



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 03:15 AM UTC

PDB ID : 2I1W / pdb_00002i1w
Title : Crystal structure of NAD kinase 1 from *Listeria monocytogenes*
Authors : Poncet-Montange, G.; Assairi, L.; Arold, S.; Pochet, S.; Labesse, G.
Deposited on : 2006-08-15
Resolution : 2.34 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 types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity	:	4-5-2 with Phenix2.0
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

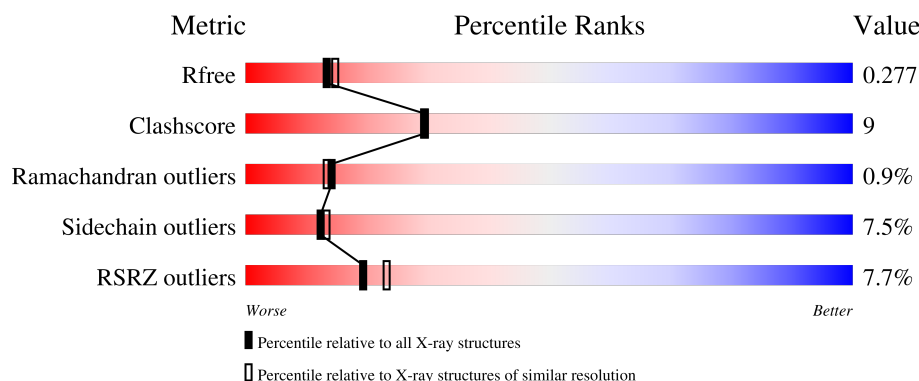
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.34 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}	180053	3031 (2.36-2.32)
Clashscore	190562	3127 (2.36-2.32)
Ramachandran outliers	187476	3095 (2.36-2.32)
Sidechain outliers	187428	3095 (2.36-2.32)
RSRZ outliers	180081	3033 (2.36-2.32)

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 of 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 $\leq 5\%$. 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	272	5% 69% 22% 6%
1	B	272	9% 81% 12% 6%
1	C	272	10% 78% 15% 5%
1	D	272	5% 73% 18% 5%

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
2	IOD	A	513	-	-	X	-
2	IOD	A	531	-	-	X	-
2	IOD	B	515	-	-	X	-
2	IOD	B	522	-	-	X	-

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 8538 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 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	257	Total	C	N	O	S	0	0	0
			2040	1309	341	381	9			
1	B	262	Total	C	N	O	S	0	0	0
			2069	1328	345	387	9			
1	C	258	Total	C	N	O	S	0	0	0
			2032	1305	341	377	9			
1	D	259	Total	C	N	O	S	0	0	0
			2016	1294	341	372	9			

There are 32 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	265	LEU	-	cloning artifact	UNP Q8Y8D7
A	266	GLU	-	cloning artifact	UNP Q8Y8D7
A	267	HIS	-	expression tag	UNP Q8Y8D7
A	268	HIS	-	expression tag	UNP Q8Y8D7
A	269	HIS	-	expression tag	UNP Q8Y8D7
A	270	HIS	-	expression tag	UNP Q8Y8D7
A	271	HIS	-	expression tag	UNP Q8Y8D7
A	272	HIS	-	expression tag	UNP Q8Y8D7
B	265	LEU	-	cloning artifact	UNP Q8Y8D7
B	266	GLU	-	cloning artifact	UNP Q8Y8D7
B	267	HIS	-	expression tag	UNP Q8Y8D7
B	268	HIS	-	expression tag	UNP Q8Y8D7
B	269	HIS	-	expression tag	UNP Q8Y8D7
B	270	HIS	-	expression tag	UNP Q8Y8D7
B	271	HIS	-	expression tag	UNP Q8Y8D7
B	272	HIS	-	expression tag	UNP Q8Y8D7
C	265	LEU	-	cloning artifact	UNP Q8Y8D7
C	266	GLU	-	cloning artifact	UNP Q8Y8D7
C	267	HIS	-	expression tag	UNP Q8Y8D7
C	268	HIS	-	expression tag	UNP Q8Y8D7
C	269	HIS	-	expression tag	UNP Q8Y8D7

