



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 14, 2017 – 04:16 pm GMT

PDB ID : 1Z0Z
Title : Crystal structure of a NAD kinase from *Archaeoglobus fulgidus* in complex with NAD
Authors : Liu, J.; Lou, Y.; Yokota, H.; Adams, P.D.; Kim, R.; Kim, S.H.; Berkeley Structural Genomics Center (BSGC)
Deposited on : 2005-03-02
Resolution : 2.85 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)	:	1.9-1692
EDS	:	trunk28620
Percentile statistics	:	20161228.v01 (using entries in the PDB archive December 28th 2016)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	recalc28949

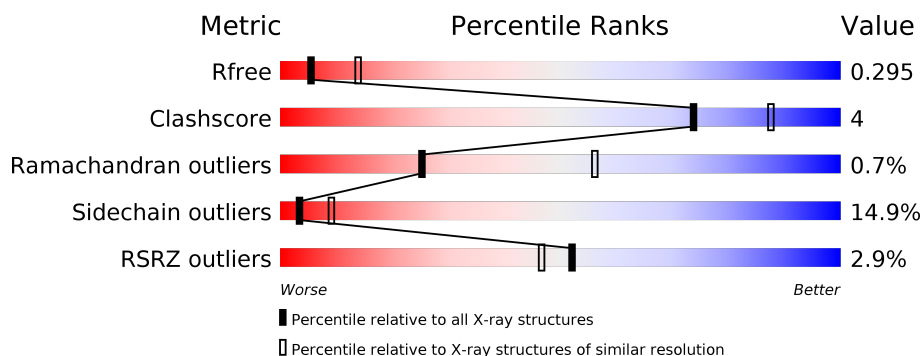
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.85 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}	100719	2469 (2.90-2.82)
Clashscore	112137	2749 (2.90-2.82)
Ramachandran outliers	110173	2687 (2.90-2.82)
Sidechain outliers	110143	2690 (2.90-2.82)
RSRZ outliers	101464	2487 (2.90-2.82)

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.

Mol	Chain	Length	Quality of chain
1	A	278	<div> <div>3%</div> <div> <div></div> <div>70%</div> <div>15%</div> <div>•</div> <div>10%</div> </div> </div>
1	B	278	<div> <div>%</div> <div> <div></div> <div>67%</div> <div>20%</div> <div>•</div> <div>10%</div> </div> </div>
1	C	278	<div> <div>4%</div> <div> <div></div> <div>71%</div> <div>17%</div> <div>•</div> <div>10%</div> </div> </div>
1	D	278	<div> <div>3%</div> <div> <div></div> <div>68%</div> <div>21%</div> <div>•</div> <div>10%</div> </div> </div>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 8150 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 Probable inorganic polyphosphate/ATP-NAD kinase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	249	Total	C	N	O	S	0	0	0
			1964	1266	334	356	8			
1	B	249	Total	C	N	O	S	0	0	0
			1964	1266	334	356	8			
1	C	249	Total	C	N	O	S	0	0	0
			1964	1266	334	356	8			
1	D	249	Total	C	N	O	S	0	0	0
			1964	1266	334	356	8			

There are 116 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-28	MET	-	cloning artifact	UNP O30297
A	-27	GLY	-	cloning artifact	UNP O30297
A	-26	SER	-	cloning artifact	UNP O30297
A	-25	SER	-	cloning artifact	UNP O30297
A	-24	HIS	-	cloning artifact	UNP O30297
A	-23	HIS	-	cloning artifact	UNP O30297
A	-22	HIS	-	cloning artifact	UNP O30297
A	-21	HIS	-	cloning artifact	UNP O30297
A	-20	HIS	-	cloning artifact	UNP O30297
A	-19	HIS	-	cloning artifact	UNP O30297
A	-18	ASP	-	cloning artifact	UNP O30297
A	-17	TYR	-	cloning artifact	UNP O30297
A	-16	ASP	-	cloning artifact	UNP O30297
A	-15	ILE	-	cloning artifact	UNP O30297
A	-14	PRO	-	cloning artifact	UNP O30297
A	-13	THR	-	cloning artifact	UNP O30297
A	-12	THR	-	cloning artifact	UNP O30297
A	-11	GLU	-	cloning artifact	UNP O30297
A	-10	ASN	-	cloning artifact	UNP O30297
A	-9	LEU	-	cloning artifact	UNP O30297
A	-8	TYR	-	cloning artifact	UNP O30297

