



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 01:13 PM UTC

PDB ID : 5YAQ / pdb_00005yaq
Title : Crystal structure of scyllo-inositol dehydrogenase with L-glucose dehydrogenase activity complexed with scyllo-inosose
Authors : Fukano, K.; Shimizu, T.; Sasaki, Y.; Nakamura, A.; Yajima, S.
Deposited on : 2017-09-01
Resolution : 1.99 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

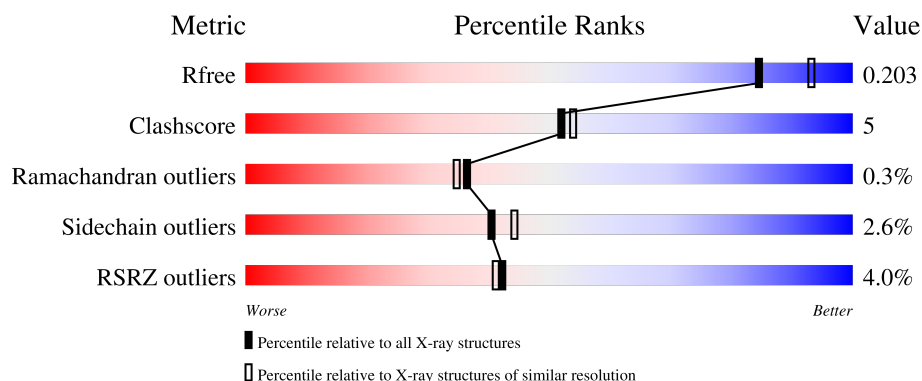
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.99 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	10052 (2.00-2.00)
Clashscore	190562	11152 (2.00-2.00)
Ramachandran outliers	187476	11031 (2.00-2.00)
Sidechain outliers	187428	11029 (2.00-2.00)
RSRZ outliers	180081	10067 (2.00-2.00)

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	380	3% 88% 8% • •
1	B	380	4% 85% 10% • •
1	C	380	6% 87% 8% • •
1	D	380	3% 89% 6% • 5%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 11787 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 Scyllo-inositol dehydrogenase with L-glucose dehydrogenase activity.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	366	Total	C	N	O	S	0	0	0
			2785	1744	510	513	18			
1	B	366	Total	C	N	O	S	0	0	0
			2785	1744	510	513	18			
1	C	366	Total	C	N	O	S	0	0	0
			2785	1744	510	513	18			
1	D	362	Total	C	N	O	S	0	0	0
			2755	1724	506	507	18			

There are 36 discrepancies between the modelled and reference sequences:

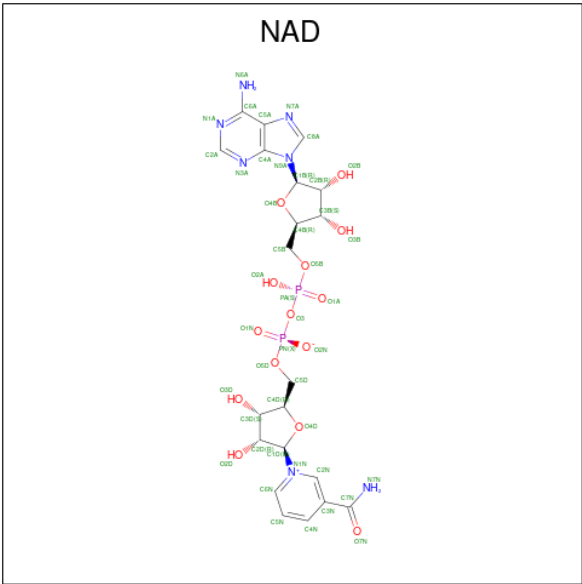
Chain	Residue	Modelled	Actual	Comment	Reference
A	72	SER	ASN	engineered mutation	UNP K7ZP76
A	373	LEU	-	expression tag	UNP K7ZP76
A	374	GLU	-	expression tag	UNP K7ZP76
A	375	HIS	-	expression tag	UNP K7ZP76
A	376	HIS	-	expression tag	UNP K7ZP76
A	377	HIS	-	expression tag	UNP K7ZP76
A	378	HIS	-	expression tag	UNP K7ZP76
A	379	HIS	-	expression tag	UNP K7ZP76
A	380	HIS	-	expression tag	UNP K7ZP76
B	72	SER	ASN	engineered mutation	UNP K7ZP76
B	373	LEU	-	expression tag	UNP K7ZP76
B	374	GLU	-	expression tag	UNP K7ZP76
B	375	HIS	-	expression tag	UNP K7ZP76
B	376	HIS	-	expression tag	UNP K7ZP76
B	377	HIS	-	expression tag	UNP K7ZP76
B	378	HIS	-	expression tag	UNP K7ZP76
B	379	HIS	-	expression tag	UNP K7ZP76
B	380	HIS	-	expression tag	UNP K7ZP76
C	72	SER	ASN	engineered mutation	UNP K7ZP76
C	373	LEU	-	expression tag	UNP K7ZP76

