



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 28, 2026 – 10:31 PM UTC

PDB ID : 1OC4 / pdb_00001oc4
Title : Lactate dehydrogenase from Plasmodium berghei
Authors : Winter, V.J.; Brady, R.L.
Deposited on : 2003-02-05
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

<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
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
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

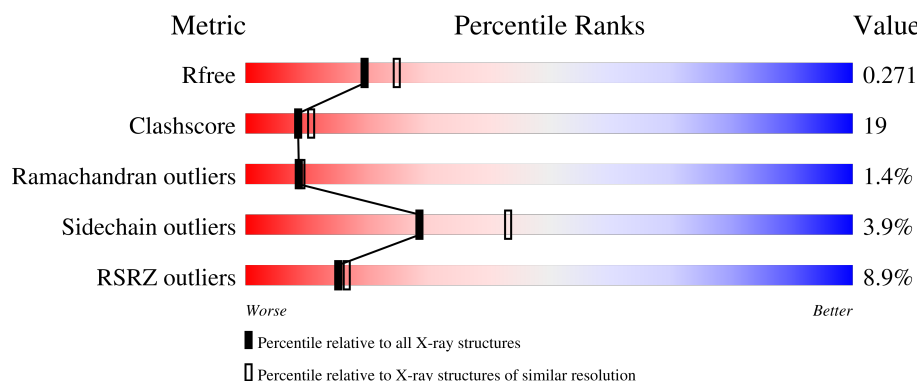
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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))
R_{free}	180053	6319 (2.30-2.30)
Clashscore	190562	6919 (2.30-2.30)
Ramachandran outliers	187476	6854 (2.30-2.30)
Sidechain outliers	187428	6854 (2.30-2.30)
RSRZ outliers	180081	6325 (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 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	322	
1	B	322	

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
4	GOL	B	1332	-	-	X	-

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 5115 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 L-LACTATE DEHYDROGENASE.

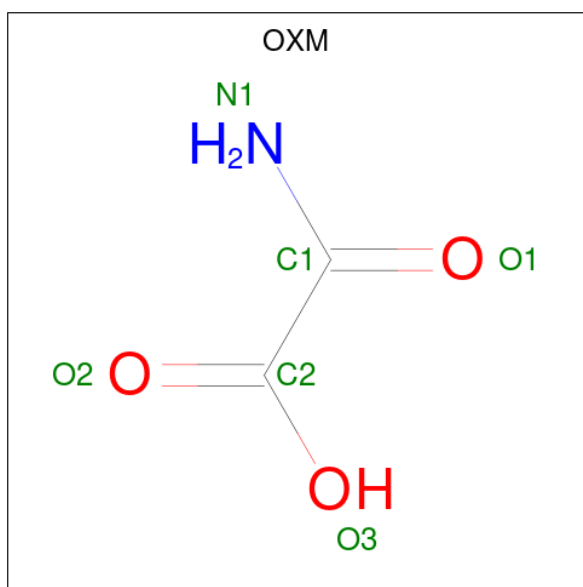
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	315	Total	C	N	O	S	57	0	0
			2404	1527	415	448	14			
1	B	315	Total	C	N	O	S	38	0	0
			2404	1527	415	448	14			

- Molecule 2 is NICOTINAMIDE-ADENINE-DINUCLEOTIDE (CCD ID: NAD) (formula: $C_{21}H_{27}N_7O_{14}P_2$).



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

- Molecule 3 is OXAMIC ACID (CCD ID: OXM) (formula: $C_2H_3NO_3$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			6	2	1	3		
3	B	1	Total	C	N	O	0	0
			6	2	1	3		

- Molecule 4 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			6	3	3		
4	B	1	Total	C	O	0	0
			6	3	3		

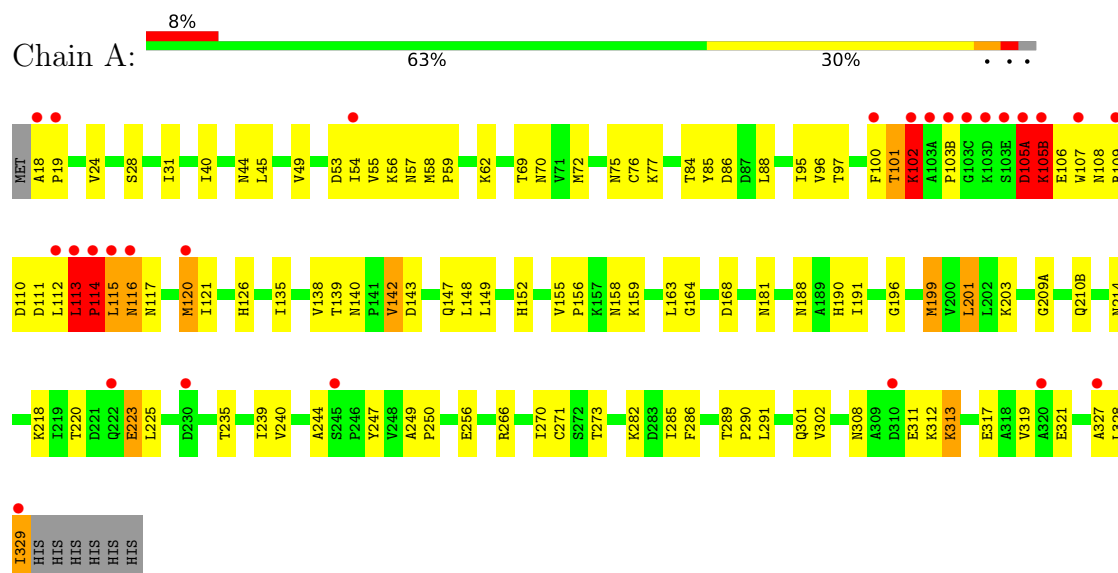
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	64	Total 64	O 64	0	0
5	B	131	Total 131	O 131	0	0

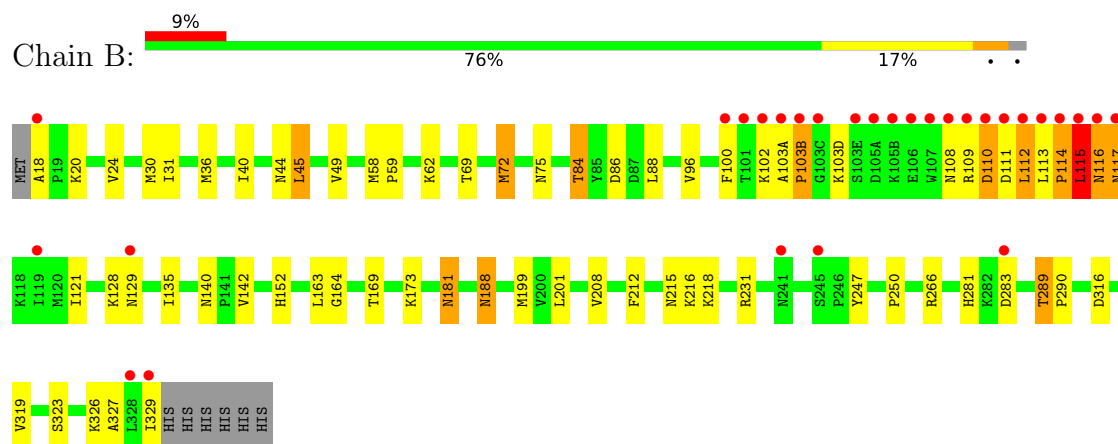
3 Residue-property plots [i](#)