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Chain	Residue	Modelled	Actual	Comment	Reference
A	-7	PHE	-	cloning artifact	UNP O30297
A	-6	GLN	-	cloning artifact	UNP O30297
A	-5	GLY	-	cloning artifact	UNP O30297
A	-4	GLY	-	cloning artifact	UNP O30297
A	-3	GLY	-	cloning artifact	UNP O30297
A	-2	GLY	-	cloning artifact	UNP O30297
A	-1	GLY	-	cloning artifact	UNP O30297
A	0	GLY	-	cloning artifact	UNP O30297
B	-28	MET	-	cloning artifact	UNP O30297
B	-27	GLY	-	cloning artifact	UNP O30297
B	-26	SER	-	cloning artifact	UNP O30297
B	-25	SER	-	cloning artifact	UNP O30297
B	-24	HIS	-	cloning artifact	UNP O30297
B	-23	HIS	-	cloning artifact	UNP O30297
B	-22	HIS	-	cloning artifact	UNP O30297
B	-21	HIS	-	cloning artifact	UNP O30297
B	-20	HIS	-	cloning artifact	UNP O30297
B	-19	HIS	-	cloning artifact	UNP O30297
B	-18	ASP	-	cloning artifact	UNP O30297
B	-17	TYR	-	cloning artifact	UNP O30297
B	-16	ASP	-	cloning artifact	UNP O30297
B	-15	ILE	-	cloning artifact	UNP O30297
B	-14	PRO	-	cloning artifact	UNP O30297
B	-13	THR	-	cloning artifact	UNP O30297
B	-12	THR	-	cloning artifact	UNP O30297
B	-11	GLU	-	cloning artifact	UNP O30297
B	-10	ASN	-	cloning artifact	UNP O30297
B	-9	LEU	-	cloning artifact	UNP O30297
B	-8	TYR	-	cloning artifact	UNP O30297
B	-7	PHE	-	cloning artifact	UNP O30297
B	-6	GLN	-	cloning artifact	UNP O30297
B	-5	GLY	-	cloning artifact	UNP O30297
B	-4	GLY	-	cloning artifact	UNP O30297
B	-3	GLY	-	cloning artifact	UNP O30297
B	-2	GLY	-	cloning artifact	UNP O30297
B	-1	GLY	-	cloning artifact	UNP O30297
B	0	GLY	-	cloning artifact	UNP O30297
C	-28	MET	-	cloning artifact	UNP O30297
C	-27	GLY	-	cloning artifact	UNP O30297
C	-26	SER	-	cloning artifact	UNP O30297
C	-25	SER	-	cloning artifact	UNP O30297
C	-24	HIS	-	cloning artifact	UNP O30297

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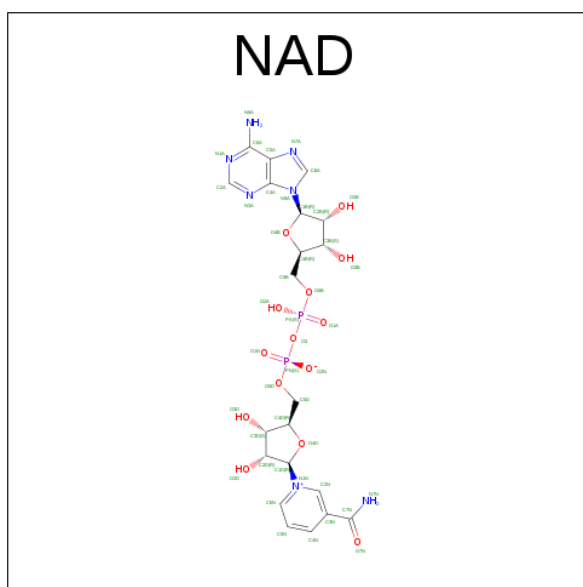
Chain	Residue	Modelled	Actual	Comment	Reference
C	-23	HIS	-	cloning artifact	UNP O30297
C	-22	HIS	-	cloning artifact	UNP O30297
C	-21	HIS	-	cloning artifact	UNP O30297
C	-20	HIS	-	cloning artifact	UNP O30297
C	-19	HIS	-	cloning artifact	UNP O30297
C	-18	ASP	-	cloning artifact	UNP O30297
C	-17	TYR	-	cloning artifact	UNP O30297
C	-16	ASP	-	cloning artifact	UNP O30297
C	-15	ILE	-	cloning artifact	UNP O30297
C	-14	PRO	-	cloning artifact	UNP O30297
C	-13	THR	-	cloning artifact	UNP O30297
C	-12	THR	-	cloning artifact	UNP O30297
C	-11	GLU	-	cloning artifact	UNP O30297
C	-10	ASN	-	cloning artifact	UNP O30297
C	-9	LEU	-	cloning artifact	UNP O30297
C	-8	TYR	-	cloning artifact	UNP O30297
C	-7	PHE	-	cloning artifact	UNP O30297
C	-6	GLN	-	cloning artifact	UNP O30297
C	-5	GLY	-	cloning artifact	UNP O30297
C	-4	GLY	-	cloning artifact	UNP O30297
C	-3	GLY	-	cloning artifact	UNP O30297
C	-2	GLY	-	cloning artifact	UNP O30297
C	-1	GLY	-	cloning artifact	UNP O30297
C	0	GLY	-	cloning artifact	UNP O30297
D	-28	MET	-	cloning artifact	UNP O30297
D	-27	GLY	-	cloning artifact	UNP O30297
D	-26	SER	-	cloning artifact	UNP O30297
D	-25	SER	-	cloning artifact	UNP O30297
D	-24	HIS	-	cloning artifact	UNP O30297
D	-23	HIS	-	cloning artifact	UNP O30297
D	-22	HIS	-	cloning artifact	UNP O30297
D	-21	HIS	-	cloning artifact	UNP O30297
D	-20	HIS	-	cloning artifact	UNP O30297
D	-19	HIS	-	cloning artifact	UNP O30297
D	-18	ASP	-	cloning artifact	UNP O30297
D	-17	TYR	-	cloning artifact	UNP O30297
D	-16	ASP	-	cloning artifact	UNP O30297
D	-15	ILE	-	cloning artifact	UNP O30297
D	-14	PRO	-	cloning artifact	UNP O30297
D	-13	THR	-	cloning artifact	UNP O30297
D	-12	THR	-	cloning artifact	UNP O30297
D	-11	GLU	-	cloning artifact	UNP O30297

