SER:OG	1:A:462:ARG:O	2.37	0.42
1:A:443:ASP:O	1:A:447:GLY:N	2.53	0.42
1:B:450:ARG:NH1	1:B:452:GLU:OE2	2.53	0.42
1:B:414:ASP:OD1	1:B:417:PHE:N	2.53	0.41
1:B:441:PHE:N	1:B:449:ARG:O	2.53	0.41
3:B:700:NAD:H8A	4:B:996:HOH:O	2.20	0.41
1:A:597:HIS:CE1	4:A:641:HOH:O	2.73	0.41
1:A:231:LYS:N	1:A:245:ASN:ND2	2.68	0.41
2:B:1611:FAD:C1'	3:B:700:NAD:H1D	2.49	0.41
1:A:417:PHE:O	1:A:450:ARG:NE	2.53	0.41
1:B:584:MET:N	1:B:585:PRO:CD	2.84	0.41
2:A:1611:FAD:C1'	3:A:700:NAD:H1D	2.51	0.41
1:B:305:ILE:O	1:B:396:ALA:N	2.53	0.41

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	439/511 (86%)	424 (97%)	15 (3%)	0	100	100
1	B	431/511 (84%)	411 (95%)	17 (4%)	3 (1%)	30	28
All	All	870/1022 (85%)	835 (96%)	32 (4%)	3 (0%)	50	54

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	197	GLY
1	B	183	ASP
1	B	182	ASP

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	365/417 (88%)	349 (96%)	16 (4%)	39	43
1	B	357/417 (86%)	337 (94%)	20 (6%)	30	29
All	All	722/834 (87%)	686 (95%)	36 (5%)	34	36

All (36) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	190	LEU
1	A	200	ARG
1	A	275	GLU
1	A	276	VAL
1	A	342	ILE
1	A	345	GLN
1	A	387	LYS
1	A	423	ASN
1	A	430	SER
1	A	432	ILE
1	A	445	LYS
1	A	449	ARG
1	A	496	LEU
1	A	557	GLU
1	A	567	LEU
1	A	598	GLU
1	B	178	LEU
1	B	182	ASP
1	B	187	THR
1	B	190	LEU
1	B	195	TRP
1	B	214	GLN
1	B	227	LEU

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Mol	Chain	Res	Type
1	B	266	LEU
1	B	267	SER
1	B	293	GLU
1	B	295	ILE
1	B	310	LEU
1	B	407	LYS
1	B	413	ILE
1	B	423	ASN
1	B	426	LEU
1	B	463	LEU
1	B	496	LEU
1	B	570	LYS
1	B	573	VAL

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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
2	FAD	A	1611	-	58,58,58	1.77	6 (10%)	85,89,89	3.13	16 (18%)
3	NAD	A	700	-	48,48,48	1.43	4 (8%)	73,73,73	2.03	13 (17%)
2	FAD	B	1611	-	58,58,58	1.89	8 (13%)	85,89,89	3.28	18 (21%)
3	NAD	B	700	-	48,48,48	1.36	4 (8%)	73,73,73	1.99	9 (12%)

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	FAD	A	1611	-	-	0/34/50/50	0/1/6/6
3	NAD	A	700	-	-	0/30/62/62	0/3/5/5
2	FAD	B	1611	-	-	0/34/50/50	0/1/6/6
3	NAD	B	700	-	-	0/30/62/62	0/3/5/5