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: L-LACTATE DEHYDROGENASE



• Molecule 1: L-LACTATE DEHYDROGENASE



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	103.59Å 71.53Å 94.54Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.49 – 2.30 29.49 – 2.30	Depositor EDS
% Data completeness (in resolution range)	95.0 (29.49-2.30) 94.9 (29.49-2.30)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.31 (at 2.29Å)	Xtriage
Refinement program	REFMAC 5.0	Depositor
R, R_{free}	0.219 , 0.268 0.218 , 0.271	Depositor DCC
R_{free} test set	1540 reflections (5.07%)	wwPDB-VP
Wilson B-factor (Å ²)	30.4	Xtriage
Anisotropy	0.101	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 40.3	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	5115	wwPDB-VP
Average B, all atoms (Å ²)	34.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.21% 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: OXM, GOL, 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	1.85	14/2441 (0.6%)	1.40	28/3305 (0.8%)
1	B	1.12	14/2441 (0.6%)	1.15	15/3305 (0.5%)
All	All	1.53	28/4882 (0.6%)	1.28	43/6610 (0.7%)

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	1	5

All (28) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	105(B)	LYS	CA-CB	-50.79	0.67	1.53
1	A	105(B)	LYS	N-CA	32.69	1.88	1.46
1	A	105(B)	LYS	CA-C	29.41	1.92	1.52
1	A	113	LEU	C-N	25.35	1.93	1.33
1	B	102	LYS	C-N	-24.15	1.08	1.33
1	A	102	LYS	N-CA	23.88	1.73	1.45
1	A	113	LEU	CA-C	21.96	1.79	1.52
1	A	101	THR	C-N	18.25	1.59	1.33
1	B	102	LYS	CA-CB	16.43	1.75	1.53
1	A	282	LYS	CB-CG	15.03	1.97	1.52
1	B	103(A)	ALA	CA-C	14.58	1.65	1.52
1	B	18	ALA	C-N	12.49	1.50	1.33
1	A	102	LYS	CA-CB	-11.88	1.34	1.53
1	B	112	LEU	CB-CG	10.96	1.75	1.53
1	B	103(B)	PRO	C-N	10.43	1.51	1.33
1	A	103(B)	PRO	N-CD	-9.24	1.34	1.47

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	77	LYS	CG-CD	-9.06	1.25	1.52
1	B	103(D)	LYS	CA-CB	-8.55	1.38	1.53
1	A	109	ARG	N-CA	-8.19	1.35	1.46
1	B	108	ASN	CB-CG	8.02	1.72	1.52
1	B	18	ALA	CA-C	7.92	1.69	1.52
1	B	18	ALA	N-CA	-7.18	1.32	1.46
1	A	105(A)	ASP	C-N	-7.12	1.17	1.34
1	B	109	ARG	C-N	-6.92	1.24	1.33
1	B	109	ARG	CA-C	6.59	1.62	1.52
1	A	106	GLU	N-CA	-6.50	1.38	1.46
1	B	72	MET	SD-CE	-5.05	1.67	1.79
1	B	115	LEU	C-O	-5.01	1.18	1.24

All (43) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	102	LYS	N-CA-CB	-21.05	79.14	111.35
1	A	105(B)	LYS	N-CA-CB	18.52	141.79	110.49
1	A	105(A)	ASP	O-C-N	-17.60	94.53	122.70
1	A	105(B)	LYS	N-CA-C	-17.39	73.76	110.80
1	A	105(B)	LYS	CB-CA-C	17.10	144.45	110.42
1	A	102	LYS	CB-CA-C	16.71	137.38	109.80
1	A	113	LEU	O-C-N	-15.97	106.27	120.48
1	A	103(B)	PRO	N-CA-CB	11.30	113.20	103.25
1	A	106	GLU	CA-C-N	-10.85	105.76	120.95
1	A	106	GLU	C-N-CA	-10.85	105.76	120.95
1	B	112	LEU	CB-CA-C	9.95	126.97	109.65
1	B	116	ASN	N-CA-C	9.54	125.28	112.30
1	A	106	GLU	N-CA-C	-8.96	102.50	112.72
1	A	113	LEU	CB-CG-CD2	8.77	137.01	110.70
1	B	102	LYS	CA-CB-CG	-8.13	97.83	114.10
1	A	106	GLU	O-C-N	7.97	132.16	122.35
1	A	113	LEU	C-N-CD	-7.23	95.34	125.00
1	B	142	VAL	N-CA-C	7.12	117.83	110.36
1	A	105(A)	ASP	CA-C-N	7.08	132.77	117.20
1	B	109	ARG	CA-C-O	-7.08	111.26	119.61
1	A	102	LYS	CA-C-O	-7.05	113.06	121.66
1	B	108	ASN	CA-CB-CG	6.91	119.51	112.60
1	B	18	ALA	N-CA-CB	6.86	120.69	110.40
1	A	103(B)	PRO	CA-N-CD	-6.52	102.87	112.00
1	A	49	VAL	N-CA-C	6.43	117.11	108.11
1	A	329	ILE	CB-CA-C	-6.22	99.17	111.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	109	ARG	O-C-N	6.19	130.38	122.33
1	A	199	MET	CB-CA-C	-6.10	101.23	109.71
1	A	218	LYS	N-CA-C	-6.08	106.02	113.50
1	A	103(B)	PRO	CB-CA-C	-6.08	103.84	111.56
1	B	110	ASP	N-CA-C	5.64	117.51	111.36
1	B	135	ILE	N-CA-C	5.59	115.94	108.11
1	A	135	ILE	N-CA-C	5.28	115.51	108.11
1	B	49	VAL	N-CA-C	5.22	115.42	108.11
1	A	49	VAL	CB-CA-C	-5.13	103.04	110.63
1	A	115	LEU	N-CA-CB	-5.09	104.32	110.53
1	B	84	THR	CB-CA-C	-5.09	100.39	109.71
1	A	256	GLU	N-CA-C	-5.09	105.73	111.28
1	A	107	TRP	N-CA-CB	5.08	117.42	109.85
1	A	142	VAL	N-CA-C	5.06	119.87	109.34
1	A	102	LYS	CA-CB-CG	-5.06	103.98	114.10
1	B	96	VAL	N-CA-C	5.05	115.04	107.51
1	B	112	LEU	N-CA-CB	-5.04	103.02	110.53