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Chain	Residue	Modelled	Actual	Comment	Reference
D	-10	ASN	-	cloning artifact	UNP O30297
D	-9	LEU	-	cloning artifact	UNP O30297
D	-8	TYR	-	cloning artifact	UNP O30297
D	-7	PHE	-	cloning artifact	UNP O30297
D	-6	GLN	-	cloning artifact	UNP O30297
D	-5	GLY	-	cloning artifact	UNP O30297
D	-4	GLY	-	cloning artifact	UNP O30297
D	-3	GLY	-	cloning artifact	UNP O30297
D	-2	GLY	-	cloning artifact	UNP O30297
D	-1	GLY	-	cloning artifact	UNP O30297
D	0	GLY	-	cloning artifact	UNP O30297

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

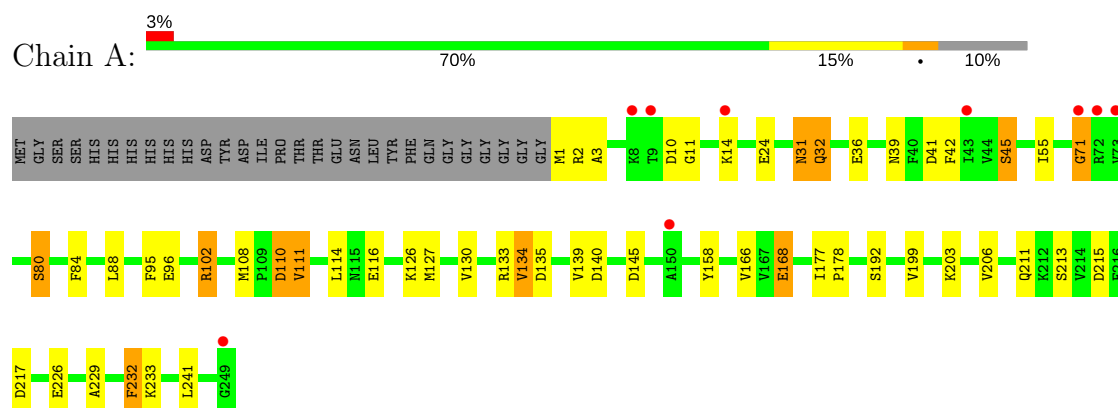
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	33	Total 33	O 33	0	0
3	B	31	Total 31	O 31	0	0
3	C	24	Total 24	O 24	0	0
3	D	30	Total 30	O 30	0	0