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Chain	Residue	Modelled	Actual	Comment	Reference
C	270	HIS	-	expression tag	UNP Q8Y8D7
C	271	HIS	-	expression tag	UNP Q8Y8D7
C	272	HIS	-	expression tag	UNP Q8Y8D7
D	265	LEU	-	cloning artifact	UNP Q8Y8D7
D	266	GLU	-	cloning artifact	UNP Q8Y8D7
D	267	HIS	-	expression tag	UNP Q8Y8D7
D	268	HIS	-	expression tag	UNP Q8Y8D7
D	269	HIS	-	expression tag	UNP Q8Y8D7
D	270	HIS	-	expression tag	UNP Q8Y8D7
D	271	HIS	-	expression tag	UNP Q8Y8D7
D	272	HIS	-	expression tag	UNP Q8Y8D7

- Molecule 2 is IODIDE ION (CCD ID: IOD) (formula: I).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	8	Total I 9 9	0	1
2	B	5	Total I 6 6	0	1
2	C	2	Total I 3 3	0	1
2	D	3	Total I 4 4	0	1

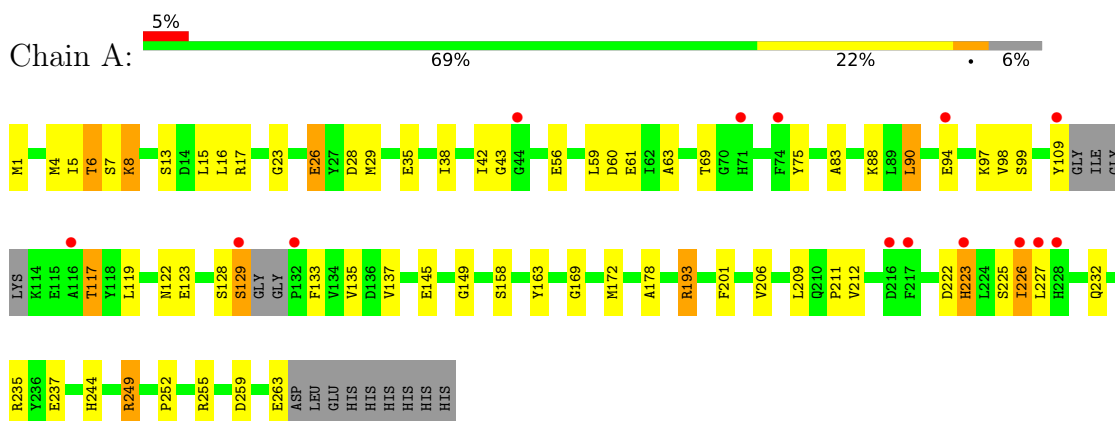
- Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	101	Total O 101 101	0	0
3	B	91	Total O 91 91	0	0
3	C	76	Total O 76 76	0	0
3	D	91	Total O 91 91	0	0

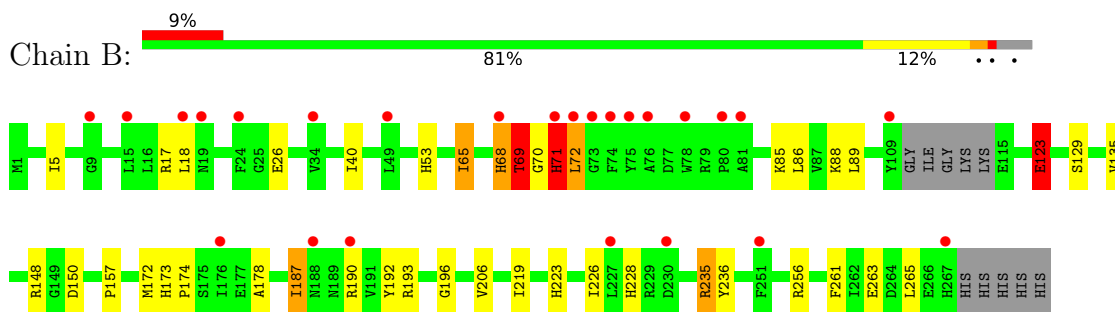
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide 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.