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
1	A	105(B)	LYS	CA

All (5) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	101	THR	Mainchain
1	A	102	LYS	Mainchain
1	A	105(A)	ASP	Peptide,Mainchain
1	A	105(B)	LYS	Mainchain

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	2404	0	2485	132	0
1	B	2404	0	2484	54	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	A	44	0	26	7	0
2	B	44	0	26	3	0
3	A	6	0	2	1	0
3	B	6	0	2	0	0
4	A	6	0	8	0	0
4	B	6	0	8	4	0
5	A	64	0	0	5	0
5	B	131	0	0	2	0
All	All	5115	0	5041	183	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 19.

All (183) 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:100:PHE:CD1	1:A:115:LEU:HG	1.35	1.54
1:A:112:LEU:C	1:A:114:PRO:HD2	1.33	1.48
1:A:100:PHE:CD1	1:A:115:LEU:CG	2.14	1.21
1:A:112:LEU:C	1:A:114:PRO:CD	2.16	1.17
1:A:113:LEU:N	1:A:114:PRO:CD	2.08	1.17
1:A:113:LEU:N	1:A:114:PRO:HD2	1.58	1.16
1:A:121:ILE:HD13	1:A:152:HIS:CE1	1.82	1.14
1:A:117:ASN:HD21	1:A:329:ILE:CD1	1.68	1.06
1:A:329:ILE:HG23	1:A:329:ILE:O	1.32	1.05
1:A:18:ALA:HB3	1:A:19:PRO:HD3	1.36	1.05
1:A:117:ASN:HD21	1:A:329:ILE:HD12	1.24	1.00
1:A:191:ILE:HG12	1:A:201:LEU:HD22	1.45	0.98
1:A:329:ILE:O	1:A:329:ILE:CG2	2.08	0.98
1:B:100:PHE:HD1	1:B:116:ASN:HD21	1.06	0.93
1:A:313:LYS:HE2	1:A:317:GLU:OE2	1.69	0.92
1:A:121:ILE:CD1	1:A:152:HIS:CE1	2.51	0.92
1:A:120:MET:HE3	1:A:120:MET:HA	1.55	0.89
1:B:121:ILE:HD12	1:B:152:HIS:NE2	1.87	0.89
1:B:121:ILE:HD12	1:B:152:HIS:CD2	2.09	0.87
1:A:220:THR:OG1	1:A:223:GLU:HG2	1.74	0.86
1:A:18:ALA:HB3	1:A:19:PRO:CD	2.05	0.86
1:A:121:ILE:HD13	1:A:152:HIS:NE2	1.94	0.82
1:A:84:THR:HG22	1:A:86:ASP:H	1.45	0.82
1:A:100:PHE:CG	1:A:115:LEU:HD23	2.17	0.80
1:A:113:LEU:N	1:A:114:PRO:HD3	1.93	0.80