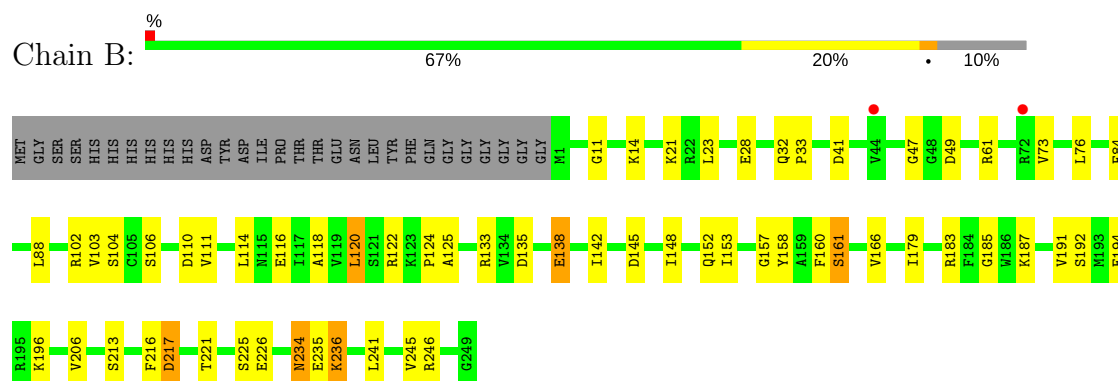
3 Residue-property plots [i](#)

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.

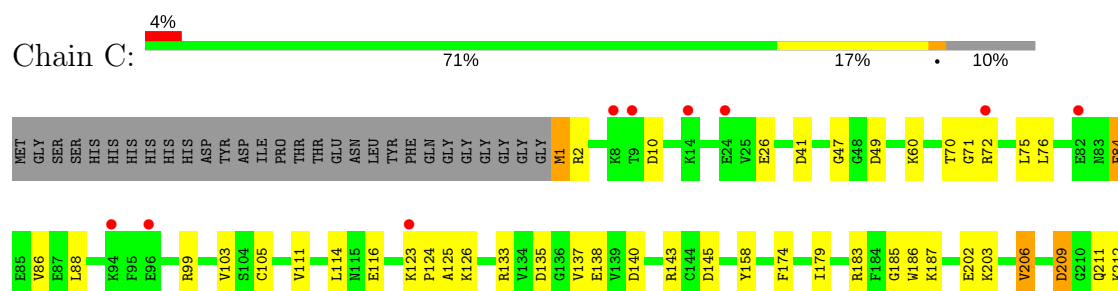
- Molecule 1: Probable inorganic polyphosphate/ATP-NAD kinase

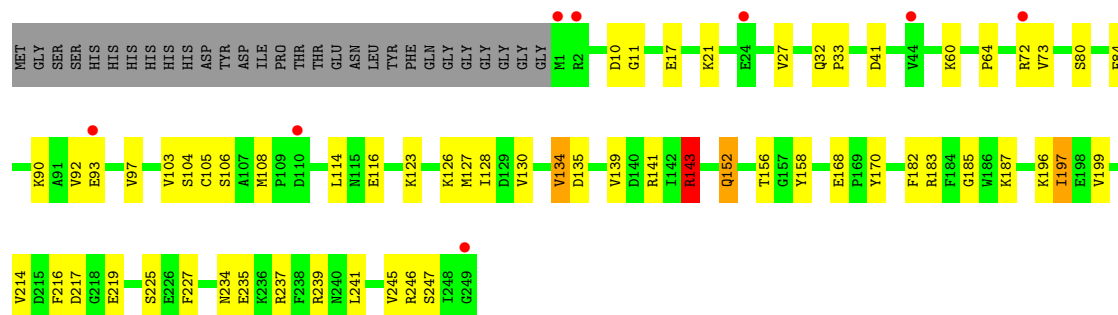


- Molecule 1: Probable inorganic polyphosphate/ATP-NAD kinase



- Molecule 1: Probable inorganic polyphosphate/ATP-NAD kinase





4 Data and refinement statistics

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, α , β , γ	122.41 Å 122.41 Å 201.87 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 2.85 20.00 – 2.85	Depositor EDS
% Data completeness (in resolution range)	(Not available) (20.00-2.85) 76.1 (20.00-2.85)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.13 (at 2.83 Å)	Xtriage
Refinement program	REFMAC 5.1.24	Depositor
R, R_{free}	0.208 , 0.275 0.255 , 0.295	Depositor DCC
R_{free} test set	1364 reflections (4.94%)	DCC
Wilson B-factor (Å ²)	67.7	Xtriage
Anisotropy	0.370	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , -3.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	8150	wwPDB-VP
Average B, all atoms (Å ²)	39.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 50.20 % of the origin peak, indicating pseudo translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo translational symmetry is equal to 6.6888e-05. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

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	0.53	0/2002	0.90	11/2699 (0.4%)
1	B	0.52	0/2002	0.79	5/2699 (0.2%)
1	C	0.53	0/2002	0.76	6/2699 (0.2%)
1	D	0.52	0/2002	0.79	5/2699 (0.2%)
All	All	0.53	0/8008	0.81	27/10796 (0.3%)

There are no bond length outliers.