All (22) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	1611	FAD	C6-C5X	-11.07	1.28	1.41
2	A	1611	FAD	C6-C5X	-10.57	1.29	1.41
3	A	700	NAD	O7N-C7N	6.43	1.39	1.24
3	B	700	NAD	O7N-C7N	6.29	1.38	1.24
2	B	1611	FAD	C4A-N9A	-3.67	1.32	1.37
2	A	1611	FAD	C9A-C5X	3.40	1.49	1.42
2	B	1611	FAD	C6-C7	-3.25	1.28	1.37
2	B	1611	FAD	C9A-C5X	3.12	1.49	1.42
2	A	1611	FAD	C6-C7	-3.10	1.29	1.37
3	A	700	NAD	C2A-N3A	3.08	1.38	1.32
3	B	700	NAD	C2A-N3A	3.07	1.38	1.32
2	A	1611	FAD	C5A-C4A	2.87	1.47	1.40
3	A	700	NAD	C2A-N1A	2.77	1.39	1.33
3	A	700	NAD	O4B-C4B	-2.76	1.38	1.45
2	B	1611	FAD	C5A-C4A	2.57	1.46	1.40
3	B	700	NAD	C2A-N1A	2.44	1.38	1.33
2	A	1611	FAD	C4A-N9A	-2.40	1.34	1.37
2	B	1611	FAD	C1'-N10	-2.34	1.45	1.48
3	B	700	NAD	O4B-C4B	-2.29	1.39	1.45
2	B	1611	FAD	C1'-C2'	2.27	1.53	1.51
2	B	1611	FAD	C2B-C1B	-2.23	1.50	1.53
2	A	1611	FAD	C1'-C2'	2.10	1.53	1.51

All (56) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	1611	FAD	C7-C6-C5X	16.37	148.38	120.91
2	A	1611	FAD	C7-C6-C5X	16.00	147.76	120.91
2	B	1611	FAD	C6-C5X-N5	13.47	134.67	118.97
2	A	1611	FAD	C6-C5X-N5	12.60	133.65	118.97
3	A	700	NAD	N3A-C2A-N1A	-11.89	118.77	128.71
3	B	700	NAD	N3A-C2A-N1A	-11.63	118.99	128.71
2	B	1611	FAD	C6-C5X-C9A	-11.12	103.61	119.02
2	A	1611	FAD	C6-C5X-C9A	-9.98	105.19	119.02
2	B	1611	FAD	N3A-C2A-N1A	-8.73	121.41	128.71
2	A	1611	FAD	C6-C7-C8	-7.75	107.48	119.88
2	B	1611	FAD	C6-C7-C8	-7.20	108.36	119.88
2	A	1611	FAD	N3A-C2A-N1A	-6.63	123.17	128.71
2	A	1611	FAD	C2-N1-C10	6.34	121.37	114.98
2	B	1611	FAD	C2-N1-C10	5.46	120.48	114.98
2	A	1611	FAD	N3A-C4A-N9A	5.13	134.70	125.43
2	B	1611	FAD	C2'-C1'-N10	-4.71	106.20	112.45
3	B	700	NAD	C3N-C7N-N7N	4.48	122.87	117.77
3	A	700	NAD	O4D-C1D-N1N	4.36	112.42	107.95
2	B	1611	FAD	N3A-C4A-N9A	4.36	133.30	125.43
2	A	1611	FAD	C4A-C5A-N7A	-3.96	106.13	109.52
2	A	1611	FAD	C9A-N10-C10	-3.87	117.97	121.77
3	B	700	NAD	O7N-C7N-C3N	-3.77	115.33	119.58
3	A	700	NAD	O4B-C1B-N9A	-3.69	105.01	108.44
2	B	1611	FAD	C8A-N9A-C4A	3.56	109.62	106.90
2	B	1611	FAD	C4A-C5A-N7A	-3.54	106.49	109.52
3	A	700	NAD	C3N-C7N-N7N	3.54	121.80	117.77
2	B	1611	FAD	C1B-N9A-C4A	-3.50	120.59	126.64
3	B	700	NAD	C8A-N9A-C4A	3.43	109.52	106.90
2	B	1611	FAD	C7M-C7-C6	3.39	128.56	120.38
2	A	1611	FAD	C7M-C7-C6	3.35	128.45	120.38
2	A	1611	FAD	C5A-C4A-N3A	-3.20	118.73	125.70
3	A	700	NAD	O5B-C5B-C4B	-3.18	97.25	108.94
3	A	700	NAD	N3A-C4A-N9A	3.18	131.17	125.43
3	B	700	NAD	N3A-C4A-N9A	3.15	131.13	125.43
2	A	1611	FAD	C2'-C1'-N10	-3.03	108.44	112.45
2	A	1611	FAD	O4B-C1B-N9A	3.01	111.24	108.44
2	A	1611	FAD	C4X-N5-C5X	2.96	120.02	116.69
3	B	700	NAD	C1B-N9A-C4A	-2.92	121.60	126.64
3	A	700	NAD	C6N-N1N-C2N	-2.83	118.85	122.04
2	B	1611	FAD	C9A-N10-C10	-2.80	119.02	121.77
3	A	700	NAD	C8A-N9A-C4A	2.76	109.01	106.90
2	A	1611	FAD	C1'-N10-C9A	2.68	121.48	118.87