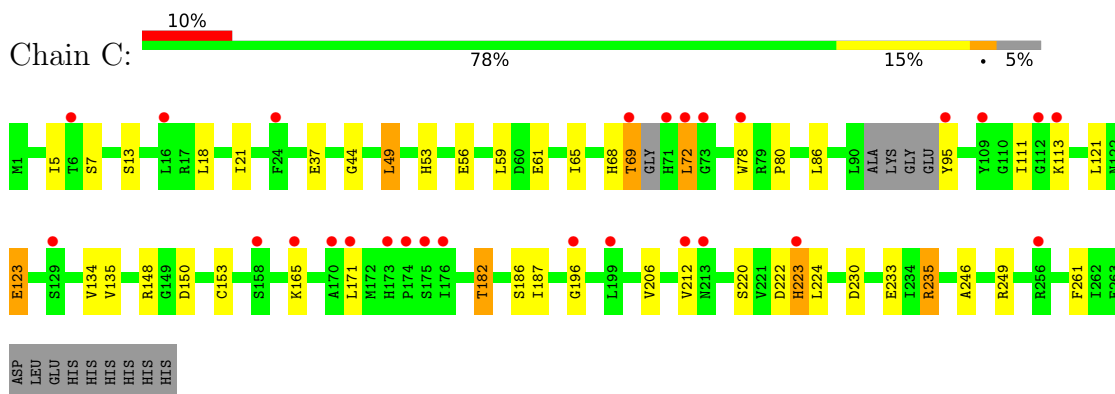
- Molecule 1: Probable inorganic polyphosphate/ATP-NAD kinase 1



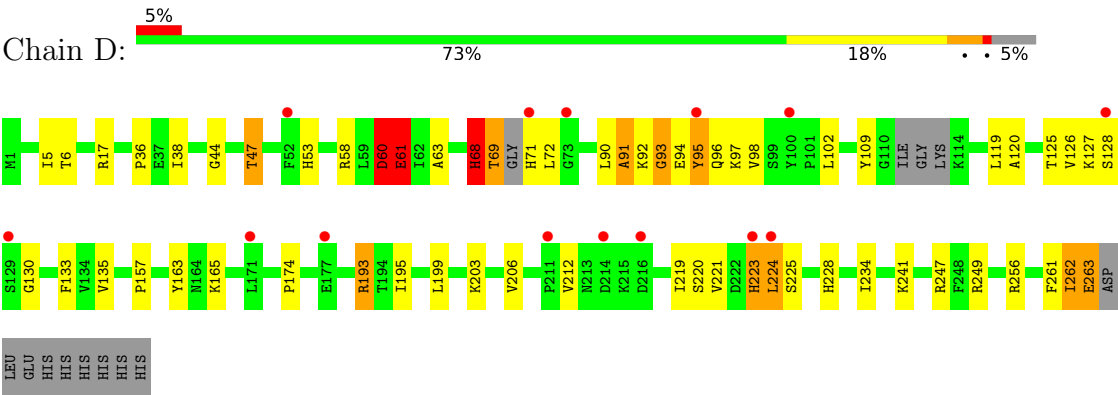
- Molecule 1: Probable inorganic polyphosphate/ATP-NAD kinase 1



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



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



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	66.61Å 119.02Å 67.94Å 90.00° 102.08° 90.00°	Depositor
Resolution (Å)	34.37 – 2.34 34.37 – 2.34	Depositor EDS
% Data completeness (in resolution range)	97.5 (34.37-2.34) 97.5 (34.37-2.34)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.38 (at 2.34Å)	Xtriage
Refinement program	REFMAC 5.2.0005	Depositor
R, R_{free}	0.194 , 0.257 0.219 , 0.277	Depositor DCC
R_{free} test set	1330 reflections (3.04%)	wwPDB-VP
Wilson B-factor (Å ²)	38.9	Xtriage
Anisotropy	0.296	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 70.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.35$	Xtriage
Estimated twinning fraction	0.008 for l,-k,h	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	8538	wwPDB-VP
Average B, all atoms (Å ²)	40.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.49% of the height of the origin peak. No significant pseudotranslation is detected.*

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: IOD

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	0.87	3/2088 (0.1%)	0.98	3/2820 (0.1%)
1	B	0.81	3/2120 (0.1%)	0.92	1/2870 (0.0%)
1	C	0.70	0/2081	0.89	0/2816
1	D	0.79	3/2063 (0.1%)	0.94	2/2788 (0.1%)
All	All	0.80	9/8352 (0.1%)	0.93	6/11294 (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
1	C	1	0
1	D	0	6
All	All	1	9

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	94	GLU	N-CA	6.86	1.56	1.46
1	B	18	LEU	C-O	6.78	1.32	1.24
1	D	68	HIS	CA-C	6.74	1.61	1.52
1	A	263	GLU	C-O	5.59	1.34	1.23
1	B	26	GLU	CD-OE2	5.57	1.35	1.25
1	A	255	ARG	C-O	5.47	1.30	1.24
1	A	259	ASP	CB-CG	5.41	1.65	1.52
1	D	93	GLY	CA-C	5.28	1.59	1.51
1	B	26	GLU	CD-OE1	5.09	1.35	1.25