All (27) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	110	ASP	CB-CA-C	-18.19	74.02	110.40
1	A	71	GLY	N-CA-C	12.10	143.35	113.10
1	D	143	ARG	NE-CZ-NH1	8.93	124.77	120.30
1	A	111	VAL	N-CA-C	-7.95	89.53	111.00
1	C	145	ASP	CB-CG-OD2	7.18	124.76	118.30
1	D	135	ASP	CB-CG-OD2	7.16	124.74	118.30
1	B	49	ASP	CB-CG-OD2	7.15	124.74	118.30
1	D	10	ASP	CB-CG-OD2	6.80	124.42	118.30
1	A	140	ASP	CB-CG-OD2	6.50	124.15	118.30
1	B	110	ASP	CB-CG-OD2	6.47	124.12	118.30
1	A	41	ASP	CB-CG-OD2	6.27	123.94	118.30
1	B	217	ASP	CB-CG-OD2	5.89	123.60	118.30
1	D	143	ARG	NE-CZ-NH2	-5.83	117.38	120.30
1	B	135	ASP	CB-CG-OD2	5.70	123.43	118.30
1	C	10	ASP	CB-CG-OD2	5.69	123.42	118.30
1	C	41	ASP	CB-CG-OD2	5.69	123.42	118.30
1	C	209	ASP	CB-CG-OD2	5.66	123.40	118.30
1	A	217	ASP	CB-CG-OD2	5.66	123.39	118.30
1	A	215	ASP	CB-CG-OD2	5.56	123.30	118.30
1	C	135	ASP	CB-CG-OD2	5.49	123.24	118.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	110	ASP	CB-CG-OD2	5.45	123.20	118.30
1	D	41	ASP	CB-CG-OD2	5.32	123.09	118.30
1	C	49	ASP	CB-CG-OD2	5.31	123.08	118.30
1	B	41	ASP	CB-CG-OD2	5.29	123.06	118.30
1	A	145	ASP	CB-CG-OD2	5.20	122.98	118.30
1	A	135	ASP	CB-CG-OD2	5.05	122.85	118.30
1	A	10	ASP	CB-CG-OD2	5.01	122.81	118.30

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	1964	0	2015	15	0
1	B	1964	0	2015	15	0
1	C	1964	0	2015	15	0
1	D	1964	0	2015	19	0
2	A	44	0	26	0	0
2	B	44	0	26	1	0
2	C	44	0	26	0	0
2	D	44	0	26	2	0
3	A	33	0	0	2	0
3	B	31	0	0	0	0
3	C	24	0	0	2	0
3	D	30	0	0	1	0
All	All	8150	0	8164	62	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 4.

All (62) 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:110:ASP:O	3:A:3100:HOH:O	1.86	0.92