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Chain	Residue	Modelled	Actual	Comment	Reference
C	374	GLU	-	expression tag	UNP K7ZP76
C	375	HIS	-	expression tag	UNP K7ZP76
C	376	HIS	-	expression tag	UNP K7ZP76
C	377	HIS	-	expression tag	UNP K7ZP76
C	378	HIS	-	expression tag	UNP K7ZP76
C	379	HIS	-	expression tag	UNP K7ZP76
C	380	HIS	-	expression tag	UNP K7ZP76
D	72	SER	ASN	engineered mutation	UNP K7ZP76
D	373	LEU	-	expression tag	UNP K7ZP76
D	374	GLU	-	expression tag	UNP K7ZP76
D	375	HIS	-	expression tag	UNP K7ZP76
D	376	HIS	-	expression tag	UNP K7ZP76
D	377	HIS	-	expression tag	UNP K7ZP76
D	378	HIS	-	expression tag	UNP K7ZP76
D	379	HIS	-	expression tag	UNP K7ZP76
D	380	HIS	-	expression tag	UNP K7ZP76

- Molecule 2 is NICOTINAMIDE-ADENINE-DINUCLEOTIDE (CCD ID: NAD) (formula: C₂₁H₂₇N₇O₁₄P₂).



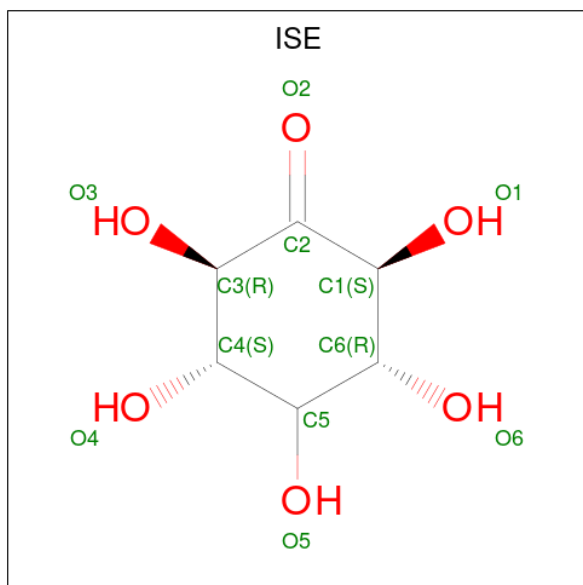
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			44	21	7	14	2		
2	B	1	Total	C	N	O	P	0	0
			44	21	7	14	2		
2	C	1	Total	C	N	O	P	0	0
			44	21	7	14	2		

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

- Molecule 3 is (2R,3S,4s,5R,6S)-2,3,4,5,6-pentahydroxycyclohexanone (CCD ID: ISE) (formula: C₆H₁₀O₆).



Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	1	Total	C O	0	0
			12	6 6		
3	C	1	Total	C O	0	0
			12	6 6		
3	D	1	Total	C O	0	0
			12	6 6		

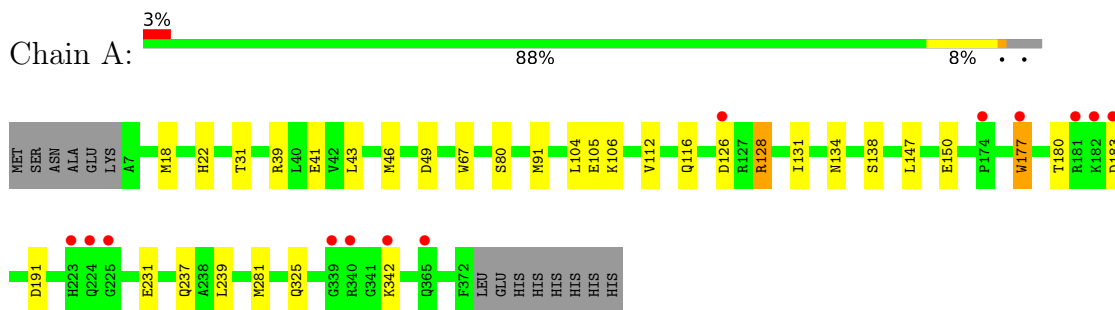
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	122	Total	O	0	0
			122	122		
4	B	124	Total	O	0	0
			124	124		
4	C	94	Total	O	0	0
			94	94		
4	D	125	Total	O	0	0
			125	125		