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:100:PHE:CD1	1:A:115:LEU:CD2	2.65	0.80
1:A:120:MET:CE	1:A:149:LEU:HD12	2.12	0.80
1:B:100:PHE:HD1	1:B:116:ASN:ND2	1.79	0.79
1:B:69:THR:HA	1:B:72:MET:HE2	1.65	0.79
1:A:112:LEU:O	1:A:114:PRO:HD2	1.80	0.78
1:A:18:ALA:CB	1:A:19:PRO:HD3	2.14	0.78
1:A:117:ASN:ND2	1:A:329:ILE:CD1	2.47	0.78
1:A:100:PHE:HD1	1:A:115:LEU:O	1.66	0.78
1:B:121:ILE:CD1	1:B:152:HIS:CD2	2.67	0.77
1:A:155:VAL:CG2	5:A:2022:HOH:O	2.33	0.77
1:A:100:PHE:CG	1:A:115:LEU:CD2	2.66	0.77
1:A:285:ILE:HD13	1:A:319:VAL:HG13	1.65	0.77
1:A:117:ASN:HD21	1:A:329:ILE:HD13	1.50	0.75
1:B:100:PHE:CG	1:B:115:LEU:HD12	2.20	0.75
1:B:69:THR:HA	1:B:72:MET:CE	2.17	0.74
1:B:327:ALA:C	1:B:329:ILE:H	1.96	0.73
1:B:121:ILE:CD1	1:B:152:HIS:NE2	2.52	0.73
1:A:327:ALA:C	1:A:329:ILE:H	1.98	0.72
1:A:105(A):ASP:C	1:A:105(B):LYS:CA	2.61	0.72
1:A:273:THR:HG21	1:A:291:LEU:HD21	1.72	0.72
1:A:117:ASN:ND2	1:A:329:ILE:HD13	2.05	0.70
1:A:155:VAL:HG22	1:A:156:PRO:HD2	1.74	0.70
1:B:103(B):PRO:HD2	1:B:115:LEU:HD21	1.73	0.69
1:A:56:LYS:O	1:A:57:ASN:HB2	1.92	0.68
1:A:70:ASN:HD21	1:A:76:CYS:H	1.41	0.68
1:A:266:ARG:HH11	1:B:181:ASN:HD21	1.42	0.67
1:B:121:ILE:HD12	1:B:152:HIS:CE1	2.30	0.67
1:A:112:LEU:CA	1:A:114:PRO:CD	2.73	0.66
1:A:121:ILE:CD1	1:A:152:HIS:NE2	2.57	0.66
1:A:291:LEU:CD2	1:A:302:VAL:HG22	2.26	0.66
1:A:53:ASP:OD1	1:A:54:ILE:N	2.29	0.66
1:A:271:CYS:HB2	1:A:291:LEU:HD12	1.78	0.65
1:A:24:VAL:HG11	1:A:88:LEU:HD23	1.79	0.65
1:A:69:THR:HA	1:A:72:MET:HE2	1.80	0.64
1:A:120:MET:HB2	1:A:148:LEU:HD23	1.78	0.64
1:A:58:MET:HB3	1:A:59:PRO:HD3	1.78	0.64
1:A:308:ASN:OD1	1:A:308:ASN:C	2.40	0.64
1:A:313:LYS:CE	1:A:317:GLU:OE2	2.46	0.63
1:A:113:LEU:CA	1:A:114:PRO:CD	2.76	0.63
1:A:28:SER:O	1:A:62:LYS:NZ	2.25	0.63
1:A:115:LEU:O	1:A:116:ASN:CG	2.42	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:44:ASN:ND2	1:A:75:ASN:H	1.97	0.62
1:B:169:THR:HB	1:B:173:LYS:HE3	1.80	0.62
1:A:110:ASP:C	1:A:112:LEU:H	2.07	0.62
1:A:100:PHE:CD1	1:A:115:LEU:CB	2.82	0.62
1:B:215:ASN:O	1:B:216:LYS:HB2	1.99	0.61
1:A:181:ASN:HD21	1:B:266:ARG:NH1	2.00	0.60
1:A:142:VAL:HG13	1:A:143:ASP:N	2.16	0.59
1:A:321:GLU:HA	1:A:321:GLU:OE1	2.02	0.59
1:B:84:THR:HG22	1:B:86:ASP:H	1.67	0.59
1:A:140:ASN:ND2	2:A:1330:NAD:O2D	2.37	0.58
1:A:120:MET:HE1	1:A:149:LEU:HD12	1.85	0.58
1:B:113:LEU:HB2	1:B:114:PRO:HD3	1.84	0.57
1:A:113:LEU:CA	1:A:114:PRO:HD2	2.31	0.57
1:A:100:PHE:CD1	1:A:115:LEU:O	2.53	0.57
1:A:58:MET:N	1:A:59:PRO:HD2	2.20	0.56
1:A:112:LEU:CA	1:A:114:PRO:HD2	2.29	0.56
1:A:155:VAL:HG23	5:A:2022:HOH:O	1.99	0.56
1:A:220:THR:H	1:A:223:GLU:HG3	1.70	0.56
1:A:210(B):GLN:HE21	1:A:214:ASN:HD21	1.54	0.55
1:B:110:ASP:C	1:B:112:LEU:H	2.14	0.55
1:A:147:GLN:HB2	1:A:286:PHE:CD1	2.42	0.55
1:B:188:ASN:HD21	4:B:1332:GOL:H2	1.71	0.55
1:A:85:TYR:CG	1:A:126:HIS:ND1	2.76	0.54
1:A:273:THR:HB	1:A:291:LEU:HD11	1.89	0.54
1:B:113:LEU:O	1:B:117:ASN:ND2	2.41	0.54
1:B:44:ASN:ND2	1:B:75:ASN:H	2.06	0.53
1:A:120:MET:HE2	1:A:149:LEU:HD12	1.90	0.53
1:B:327:ALA:C	1:B:329:ILE:N	2.63	0.53
1:A:54:ILE:HG13	1:A:55:VAL:N	2.23	0.52
1:A:201:LEU:HD23	1:A:201:LEU:N	2.24	0.52
1:A:70:ASN:ND2	1:A:76:CYS:H	2.04	0.52
1:A:100:PHE:CD2	1:A:115:LEU:HD21	2.44	0.52
1:A:110:ASP:O	1:A:112:LEU:N	2.43	0.52
1:A:117:ASN:OD1	1:A:117:ASN:C	2.53	0.52
1:A:108:ASN:OD1	1:A:110:ASP:N	2.33	0.51
1:A:108:ASN:OD1	1:A:110:ASP:HB2	2.11	0.51
1:A:190:HIS:CD2	1:A:270:ILE:HD11	2.46	0.51
1:A:100:PHE:CD2	1:A:115:LEU:CD2	2.93	0.51
1:A:158:ASN:OD1	1:A:159:LYS:HG2	2.10	0.51
1:B:31:ILE:HG22	1:B:36:MET:HE2	1.93	0.51
1:A:285:ILE:HD13	1:A:319:VAL:CG1	2.40	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:220:THR:OG1	1:A:223:GLU:CG	2.56	0.50
1:A:239:ILE:HG22	1:A:244:ALA:O	2.11	0.50
1:A:289:THR:OG1	1:A:290:PRO:HD2	2.12	0.49
1:A:138:VAL:O	2:A:1330:NAD:H2N	2.13	0.49
1:A:155:VAL:CG2	1:A:156:PRO:HD2	2.42	0.49
1:A:291:LEU:HD23	1:A:302:VAL:HA	1.95	0.49
1:A:117:ASN:ND2	1:A:329:ILE:HG21	2.28	0.49
1:A:112:LEU:HA	1:A:114:PRO:CG	2.43	0.49
1:A:113:LEU:CA	1:A:114:PRO:N	2.76	0.49
1:B:58:MET:SD	1:B:62:LYS:HE3	2.53	0.49
1:B:140:ASN:OD1	2:B:1330:NAD:C2N	2.61	0.49
1:A:191:ILE:CG1	1:A:201:LEU:HD22	2.32	0.49
2:A:1330:NAD:C4N	3:A:1331:OXM:C1	2.91	0.48
1:B:84:THR:HG23	5:B:2013:HOH:O	2.12	0.48
1:B:247:TYR:HA	1:B:250:PRO:HG2	1.96	0.48
1:A:329:ILE:O	1:A:329:ILE:HG12	2.09	0.48
1:A:18:ALA:CB	1:A:19:PRO:CD	2.74	0.48
1:B:112:LEU:CG	1:B:112:LEU:CA	2.92	0.48
1:B:281:HIS:CD2	1:B:319:VAL:HG11	2.48	0.48
1:A:85:TYR:CD2	1:A:126:HIS:ND1	2.81	0.48
1:A:196:GLY:O	1:A:199:MET:HG3	2.13	0.48
1:B:327:ALA:O	1:B:329:ILE:HG23	2.14	0.47
1:A:142:VAL:CG1	1:A:143:ASP:N	2.77	0.47
1:A:181:ASN:HD21	1:B:266:ARG:HH11	1.61	0.47
1:A:203:LYS:NZ	5:A:2036:HOH:O	2.47	0.47
1:A:85:TYR:CD1	1:A:126:HIS:CE1	3.03	0.46
1:A:210(B):GLN:HE21	1:A:214:ASN:ND2	2.13	0.46
1:B:113:LEU:HB2	1:B:114:PRO:CD	2.46	0.46
1:A:247:TYR:C	1:A:250:PRO:HD2	2.40	0.46
1:B:30:MET:HE2	1:B:247:TYR:CZ	2.50	0.46
1:A:311:GLU:OE2	5:A:2061:HOH:O	2.21	0.46
1:A:110:ASP:C	1:A:112:LEU:N	2.73	0.46
1:B:283:ASP:O	1:B:326:LYS:NZ	2.48	0.46
1:A:249:ALA:N	1:A:250:PRO:CD	2.78	0.46
1:B:163:LEU:HD23	2:B:1330:NAD:C7N	2.46	0.46
1:B:100:PHE:CD1	1:B:116:ASN:ND2	2.70	0.45
1:A:327:ALA:C	1:A:329:ILE:N	2.67	0.45
1:B:208:VAL:HB	1:B:212:PHE:CE1	2.52	0.45
1:B:215:ASN:O	1:B:216:LYS:CB	2.62	0.45
1:A:58:MET:N	1:A:59:PRO:CD	2.79	0.44
1:A:285:ILE:HG12	1:A:286:PHE:N	2.32	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:44:ASN:HD21	1:A:75:ASN:H	1.65	0.44
1:A:266:ARG:NH1	1:B:181:ASN:HD21	2.11	0.44
1:B:128:LYS:HB2	1:B:128:LYS:HE3	1.85	0.44
1:A:301:GLN:NE2	1:B:218:LYS:HE2	2.33	0.44
1:A:113:LEU:H	1:A:114:PRO:HD3	1.80	0.44
1:A:54:ILE:HG23	2:A:1330:NAD:C5A	2.48	0.43
1:B:188:ASN:HD21	4:B:1332:GOL:C2	2.29	0.43
1:B:289:THR:OG1	1:B:290:PRO:HD2	2.17	0.43
1:A:163:LEU:HD23	2:A:1330:NAD:C7N	2.48	0.43
1:A:209(A):GLY:O	4:B:1332:GOL:H32	2.19	0.43
1:A:100:PHE:CD1	1:A:115:LEU:HD23	2.45	0.42
1:A:58:MET:HB3	1:A:59:PRO:CD	2.46	0.42
1:A:97:THR:HB	2:A:1330:NAD:H51N	2.00	0.42
1:B:113:LEU:O	1:B:117:ASN:CG	2.62	0.42
1:A:40:ILE:HG23	1:A:45:LEU:HB3	2.00	0.42
1:B:24:VAL:HG11	1:B:88:LEU:HD23	2.01	0.42
1:A:53:ASP:HB3	1:A:59:PRO:HG3	2.00	0.42
1:B:40:ILE:HG23	1:B:45:LEU:HB3	2.01	0.42
1:B:188:ASN:ND2	4:B:1332:GOL:H2	2.34	0.42
1:A:96:VAL:HG11	1:A:120:MET:HE1	2.02	0.42
1:A:112:LEU:HA	1:A:114:PRO:CD	2.46	0.41
1:A:138:VAL:O	1:A:139:THR:C	2.63	0.41
1:A:203:LYS:HE3	1:A:225:LEU:HD23	2.01	0.41
1:A:220:THR:H	1:A:223:GLU:CG	2.31	0.41
1:B:212:PHE:HD2	1:B:218:LYS:HD3	1.85	0.41
1:A:58:MET:SD	1:A:62:LYS:HD2	2.60	0.41
2:B:1330:NAD:H6N	5:B:2097:HOH:O	2.20	0.41
1:A:329:ILE:HD12	5:A:2021:HOH:O	2.20	0.41
1:A:168:ASP:OD1	1:A:199:MET:HE2	2.21	0.41
1:B:58:MET:N	1:B:59:PRO:HD2	2.36	0.41
1:A:31:ILE:HG13	2:A:1330:NAD:PN	2.61	0.41
1:B:20:LYS:HZ2	1:B:75:ASN:HB3	1.85	0.41
1:B:199:MET:HG3	1:B:201:LEU:CD1	2.50	0.41
1:B:281:HIS:CE1	1:B:316:ASP:OD1	2.74	0.40
1:B:181:ASN:HD22	1:B:181:ASN:HA	1.71	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	313/322 (97%)	292 (93%)	15 (5%)	6 (2%)	6	5
1	B	313/322 (97%)	298 (95%)	12 (4%)	3 (1%)	12	15
All	All	626/644 (97%)	590 (94%)	27 (4%)	9 (1%)	9	9