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:183:ARG:HD2	1:D:185:GLY:O	1.85	0.76
1:D:123:LYS:NZ	3:D:3089:HOH:O	2.29	0.66
1:A:1:MET:N	3:A:3093:HOH:O	1.89	0.64
1:A:166:VAL:HG23	1:B:187:LYS:HG2	1.79	0.64
1:C:70:THR:HG23	1:C:71:GLY:H	1.62	0.63
1:C:209:ASP:O	1:C:211:GLN:HG2	2.00	0.62
1:C:183:ARG:HD2	1:C:185:GLY:O	2.02	0.59
1:B:183:ARG:HD2	1:B:185:GLY:O	2.03	0.59
1:C:236:LYS:NZ	3:C:3083:HOH:O	2.31	0.58
1:B:133:ARG:HG2	1:B:138:GLU:HA	1.87	0.57
1:A:31:ASN:HD22	1:A:32:GLN:HG2	1.70	0.56
1:D:143:ARG:HG3	1:D:143:ARG:HH11	1.69	0.56
1:D:152:GLN:HE21	1:D:152:GLN:H	1.52	0.55
1:D:127:MET:HE1	1:D:182:PHE:HB2	1.89	0.55
1:A:134:VAL:HG13	1:A:139:VAL:HG21	1.89	0.54
1:C:124:PRO:O	1:C:125:ALA:HB3	2.08	0.54
1:C:1:MET:N	3:C:3078:HOH:O	2.39	0.52
1:D:32:GLN:HB2	1:D:33:PRO:CD	2.40	0.51
1:D:106:SER:O	1:D:106:SER:OG	2.25	0.51
1:C:140:ASP:OD1	1:C:183:ARG:NH2	2.41	0.51
1:C:202:GLU:OE1	1:C:217:ASP:HB3	2.10	0.50
1:A:232:PHE:O	1:A:233:LYS:C	2.50	0.49
1:C:103:VAL:HG11	1:C:174:PHE:HE2	1.77	0.49
1:C:137:VAL:HG21	1:D:239:ARG:HA	1.94	0.49
1:B:234:ASN:HD22	1:B:236:LYS:H	1.60	0.48
1:A:206:VAL:O	1:A:213:SER:HA	2.13	0.48
1:A:45:SER:HB2	1:A:55:ILE:CD1	2.43	0.48
1:B:142:ILE:HD12	1:B:183:ARG:CZ	2.44	0.48
1:A:168:GLU:HG2	1:B:191:VAL:HA	1.96	0.47
1:C:186:TRP:O	1:C:187:LYS:HG2	2.13	0.47
1:B:234:ASN:C	1:B:234:ASN:HD22	2.17	0.47
1:A:102:ARG:HD3	1:A:229:ALA:HB2	1.96	0.46
1:C:202:GLU:O	1:C:203:LYS:C	2.54	0.46
1:A:2:ARG:NH1	1:A:39:ASN:O	2.48	0.46
1:C:84:PHE:CD1	1:C:84:PHE:C	2.89	0.45
1:D:134:VAL:CG1	1:D:139:VAL:HG11	2.46	0.45
1:A:80:SER:O	1:A:84:PHE:HB3	2.17	0.45
1:B:124:PRO:O	1:B:125:ALA:HB3	2.17	0.44
1:B:157:GLY:O	1:B:160:PHE:HB3	2.18	0.44
1:D:17:GLU:HG2	1:D:27:VAL:HG21	1.99	0.44
1:B:161:SER:OG	2:B:3076:NAD:N7N	2.51	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:197:ILE:CD1	1:D:197:ILE:N	2.82	0.43
1:D:168:GLU:HG2	1:D:170:TYR:H	1.82	0.43
1:B:61:ARG:CZ	1:B:61:ARG:HB3	2.48	0.43
2:D:3078:NAD:O2A	2:D:3078:NAD:H51N	2.18	0.43
1:D:72:ARG:NH2	1:D:247:SER:O	2.52	0.43
1:D:156:THR:O	2:D:3078:NAD:C2A	2.67	0.43
1:C:206:VAL:O	1:C:213:SER:HA	2.18	0.43
1:D:143:ARG:CG	1:D:143:ARG:HH11	2.31	0.43
1:D:234:ASN:HB3	1:D:237:ARG:HB2	2.00	0.42
1:A:3:ALA:HB2	1:A:42:PHE:CE2	2.54	0.42
1:D:152:GLN:H	1:D:152:GLN:NE2	2.17	0.42
1:A:130:VAL:HG13	1:A:199:VAL:HG13	2.01	0.41
1:A:177:ILE:HG13	1:A:178:PRO:HD2	2.02	0.41
1:B:152:GLN:HG2	1:B:153:ILE:N	2.35	0.41
1:D:134:VAL:HA	1:D:196:LYS:O	2.19	0.41
1:B:32:GLN:HB2	1:B:33:PRO:HD2	2.02	0.41
1:B:148:ILE:HD11	1:B:179:ILE:HD11	2.02	0.40
1:B:118:ALA:HB1	1:B:120:LEU:HD21	2.04	0.40
1:D:130:VAL:HG13	1:D:199:VAL:HG13	2.04	0.40
1:C:47:GLY:HA2	1:C:70:THR:HG22	2.03	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	247/278 (89%)	226 (92%)	19 (8%)	2 (1%)	22	52
1	B	247/278 (89%)	228 (92%)	17 (7%)	2 (1%)	22	52
1	C	247/278 (89%)	224 (91%)	23 (9%)	0	100	100
1	D	247/278 (89%)	225 (91%)	19 (8%)	3 (1%)	15	42
All	All	988/1112 (89%)	903 (91%)	78 (8%)	7 (1%)	25	56

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	11	GLY
1	A	71	GLY
1	B	11	GLY
1	D	64	PRO
1	D	128	ILE
1	B	47	GLY
1	D	11	GLY

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	213/235 (91%)	186 (87%)	27 (13%)	5	13
1	B	213/235 (91%)	175 (82%)	38 (18%)	2	5
1	C	213/235 (91%)	184 (86%)	29 (14%)	4	11
1	D	213/235 (91%)	180 (84%)	33 (16%)	3	8
All	All	852/940 (91%)	725 (85%)	127 (15%)	3	9

All (127) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	14	LYS
1	A	24	GLU
1	A	31	ASN
1	A	32	GLN
1	A	36	GLU
1	A	45	SER
1	A	80	SER
1	A	88	LEU
1	A	95	PHE
1	A	96	GLU
1	A	102	ARG
1	A	108	MET
1	A	111	VAL