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	700	NAD	O5B-C5B-C4B	-2.63	99.27	108.94
3	A	700	NAD	O7N-C7N-C3N	-2.61	116.64	119.58
2	B	1611	FAD	C2A-N1A-C6A	2.51	123.31	118.77
3	B	700	NAD	N7A-C8A-N9A	-2.43	107.49	114.36
2	B	1611	FAD	C4X-N5-C5X	2.35	119.33	116.69
3	A	700	NAD	C1B-N9A-C4A	-2.33	122.61	126.64
3	A	700	NAD	C2N-C3N-C4N	2.33	120.95	118.31
2	B	1611	FAD	C5A-C4A-N3A	-2.30	120.69	125.70
2	A	1611	FAD	C4X-C10-N1	-2.29	120.44	122.73
3	B	700	NAD	C3B-C2B-C1B	2.22	104.39	100.91
3	A	700	NAD	C4D-O4D-C1D	2.20	112.14	109.75
2	B	1611	FAD	O2A-PA-O1A	2.11	124.02	112.21
3	A	700	NAD	N7A-C8A-N9A	-2.11	108.39	114.36
2	B	1611	FAD	C4-N3-C2	-2.03	121.22	125.39

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 ⓘ

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	445/511 (87%)	0.25	25 (5%) 24 27	28, 46, 72, 84	0
1	B	435/511 (85%)	0.30	20 (4%) 31 34	27, 46, 79, 93	0
All	All	880/1022 (86%)	0.27	45 (5%) 27 30	27, 46, 77, 93	0

All (45) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	195	TRP	11.2
1	A	195	TRP	8.4
1	B	182	ASP	5.9
1	A	510	ALA	5.7
1	A	128	PRO	5.6
1	A	197	GLY	5.3
1	A	196	ASN	5.0
1	B	197	GLY	4.8
1	B	184	PRO	4.5
1	A	509	LYS	4.2
1	A	198	LYS	4.0
1	B	274	ALA	3.9
1	B	509	LYS	3.7
1	A	375	GLY	3.6
1	B	271	ARG	3.5
1	A	194	GLN	3.5
1	B	557	GLU	3.4
1	A	536	GLU	3.3
1	B	324	ALA	3.3
1	A	610	HIS	3.2
1	A	598	GLU	3.1
1	B	209	PHE	3.0
1	B	272	ALA	2.9
1	B	368	VAL	2.9

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Mol	Chain	Res	Type	RSRZ
1	A	543	PRO	2.8
1	A	220	GLU	2.8
1	A	222	GLY	2.8
1	B	198	LYS	2.8
1	A	300	LYS	2.7
1	A	221	ASN	2.6
1	B	369	GLN	2.6
1	B	598	GLU	2.5
1	B	268	ALA	2.5
1	B	196	ASN	2.4
1	A	199	GLU	2.4
1	B	570	LYS	2.4
1	A	209	PHE	2.4
1	B	193	ARG	2.2
1	A	256	LEU	2.2
1	A	154	PRO	2.2
1	B	367	ILE	2.1
1	A	208	SER	2.1
1	A	257	ILE	2.1
1	A	434	VAL	2.1
1	A	219	ILE	2.0

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

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

6.4 Ligands ⓘ

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	RSR	LLDF	B-factors(Å ²)	Q<0.9
3	NAD	B	700	44/44	0.18	0.49	30,39,56,57	0

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Mol	Type	Chain	Res	Atoms	RSR	LLDF	B-factors(\AA^2)	Q<0.9
2	FAD	B	1611	53/53	0.17	0.38	32,35,37,41	0
2	FAD	A	1611	53/53	0.14	-0.13	33,36,42,44	0
3	NAD	A	700	44/44	0.12	-0.46	30,37,50,51	0

6.5 Other polymers ⓘ

There are no such residues in this entry.