All (9) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	105(A)	ASP
1	A	114	PRO
1	A	116	ASN
1	A	111	ASP
1	A	328	LEU
1	B	111	ASP
1	A	164	GLY
1	B	164	GLY
1	B	114	PRO

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	266/273 (97%)	254 (96%)	12 (4%)	24	37
1	B	266/273 (97%)	257 (97%)	9 (3%)	32	49
All	All	532/546 (97%)	511 (96%)	21 (4%)	28	43

All (21) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	95	ILE
1	A	102	LYS
1	A	113	LEU
1	A	114	PRO
1	A	120	MET
1	A	188	ASN
1	A	201	LEU
1	A	223	GLU
1	A	235	THR
1	A	240	VAL
1	A	312	LYS
1	A	313	LYS
1	B	45	LEU
1	B	115	LEU
1	B	117	ASN
1	B	129	ASN
1	B	181	ASN
1	B	188	ASN
1	B	231	ARG
1	B	289	THR
1	B	323	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (23) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	44	ASN
1	A	57	ASN
1	A	70	ASN
1	A	75	ASN
1	A	117	ASN
1	A	140	ASN
1	A	178	GLN
1	A	181	ASN
1	A	190	HIS
1	A	210(B)	GLN
1	A	214	ASN
1	B	44	ASN
1	B	117	ASN
1	B	126	HIS
1	B	129	ASN
1	B	178	GLN

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Mol	Chain	Res	Type
1	B	181	ASN
1	B	188	ASN
1	B	210(B)	GLN
1	B	222	GLN
1	B	234	ASN
1	B	241	ASN
1	B	281	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 ⓘ

6 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	B	1330	-	46,48,48	1.68	4 (8%)	64,73,73	1.68	12 (18%)
3	OXM	A	1331	-	5,5,5	3.06	2 (40%)	2,6,6	0.67	0
3	OXM	B	1331	-	5,5,5	2.71	1 (20%)	2,6,6	0.78	0
4	GOL	A	1332	-	5,5,5	0.21	0	5,5,5	0.67	0
4	GOL	B	1332	-	5,5,5	0.45	0	5,5,5	0.83	0
2	NAD	A	1330	-	46,48,48	1.65	4 (8%)	64,73,73	1.76	12 (18%)

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	B	1330	-	-	6/30/62/62	0/5/5/5
3	OXM	A	1331	-	-	1/4/4/4	-
3	OXM	B	1331	-	-	4/4/4/4	-
4	GOL	A	1332	-	-	0/4/4/4	-
4	GOL	B	1332	-	-	2/4/4/4	-
2	NAD	A	1330	-	-	7/30/62/62	0/5/5/5