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Mol	Chain	Res	Type
1	A	114	LEU
1	A	116	GLU
1	A	126	LYS
1	A	127	MET
1	A	133	ARG
1	A	134	VAL
1	A	158	TYR
1	A	168	GLU
1	A	192	SER
1	A	203	LYS
1	A	211	GLN
1	A	226	GLU
1	A	232	PHE
1	A	241	LEU
1	B	14	LYS
1	B	21	LYS
1	B	23	LEU
1	B	28	GLU
1	B	73	VAL
1	B	76	LEU
1	B	84	PHE
1	B	88	LEU
1	B	102	ARG
1	B	103	VAL
1	B	104	SER
1	B	106	SER
1	B	111	VAL
1	B	114	LEU
1	B	116	GLU
1	B	120	LEU
1	B	122	ARG
1	B	138	GLU
1	B	145	ASP
1	B	158	TYR
1	B	161	SER
1	B	166	VAL
1	B	192	SER
1	B	194	GLU
1	B	196	LYS
1	B	206	VAL
1	B	213	SER
1	B	216	PHE

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Mol	Chain	Res	Type
1	B	217	ASP
1	B	221	THR
1	B	225	SER
1	B	226	GLU
1	B	234	ASN
1	B	235	GLU
1	B	236	LYS
1	B	241	LEU
1	B	245	VAL
1	B	246	ARG
1	C	1	MET
1	C	2	ARG
1	C	26	GLU
1	C	60	LYS
1	C	72	ARG
1	C	75	LEU
1	C	76	LEU
1	C	84	PHE
1	C	86	VAL
1	C	88	LEU
1	C	99	ARG
1	C	105	CYS
1	C	111	VAL
1	C	114	LEU
1	C	116	GLU
1	C	123	LYS
1	C	126	LYS
1	C	133	ARG
1	C	138	GLU
1	C	143	ARG
1	C	158	TYR
1	C	179	ILE
1	C	206	VAL
1	C	212	LYS
1	C	213	SER
1	C	226	GLU
1	C	239	ARG
1	C	241	LEU
1	C	245	VAL
1	D	21	LYS
1	D	60	LYS
1	D	73	VAL

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Mol	Chain	Res	Type
1	D	80	SER
1	D	84	PHE
1	D	90	LYS
1	D	92	VAL
1	D	93	GLU
1	D	97	VAL
1	D	103	VAL
1	D	104	SER
1	D	105	CYS
1	D	108	MET
1	D	114	LEU
1	D	116	GLU
1	D	126	LYS
1	D	134	VAL
1	D	141	ARG
1	D	143	ARG
1	D	152	GLN
1	D	158	TYR
1	D	187	LYS
1	D	197	ILE
1	D	214	VAL
1	D	216	PHE
1	D	217	ASP
1	D	219	GLU
1	D	225	SER
1	D	227	PHE
1	D	235	GLU
1	D	241	LEU
1	D	245	VAL
1	D	246	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	31	ASN
1	A	78	HIS
1	B	78	HIS
1	B	83	ASN
1	B	234	ASN
1	C	39	ASN
1	D	32	GLN
1	D	152	GLN

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 ⓘ

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	NAD	A	3075	-	41,48,48	1.79	3 (7%)	43,73,73	2.00	6 (13%)
2	NAD	B	3076	-	41,48,48	1.77	4 (9%)	43,73,73	1.83	7 (16%)
2	NAD	C	3077	-	41,48,48	1.73	3 (7%)	43,73,73	1.98	5 (11%)
2	NAD	D	3078	-	41,48,48	1.77	4 (9%)	43,73,73	1.99	4 (9%)

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	3075	-	-	0/22/62/62	0/5/5/5
2	NAD	B	3076	-	-	0/22/62/62	0/5/5/5
2	NAD	C	3077	-	-	0/22/62/62	0/5/5/5
2	NAD	D	3078	-	-	0/22/62/62	0/5/5/5

All (14) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	3076	NAD	C2N-C3N	2.07	1.42	1.39
2	D	3078	NAD	O4B-C1B	2.16	1.44	1.41
2	D	3078	NAD	C2A-N1A	2.86	1.39	1.33
2	C	3077	NAD	C2A-N1A	3.16	1.39	1.33
2	B	3076	NAD	C2A-N1A	3.28	1.40	1.33
2	A	3075	NAD	C2A-N1A	3.32	1.40	1.33
2	D	3078	NAD	C2A-N3A	3.88	1.38	1.32
2	A	3075	NAD	C2A-N3A	4.02	1.38	1.32
2	C	3077	NAD	C2A-N3A	4.14	1.39	1.32
2	B	3076	NAD	C2A-N3A	4.66	1.39	1.32
2	B	3076	NAD	O7N-C7N	8.43	1.41	1.24
2	C	3077	NAD	O7N-C7N	8.73	1.42	1.24
2	D	3078	NAD	O7N-C7N	8.88	1.42	1.24
2	A	3075	NAD	O7N-C7N	9.03	1.42	1.24