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	223	HIS	N-CA-C	8.20	128.27	110.80
1	A	15	LEU	N-CA-C	8.19	121.32	111.82
1	B	123	GLU	CB-CA-C	-5.77	99.86	111.17
1	D	262	ILE	N-CA-C	5.42	120.62	109.34
1	A	169	GLY	N-CA-C	5.30	118.02	111.93
1	D	94	GLU	N-CA-C	-5.18	96.94	107.41

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
1	C	95	TYR	CA

All (9) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	222	ASP	Peptide
1	B	68	HIS	Peptide
1	B	69	THR	Peptide
1	D	262	ILE	Peptide
1	D	60	ASP	Peptide
1	D	68	HIS	Peptide
1	D	92	LYS	Peptide
1	D	93	GLY	Peptide
1	D	95	TYR	Peptide

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	2040	0	1983	44	0
1	B	2069	0	1997	37	0
1	C	2032	0	1961	30	0
1	D	2016	0	1948	31	0
2	A	9	0	0	9	0
2	B	6	0	0	6	0
2	C	3	0	0	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	D	4	0	0	1	0
3	A	101	0	0	4	1
3	B	91	0	0	7	4
3	C	76	0	0	3	2
3	D	91	0	0	1	3
All	All	8538	0	7889	138	5

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 9.