All (11) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	1330	NAD	O7N-C7N	8.96	1.40	1.24
2	B	1330	NAD	O7N-C7N	8.94	1.40	1.24
3	A	1331	OXM	C1-C2	-6.38	1.47	1.55
3	B	1331	OXM	C1-C2	-5.67	1.48	1.55
2	B	1330	NAD	C8A-N7A	2.82	1.37	1.31
2	A	1330	NAD	C8A-N7A	2.60	1.36	1.31
2	A	1330	NAD	C2A-N1A	2.55	1.38	1.33
2	B	1330	NAD	C2A-N3A	2.48	1.38	1.33
2	A	1330	NAD	C2A-N3A	2.37	1.38	1.33
2	B	1330	NAD	C2A-N1A	2.31	1.38	1.33
3	A	1331	OXM	O3-C2	-2.19	1.24	1.30

All (24) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1330	NAD	N3A-C2A-N1A	-5.57	120.15	128.58
2	B	1330	NAD	N3A-C2A-N1A	-5.50	120.25	128.58
2	A	1330	NAD	C5A-C4A-N3A	-4.46	120.57	126.72
2	A	1330	NAD	N9A-C8A-N7A	-4.17	108.02	113.94
2	B	1330	NAD	O2A-PA-O3	4.02	118.13	107.27
2	A	1330	NAD	C5A-N7A-C8A	3.92	109.61	103.45
2	B	1330	NAD	N9A-C8A-N7A	-3.92	108.38	113.94
2	B	1330	NAD	C5A-C4A-N3A	-3.70	121.62	126.72
2	A	1330	NAD	C2A-N3A-C4A	3.41	120.16	111.83
2	A	1330	NAD	C4A-C5A-N7A	-3.27	106.85	110.58
2	B	1330	NAD	C5A-N7A-C8A	3.24	108.54	103.45
2	B	1330	NAD	C2A-N3A-C4A	3.13	119.47	111.83
2	B	1330	NAD	C4A-C5A-N7A	-2.97	107.19	110.58

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	1330	NAD	C2B-C1B-N9A	2.76	120.17	113.30
2	B	1330	NAD	C5A-C4A-N9A	2.61	108.66	105.81
2	B	1330	NAD	O3-PA-O1A	-2.54	103.08	110.70
2	B	1330	NAD	C4A-N9A-C1B	-2.53	120.71	126.63
2	A	1330	NAD	N3A-C4A-N9A	2.42	131.29	127.17
2	A	1330	NAD	C3N-C7N-N7N	2.41	120.71	117.74
2	A	1330	NAD	O4B-C1B-N9A	2.35	112.59	108.09
2	B	1330	NAD	C1B-N9A-C8A	2.23	132.05	127.09
2	A	1330	NAD	C5A-C4A-N9A	2.14	108.14	105.81
2	A	1330	NAD	O3-PN-O1N	-2.13	104.30	110.70
2	A	1330	NAD	C4B-O4B-C1B	-2.01	105.03	109.47

There are no chirality outliers.

All (20) torsion outliers are listed below:

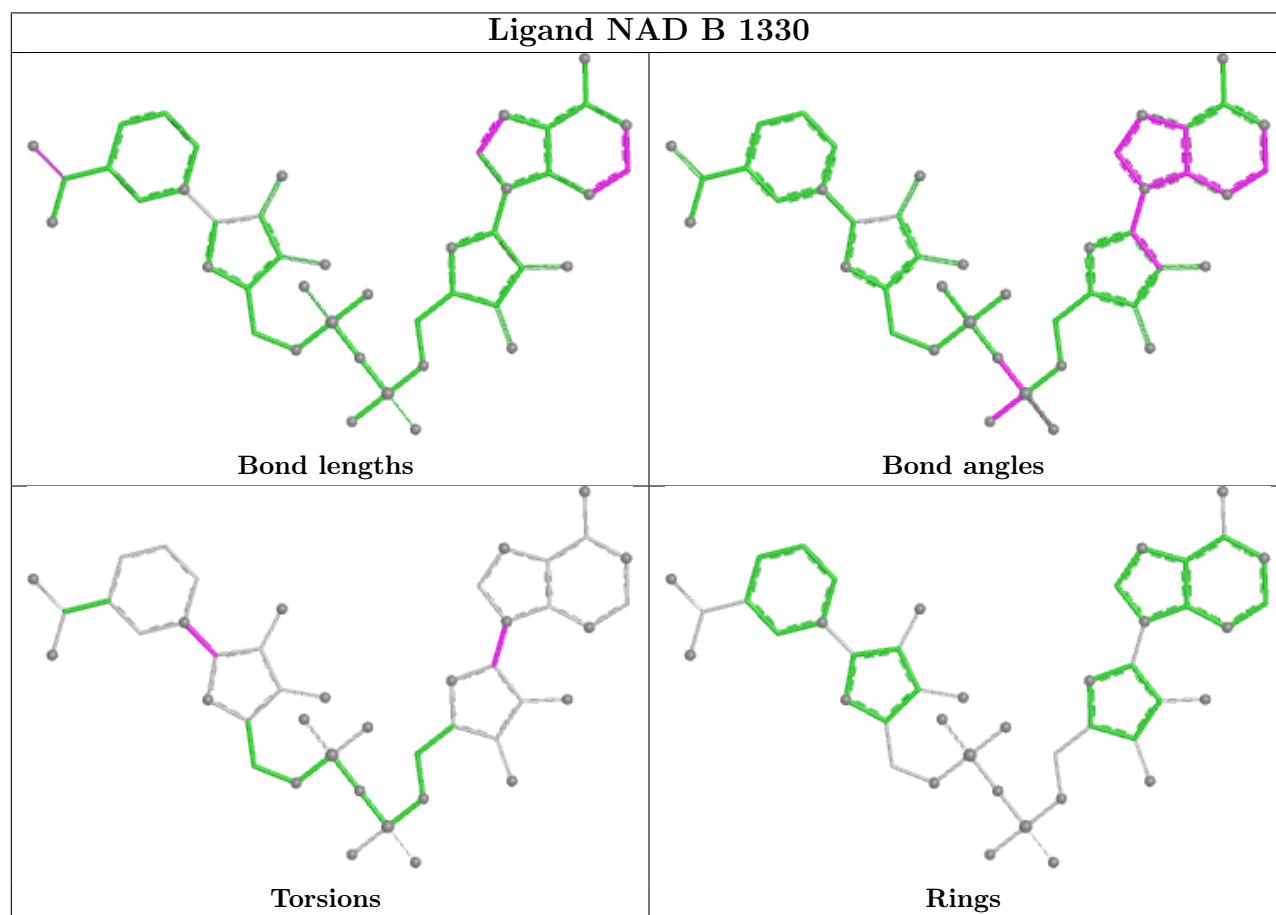
Mol	Chain	Res	Type	Atoms
2	A	1330	NAD	O4D-C1D-N1N-C2N
2	A	1330	NAD	O4D-C1D-N1N-C6N
2	A	1330	NAD	C2D-C1D-N1N-C2N
2	A	1330	NAD	C2D-C1D-N1N-C6N
2	B	1330	NAD	O4D-C1D-N1N-C2N
2	B	1330	NAD	O4D-C1D-N1N-C6N
2	B	1330	NAD	C2D-C1D-N1N-C2N
2	B	1330	NAD	C2D-C1D-N1N-C6N
3	A	1331	OXM	O1-C1-C2-O3
3	B	1331	OXM	N1-C1-C2-O2
3	B	1331	OXM	N1-C1-C2-O3
3	B	1331	OXM	O1-C1-C2-O2
3	B	1331	OXM	O1-C1-C2-O3
4	B	1332	GOL	C1-C2-C3-O3
4	B	1332	GOL	O2-C2-C3-O3
2	A	1330	NAD	PN-O3-PA-O5B
2	B	1330	NAD	C2B-C1B-N9A-C8A
2	A	1330	NAD	C2B-C1B-N9A-C8A
2	B	1330	NAD	O4B-C1B-N9A-C8A
2	A	1330	NAD	O4B-C1B-N9A-C8A