All (22) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	3078	NAD	N3A-C2A-N1A	-10.96	119.31	128.86
2	C	3077	NAD	N3A-C2A-N1A	-10.73	119.51	128.86
2	A	3075	NAD	N3A-C2A-N1A	-10.25	119.93	128.86
2	B	3076	NAD	N3A-C2A-N1A	-9.18	120.86	128.86
2	C	3077	NAD	O7N-C7N-N7N	-3.47	117.65	122.58
2	A	3075	NAD	C1B-N9A-C4A	-3.47	120.65	126.64
2	B	3076	NAD	O7N-C7N-N7N	-3.11	118.16	122.58
2	B	3076	NAD	O5B-PA-O1A	-3.09	96.77	109.25
2	D	3078	NAD	C1B-N9A-C4A	-2.95	121.54	126.64
2	D	3078	NAD	O7N-C7N-N7N	-2.63	118.84	122.58
2	B	3076	NAD	C4A-C5A-N7A	-2.47	107.03	109.41
2	A	3075	NAD	O7N-C7N-N7N	-2.44	119.11	122.58
2	C	3077	NAD	C1B-N9A-C4A	-2.14	122.93	126.64
2	B	3076	NAD	C1B-N9A-C4A	-2.01	123.16	126.64
2	A	3075	NAD	N6A-C6A-N1A	2.07	122.87	118.77
2	D	3078	NAD	O2A-PA-O5B	2.12	118.16	108.14
2	B	3076	NAD	O7N-C7N-C3N	2.13	122.11	119.62
2	C	3077	NAD	O2N-PN-O1N	2.29	124.12	112.28
2	A	3075	NAD	O2N-PN-O1N	2.39	124.66	112.28
2	B	3076	NAD	O2A-PA-O5B	2.43	119.64	108.14
2	C	3077	NAD	C3N-C7N-N7N	2.58	120.72	117.77
2	A	3075	NAD	C3N-C7N-N7N	3.23	121.46	117.77

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	3076	NAD	1	0
2	D	3078	NAD	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

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	249/278 (89%)	0.23	9 (3%) 43 37	34, 38, 41, 44	0
1	B	249/278 (89%)	0.12	2 (0%) 86 85	34, 38, 41, 47	0
1	C	249/278 (89%)	0.18	10 (4%) 39 33	34, 39, 41, 44	0
1	D	249/278 (89%)	0.15	8 (3%) 48 42	35, 38, 41, 45	0
All	All	996/1112 (89%)	0.17	29 (2%) 52 46	34, 38, 41, 47	0

All (29) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	71	GLY	5.0
1	B	72	ARG	3.5
1	D	249	GLY	3.4
1	C	14	LYS	3.0
1	A	72	ARG	3.0
1	C	9	THR	2.9
1	C	8	LYS	2.9
1	A	8	LYS	2.7
1	D	44	VAL	2.6
1	C	72	ARG	2.6
1	A	14	LYS	2.5
1	A	249	GLY	2.4
1	A	73	VAL	2.3
1	A	43	ILE	2.3
1	C	24	GLU	2.3
1	C	82	GLU	2.3
1	C	94	LYS	2.3
1	C	96	GLU	2.3
1	D	2	ARG	2.3
1	D	1	MET	2.2
1	D	72	ARG	2.2

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Mol	Chain	Res	Type	RSRZ
1	B	44	VAL	2.2
1	A	9	THR	2.2
1	C	236	LYS	2.2
1	D	24	GLU	2.1
1	D	110	ASP	2.1
1	A	150	ALA	2.1
1	C	123	LYS	2.1
1	D	93	GLU	2.1

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

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

6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

6.4 Ligands [i](#)

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	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
2	NAD	A	3075	44/44	0.88	0.23	0.32	75,79,89,92	0
2	NAD	C	3077	44/44	0.87	0.22	0.25	74,79,89,90	0
2	NAD	D	3078	44/44	0.88	0.22	0.06	67,73,86,88	0
2	NAD	B	3076	44/44	0.90	0.21	-0.26	67,73,80,81	0

6.5 Other polymers [i](#)

There are no such residues in this entry.