All (138) 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)
2:A:530:IOD:I	3:A:596:HOH:O	2.29	1.19
2:A:531:IOD:I	3:A:535:HOH:O	2.43	1.06
1:C:196:GLY:O	2:C:523:IOD:I	2.56	0.94
1:D:69:THR:O	1:D:71:HIS:N	2.00	0.93
1:B:256:ARG:HD3	3:B:559:HOH:O	1.70	0.91
1:A:225:SER:C	1:A:226:ILE:HD13	1.97	0.88
1:A:128:SER:O	1:A:129:SER:HB3	1.71	0.88
1:B:196:GLY:O	2:B:528:IOD:I	2.63	0.87
1:A:6:THR:HG21	2:A:513:IOD:I	2.44	0.87
1:B:265:LEU:HG	2:B:522:IOD:I	2.49	0.82
1:A:117:THR:CG2	2:A:531:IOD:I	3.01	0.79
1:C:235:ARG:HH11	1:C:235:ARG:HB2	1.47	0.78
1:D:58:ARG:O	1:D:61:GLU:HB3	1.89	0.73
1:C:153:CYS:HB3	1:C:182:THR:HG22	1.70	0.72
1:D:60:ASP:OD1	3:D:594:HOH:O	2.09	0.70
1:B:193:ARG:NH2	1:D:261:PHE:O	2.25	0.70
1:A:193:ARG:NH2	1:C:261:PHE:O	2.25	0.69
1:C:235:ARG:HH11	1:C:235:ARG:CG	2.06	0.68
1:D:38:ILE:HD13	1:D:90:LEU:HD22	1.75	0.68
1:C:235:ARG:HH11	1:C:235:ARG:CB	2.06	0.68
1:A:63:ALA:HB2	1:A:97:LYS:HE2	1.76	0.68
1:B:235:ARG:HH11	1:B:235:ARG:HG3	1.59	0.68
1:D:247:ARG:NH1	2:D:510:IOD:I	2.98	0.67
1:A:158:SER:O	3:A:586:HOH:O	2.14	0.65
1:C:53:HIS:HE1	1:C:222:ASP:OD2	1.80	0.64
1:B:190:ARG:CB	1:D:263:GLU:HB2	2.28	0.63
1:A:117:THR:HG22	2:A:531:IOD:I	2.69	0.63
1:A:7:SER:HB2	1:A:13:SER:HB3	1.81	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:69:THR:HA	1:C:80:PRO:HG3	1.83	0.61
1:A:226:ILE:HD13	1:A:226:ILE:N	2.16	0.60
1:C:235:ARG:HB2	1:C:235:ARG:NH1	2.15	0.60
1:B:88:LYS:HG3	1:B:89:LEU:N	2.18	0.59
1:B:123:GLU:CG	3:B:596:HOH:O	2.50	0.59
1:A:249:ARG:HH12	1:A:252:PRO:HD3	1.67	0.58
1:C:49:LEU:HD12	1:C:121:LEU:HD23	1.85	0.58
1:A:4:MET:HE1	1:A:6:THR:HG23	1.87	0.57
1:A:193:ARG:HH11	1:A:193:ARG:HG3	1.70	0.57
1:D:126:VAL:HG11	1:D:234:ILE:HD13	1.85	0.57
1:A:193:ARG:HH11	1:A:193:ARG:CG	2.17	0.57
1:B:261:PHE:O	1:D:193:ARG:NH2	2.35	0.57
1:D:126:VAL:HG12	1:D:219:ILE:HG13	1.86	0.56
1:B:172:MET:HE1	1:B:178:ALA:HB3	1.88	0.56
1:D:128:SER:HB2	1:D:133:PHE:HB2	1.88	0.56
1:D:63:ALA:HB2	1:D:97:LYS:HE2	1.87	0.55
1:B:68:HIS:HB3	1:B:72:LEU:HA	1.88	0.55
1:C:235:ARG:HH11	1:C:235:ARG:HG3	1.71	0.55
1:A:29:MET:HE1	1:A:90:LEU:HD13	1.89	0.54
1:A:99:SER:OG	1:A:244:HIS:ND1	2.41	0.54
1:A:249:ARG:HH11	1:A:249:ARG:CG	2.20	0.54
1:C:134:VAL:HG22	1:C:148:ARG:HG3	1.88	0.54
1:A:128:SER:HB2	1:A:133:PHE:HB2	1.90	0.54
1:A:8:LYS:HG2	1:A:43:GLY:HA3	1.89	0.54
1:A:7:SER:HA	1:A:42:ILE:O	2.07	0.54
1:C:65:ILE:HG12	1:C:246:ALA:HB3	1.91	0.53
1:B:123:GLU:HG3	3:B:596:HOH:O	2.07	0.53
1:A:99:SER:OG	1:A:244:HIS:CE1	2.61	0.53
1:D:219:ILE:HD13	1:D:228:HIS:CD2	2.44	0.53
1:B:263:GLU:HG3	2:B:522:IOD:I	2.79	0.52
1:B:265:LEU:CG	2:B:522:IOD:I	3.27	0.52
1:B:235:ARG:HH11	1:B:235:ARG:CG	2.23	0.51
1:A:172:MET:HE1	1:A:178:ALA:HB3	1.91	0.51
1:B:68:HIS:HD2	1:B:69:THR:H	1.59	0.51
1:A:28:ASP:O	1:A:29:MET:HG3	2.11	0.50
1:B:5:ILE:CG2	1:B:17:ARG:HG3	2.41	0.50
1:A:75:TYR:OH	1:A:122:ASN:ND2	2.44	0.50
1:A:117:THR:HG23	2:A:531:IOD:I	2.82	0.50
1:B:223:HIS:HB3	3:B:593:HOH:O	2.12	0.50
1:D:120:ALA:HB2	1:D:221:VAL:HG13	1.92	0.49