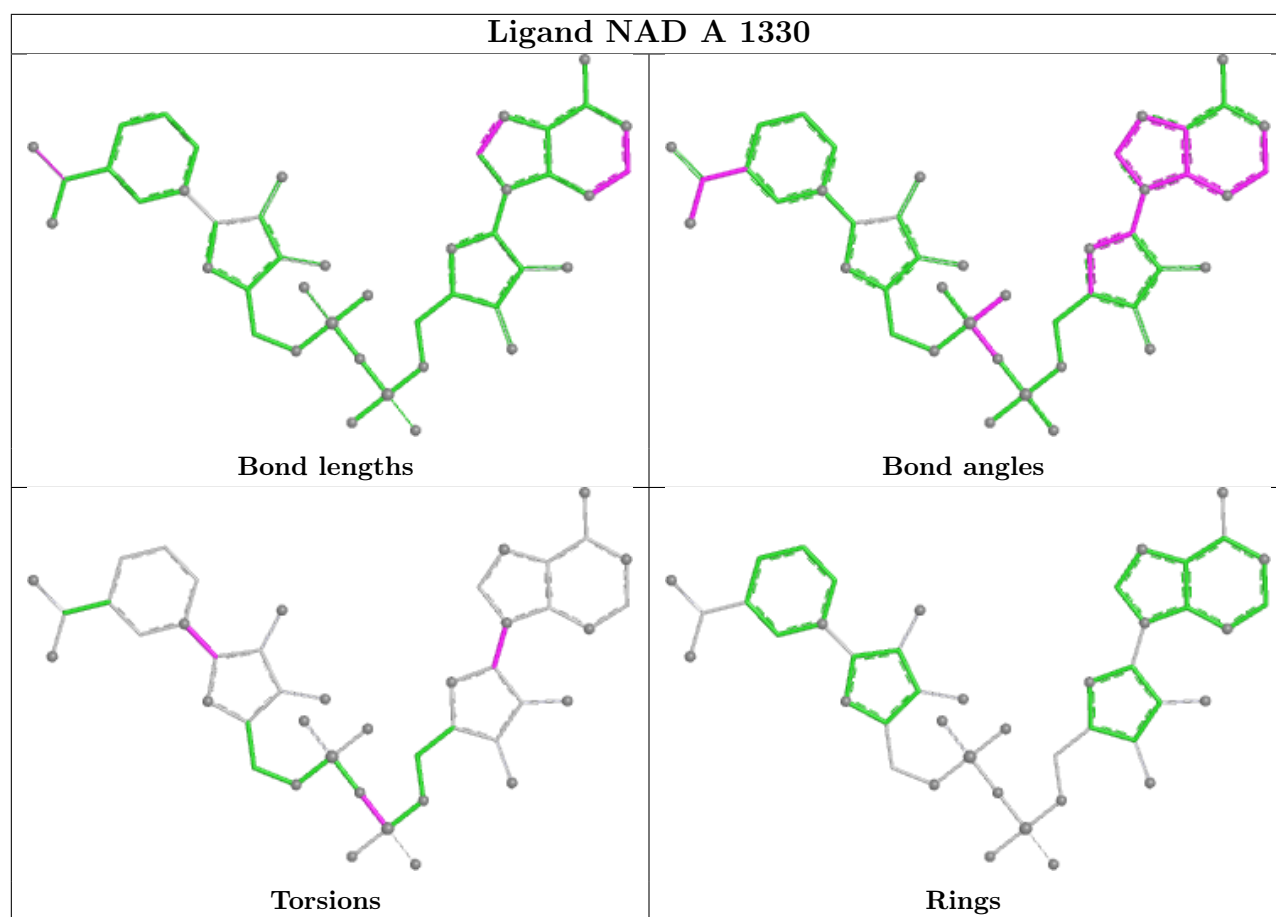
There are no ring outliers.

4 monomers are involved in 14 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	1330	NAD	3	0
3	A	1331	OXM	1	0
4	B	1332	GOL	4	0
2	A	1330	NAD	7	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.





5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
1	A	2
1	B	1

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	A	113:LEU	C	114:PRO	N	1.93
1	A	105(A):ASP	C	105(B):LYS	N	1.17
1	B	102:LYS	C	103(A):ALA	N	1.08

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	314/322 (97%)	0.55	27 (8%)	16 17	9, 32, 54, 94	13 (4%)
1	B	315/322 (97%)	0.43	29 (9%)	14 16	20, 30, 53, 94	12 (3%)
All	All	629/644 (97%)	0.49	56 (8%)	15 17	9, 31, 54, 94	25 (3%)

All (56) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	105(B)	LYS	18.3
1	A	115	LEU	9.9
1	A	113	LEU	7.3
1	B	112	LEU	6.1
1	B	115	LEU	5.8
1	B	114	PRO	5.8
1	A	114	PRO	5.7
1	B	113	LEU	5.5
1	B	329	ILE	5.3
1	B	103(E)	SER	5.1
1	A	107	TRP	5.0
1	B	107	TRP	5.0
1	B	105(A)	ASP	4.9
1	A	329	ILE	4.6
1	A	103(A)	ALA	4.3
1	A	102	LYS	4.2
1	B	116	ASN	4.2
1	B	18	ALA	4.0
1	B	103(B)	PRO	3.6
1	A	112	LEU	3.6
1	B	117	ASN	3.5
1	A	103(E)	SER	3.5
1	B	105(B)	LYS	3.3
1	B	129	ASN	3.2