1:A:56:GLU:HA	1:A:59:LEU:HG	1.94	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:135:VAL:HG21	1:A:209:LEU:HB3	1.94	0.49
1:D:72:LEU:HD13	1:D:256:ARG:NH2	2.28	0.49
1:C:148:ARG:HD3	1:C:187:ILE:HD12	1.94	0.49
1:D:36:PRO:O	1:D:58:ARG:NH1	2.45	0.49
1:C:7:SER:HB2	1:C:13:SER:HB3	1.95	0.48
1:A:133:PHE:HB3	1:A:149:GLY:O	2.14	0.48
1:B:53:HIS:CD2	1:B:223:HIS:HE1	2.32	0.48
1:C:44:GLY:HA2	1:C:68:HIS:CD2	2.49	0.47
1:A:23:GLY:O	1:A:26:GLU:HG3	2.14	0.47
1:A:60:ASP:HB3	1:A:61:GLU:HG2	1.96	0.47
1:C:78:TRP:CE2	1:C:86:LEU:HD13	2.49	0.47
1:A:163:TYR:OH	1:C:150:ASP:OD2	2.32	0.47
1:D:44:GLY:H	1:D:47:THR:HG23	1.81	0.46
1:D:95:TYR:C	1:D:95:TYR:CD1	2.93	0.46
1:B:65:ILE:CD1	1:B:86:LEU:HD21	2.46	0.46
1:A:193:ARG:CG	1:A:193:ARG:NH1	2.77	0.46
1:B:68:HIS:CB	1:B:72:LEU:HA	2.44	0.46
1:B:70:GLY:O	1:B:71:HIS:O	2.33	0.46
2:B:515:IOD:I	1:D:165:LYS:HD3	2.85	0.46
1:A:211:PRO:HD2	1:A:232:GLN:O	2.15	0.46
1:C:49:LEU:HB3	1:C:223:HIS:CE1	2.51	0.46
1:C:123:GLU:CG	3:C:526:HOH:O	2.65	0.45
1:D:53:HIS:CD2	1:D:223:HIS:CE1	3.04	0.45
1:D:125:THR:OG1	1:D:220:SER:OG	2.35	0.45
1:D:58:ARG:HG2	1:D:61:GLU:HG2	1.98	0.45
1:B:148:ARG:HD3	1:B:192:TYR:CE2	2.52	0.45
1:B:68:HIS:CG	1:B:72:LEU:HA	2.52	0.45
1:C:165:LYS:HG3	1:C:171:LEU:HD21	1.98	0.45
1:D:95:TYR:CD1	1:D:95:TYR:O	2.70	0.45
1:A:235:ARG:NH1	1:A:237:GLU:OE1	2.49	0.45
1:B:150:ASP:OD2	1:D:163:TYR:OH	2.31	0.45
1:B:88:LYS:HG3	1:B:89:LEU:H	1.81	0.45
1:D:90:LEU:O	1:D:91:ALA:C	2.60	0.45
1:B:173:HIS:HE1	3:B:573:HOH:O	2.00	0.45
1:C:5:ILE:HD12	1:C:21:ILE:HG13	2.00	0.44
1:B:71:HIS:HB2	1:B:72:LEU:H	1.67	0.44
1:C:123:GLU:HG2	3:C:526:HOH:O	2.17	0.44
1:C:78:TRP:CZ2	1:C:86:LEU:HD13	2.52	0.43
1:D:157:PRO:HB2	1:D:174:PRO:HA	1.99	0.43
1:B:173:HIS:CE1	3:B:573:HOH:O	2.71	0.43
1:D:195:ILE:HG21	1:D:199:LEU:HD21	1.99	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:109:TYR:C	3:A:635:HOH:O	2.62	0.43
1:C:235:ARG:CG	1:C:235:ARG:NH1	2.75	0.43
1:C:56:GLU:HA	1:C:59:LEU:HG	2.01	0.42
1:A:137:VAL:HB	1:A:145:GLU:HG2	2.01	0.42
1:B:40:ILE:HG12	1:B:65:ILE:HG23	2.01	0.42
1:D:109:TYR:HH	1:D:228:HIS:CE1	2.35	0.42
1:A:16:LEU:HD11	1:A:83:ALA:CB	2.49	0.42
1:C:53:HIS:CE1	1:C:222:ASP:OD2	2.67	0.42
1:B:5:ILE:HG21	1:B:17:ARG:HG3	2.02	0.42
1:C:111:ILE:C	1:C:113:LYS:H	2.28	0.42
1:B:85:LYS:HG2	1:B:88:LYS:HE2	2.01	0.41
1:D:5:ILE:O	1:D:17:ARG:HD2	2.20	0.41
1:B:219:ILE:HB	1:B:228:HIS:CD2	2.54	0.41
1:A:178:ALA:HA	1:A:201:PHE:O	2.21	0.41
1:A:225:SER:O	1:A:226:ILE:HD13	2.17	0.41
1:D:220:SER:HA	1:D:224:LEU:O	2.20	0.41
1:C:233:GLU:CD	1:C:235:ARG:HH12	2.29	0.41
1:B:236:TYR:OH	3:B:588:HOH:O	2.22	0.41
1:D:119:LEU:HD11	1:D:241:LYS:HD3	2.03	0.41
1:A:119:LEU:HG	2:A:531:IOD:I	2.91	0.41
1:B:187:ILE:HA	2:B:515:IOD:I	2.91	0.41
1:A:5:ILE:HB	1:A:17:ARG:HG3	2.03	0.40
1:A:38:ILE:HG21	1:A:90:LEU:HD11	2.02	0.40
1:A:98:VAL:HG11	2:A:533:IOD:I	2.91	0.40
1:B:157:PRO:HB2	1:B:174:PRO:HA	2.02	0.40
1:A:6:THR:CG2	2:A:513:IOD:I	3.30	0.40
1:B:68:HIS:HD2	1:B:69:THR:N	2.20	0.40
1:C:182:THR:HB	3:C:533:HOH:O	2.21	0.40

All (5) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:B:611:HOH:O	3:C:584:HOH:O[1_556]	1.34	0.86
3:B:604:HOH:O	3:D:589:HOH:O[1_556]	1.63	0.57
3:B:613:HOH:O	3:D:586:HOH:O[1_556]	2.05	0.15
3:A:588:HOH:O	3:C:580:HOH:O[1_556]	2.11	0.09
3:B:612:HOH:O	3:D:587:HOH:O[1_556]	2.17	0.03

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	251/272 (92%)	238 (95%)	12 (5%)	1 (0%)	30	32
1	B	258/272 (95%)	243 (94%)	13 (5%)	2 (1%)	16	16
1	C	252/272 (93%)	240 (95%)	11 (4%)	1 (0%)	30	32
1	D	253/272 (93%)	231 (91%)	17 (7%)	5 (2%)	6	3
All	All	1014/1088 (93%)	952 (94%)	53 (5%)	9 (1%)	14	13

All (9) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	223	HIS
1	B	71	HIS
1	C	72	LEU
1	D	61	GLU
1	D	68	HIS
1	D	91	ALA
1	D	96	GLN
1	B	129	SER
1	D	130	GLY

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	219/237 (92%)	201 (92%)	18 (8%)	10	11
1	B	221/237 (93%)	211 (96%)	10 (4%)	24	32

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	C	216/237 (91%)	197 (91%)	19 (9%)	9	9
1	D	212/237 (90%)	194 (92%)	18 (8%)	10	10
All	All	868/948 (92%)	803 (92%)	65 (8%)	12	13

All (65) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	MET
1	A	6	THR
1	A	8	LYS
1	A	26	GLU
1	A	35	GLU
1	A	69	THR
1	A	88	LYS
1	A	90	LEU
1	A	94	GLU
1	A	117	THR
1	A	123	GLU
1	A	129	SER
1	A	193	ARG
1	A	206	VAL
1	A	212	VAL
1	A	226	ILE
1	A	227	LEU
1	A	249	ARG
1	B	65	ILE
1	B	69	THR
1	B	71	HIS
1	B	72	LEU
1	B	123	GLU
1	B	135	VAL
1	B	187	ILE
1	B	206	VAL
1	B	226	ILE
1	B	235	ARG
1	C	18	LEU
1	C	37	GLU
1	C	49	LEU
1	C	61	GLU
1	C	69	THR
1	C	72	LEU

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Mol	Chain	Res	Type
1	C	95	TYR
1	C	123	GLU
1	C	135	VAL
1	C	182	THR
1	C	186	SER
1	C	206	VAL
1	C	212	VAL
1	C	220	SER
1	C	223	HIS
1	C	224	LEU
1	C	230	ASP
1	C	235	ARG
1	C	249	ARG
1	D	6	THR
1	D	47	THR
1	D	60	ASP
1	D	61	GLU
1	D	69	THR
1	D	98	VAL
1	D	102	LEU
1	D	127	LYS
1	D	135	VAL
1	D	193	ARG
1	D	203	LYS
1	D	206	VAL
1	D	212	VAL
1	D	223	HIS
1	D	224	LEU
1	D	225	SER
1	D	249	ARG
1	D	263	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (14) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	19	ASN
1	A	122	ASN
1	A	210	GLN
1	B	173	HIS
1	B	210	GLN
1	B	223	HIS
1	C	53	HIS

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Mol	Chain	Res	Type
1	C	173	HIS
1	C	204	HIS
1	C	210	GLN
1	C	244	HIS
1	D	143	HIS
1	D	188	ASN
1	D	244	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 oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

Of 22 ligands modelled in this entry, 22 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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	257/272 (94%)	0.71	14 (5%)	31 37	28, 40, 51, 55	14 (5%)
1	B	262/272 (96%)	0.96	25 (9%)	14 17	23, 40, 48, 55	13 (4%)
1	C	258/272 (94%)	1.01	27 (10%)	11 13	32, 40, 47, 49	8 (3%)
1	D	259/272 (95%)	0.83	14 (5%)	31 37	31, 40, 47, 50	10 (3%)
All	All	1036/1088 (95%)	0.87	80 (7%)	19 23	23, 40, 48, 55	45 (4%)

All (80) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	223	HIS	4.0
1	A	217	PHE	3.6
1	A	227	LEU	3.5
1	B	230	ASP	3.4
1	C	109	TYR	3.3
1	B	81	ALA	3.2
1	B	73	GLY	3.2
1	C	223	HIS	3.2
1	D	100	TYR	3.1
1	C	171	LEU	3.1
1	A	71	HIS	3.1
1	D	129	SER	3.0
1	B	74	PHE	3.0
1	A	116	ALA	2.9
1	B	75	TYR	2.9
1	A	132	PRO	2.9
1	D	216	ASP	2.9
1	D	73	GLY	2.8
1	C	129	SER	2.7
1	A	129	SER	2.7
1	B	80	PRO	2.7