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Mol	Chain	Res	Type	RSRZ
1	A	103(B)	PRO	3.1
1	B	328	LEU	3.1
1	A	18	ALA	3.0
1	B	108	ASN	2.9
1	A	100	PHE	2.7
1	B	102	LYS	2.7
1	B	103(C)	GLY	2.6
1	B	103(A)	ALA	2.6
1	B	119	ILE	2.6
1	A	109	ARG	2.6
1	A	103(C)	GLY	2.5
1	B	109	ARG	2.5
1	A	310	ASP	2.5
1	B	100	PHE	2.4
1	B	110	ASP	2.4
1	A	327	ALA	2.4
1	A	105(A)	ASP	2.4
1	B	111	ASP	2.3
1	A	19	PRO	2.3
1	B	101	THR	2.2
1	A	120	MET	2.1
1	B	106	GLU	2.1
1	A	116	ASN	2.1
1	B	241	ASN	2.1
1	A	320	ALA	2.1
1	B	245	SER	2.1
1	A	222	GLN	2.1
1	A	54	ILE	2.1
1	A	230	ASP	2.1
1	B	283	ASP	2.1
1	A	103(D)	LYS	2.0
1	A	245	SER	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 ⓘ

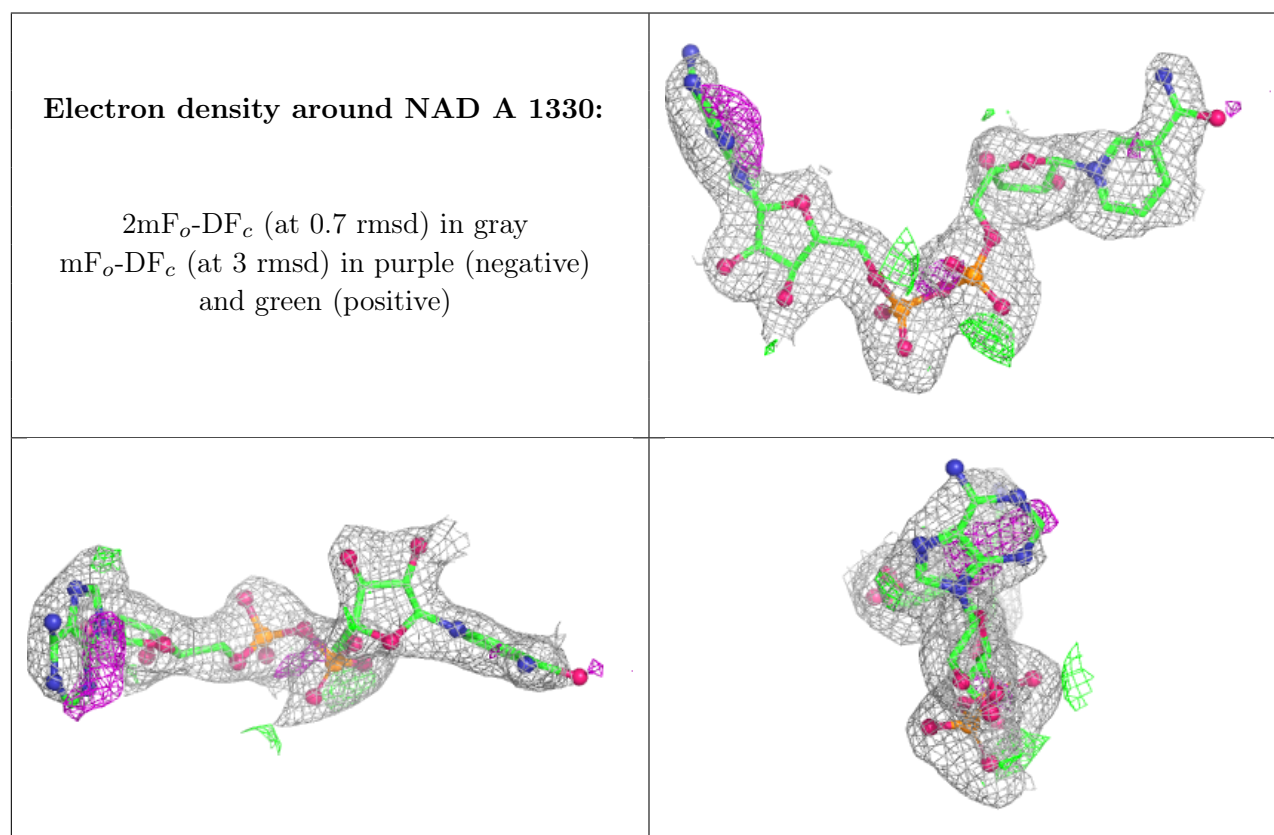
There are no oligosaccharides in this entry.

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.

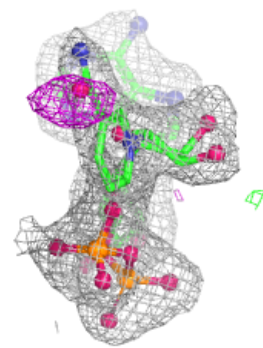
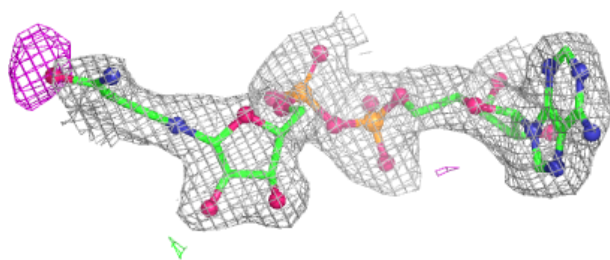
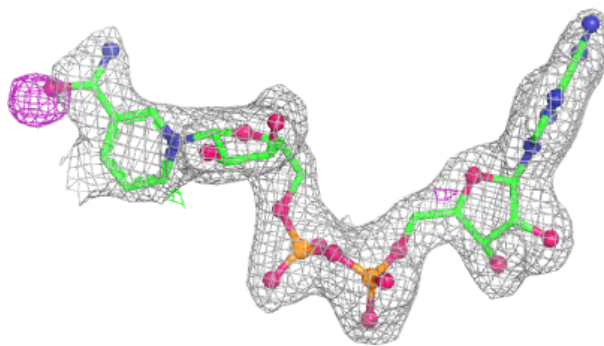
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	OXM	A	1331	6/6	0.84	0.14	49,50,50,50	0
2	NAD	A	1330	44/44	0.85	0.13	34,42,50,52	0
4	GOL	A	1332	6/6	0.91	0.09	24,24,26,28	0
3	OXM	B	1331	6/6	0.93	0.09	43,43,43,43	0
2	NAD	B	1330	44/44	0.96	0.07	18,33,37,40	0
4	GOL	B	1332	6/6	0.97	0.05	25,26,27,27	0

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.



Electron density around NAD B 1330:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



6.5 Other polymers [i](#)

There are no such residues in this entry.