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Mol	Chain	Res	Type	RSRZ
1	A	109	TYR	2.6
1	C	95	TYR	2.6
1	C	71	HIS	2.6
1	B	71	HIS	2.6
1	B	18	LEU	2.6
1	B	227	LEU	2.6
1	C	73	GLY	2.6
1	A	44	GLY	2.5
1	D	128	SER	2.5
1	B	267	HIS	2.5
1	C	112	GLY	2.5
1	C	174	PRO	2.5
1	C	24	PHE	2.4
1	C	212	VAL	2.4
1	B	76	ALA	2.4
1	C	196	GLY	2.4
1	B	68	HIS	2.4
1	D	214	ASP	2.4
1	B	251	PHE	2.4
1	C	72	LEU	2.4
1	D	52	PHE	2.4
1	D	171	LEU	2.4
1	D	177	GLU	2.4
1	C	69	THR	2.3
1	B	24	PHE	2.3
1	B	109	TYR	2.3
1	A	223	HIS	2.3
1	D	71	HIS	2.3
1	B	49	LEU	2.3
1	B	190	ARG	2.3
1	C	78	TRP	2.3
1	B	19	ASN	2.2
1	C	213	ASN	2.2
1	D	211	PRO	2.2
1	C	165	LYS	2.2
1	A	226	ILE	2.2
1	B	188	ASN	2.2
1	C	113	LYS	2.2
1	C	158	SER	2.2
1	C	170	ALA	2.2
1	C	175	SER	2.2
1	A	216	ASP	2.1

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Mol	Chain	Res	Type	RSRZ
1	B	34	VAL	2.1
1	D	95	TYR	2.1
1	C	16	LEU	2.1
1	A	228	HIS	2.1
1	A	74	PHE	2.1
1	C	256	ARG	2.1
1	C	173	HIS	2.0
1	B	15	LEU	2.0
1	B	72	LEU	2.0
1	B	9	GLY	2.0
1	B	78	TRP	2.0
1	A	94	GLU	2.0
1	C	6	THR	2.0
1	C	199	LEU	2.0
1	D	224	LEU	2.0
1	B	176	ILE	2.0
1	C	176	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 oligosaccharides 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. 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	B-factors(Å ²)	Q<0.9
2	IOD	C	504[A]	1/1	0.90	0.16	33,33,33,33	1
2	IOD	C	504[B]	1/1	0.90	0.16	35,35,35,35	1
2	IOD	B	515	1/1	0.92	0.11	49,49,49,49	1
2	IOD	B	522	1/1	0.93	0.15	40,40,40,40	1
2	IOD	B	535	1/1	0.93	0.11	33,33,33,33	1
2	IOD	A	533	1/1	0.93	0.17	48,48,48,48	1

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	IOD	A	513	1/1	0.93	0.09	49,49,49,49	1
2	IOD	D	506[A]	1/1	0.93	0.15	33,33,33,33	1
2	IOD	D	506[B]	1/1	0.93	0.15	39,39,39,39	1
2	IOD	A	530	1/1	0.94	0.20	45,45,45,45	1
2	IOD	D	520	1/1	0.94	0.16	41,41,41,41	1
2	IOD	A	531	1/1	0.95	0.17	46,46,46,46	1
2	IOD	A	534	1/1	0.95	0.26	39,39,39,39	1
2	IOD	C	523	1/1	0.95	0.11	40,40,40,40	1
2	IOD	A	532	1/1	0.96	0.17	41,41,41,41	1
2	IOD	B	508[A]	1/1	0.96	0.16	37,37,37,37	1
2	IOD	D	510	1/1	0.96	0.24	40,40,40,40	1
2	IOD	B	508[B]	1/1	0.96	0.16	41,41,41,41	1
2	IOD	A	502[A]	1/1	0.97	0.13	43,43,43,43	1
2	IOD	A	502[B]	1/1	0.97	0.13	35,35,35,35	1
2	IOD	A	501	1/1	0.97	0.08	36,36,36,36	1
2	IOD	B	528	1/1	0.97	0.09	31,31,31,31	1

6.5 Other polymers

There are no such residues in this entry.