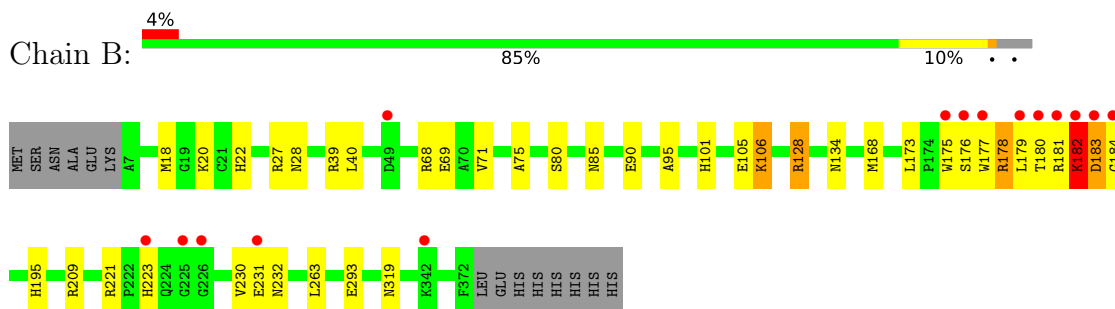
3 Residue-property plots

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.

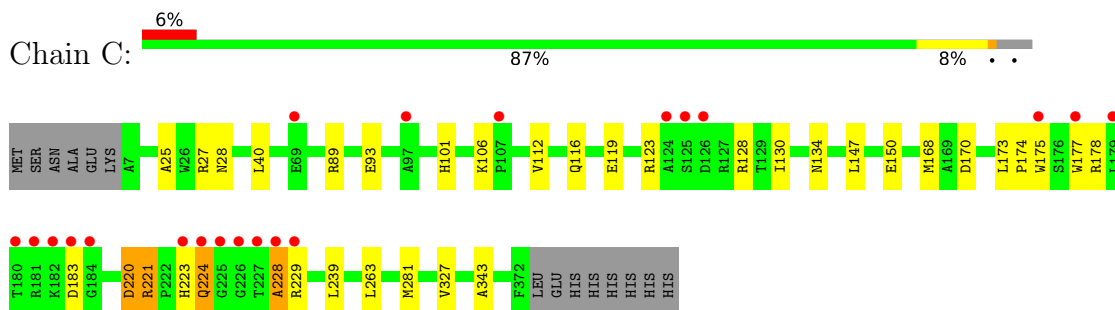
- Molecule 1: Scyllo-inositol dehydrogenase with L-glucose dehydrogenase activity



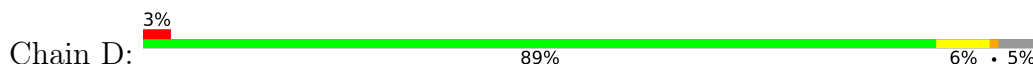
- Molecule 1: Scyllo-inositol dehydrogenase with L-glucose dehydrogenase activity

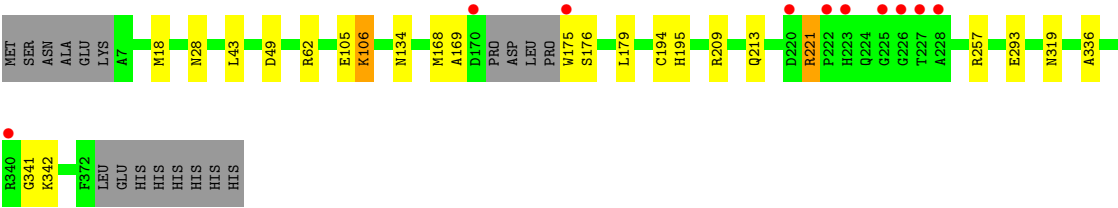


- Molecule 1: Scyllo-inositol dehydrogenase with L-glucose dehydrogenase activity



- Molecule 1: Scyllo-inositol dehydrogenase with L-glucose dehydrogenase activity





4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	86.55Å 121.24Å 140.40Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	39.00 – 1.99 39.00 – 1.99	Depositor EDS
% Data completeness (in resolution range)	99.5 (39.00-1.99) 99.5 (39.00-1.99)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.69 (at 2.00Å)	Xtriage
Refinement program	REFMAC 5.8.0107	Depositor
R, R_{free}	0.180 , 0.194 0.189 , 0.203	Depositor DCC
R_{free} test set	5054 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	21.5	Xtriage
Anisotropy	0.394	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 30.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	11787	wwPDB-VP
Average B, all atoms (Å ²)	23.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.87% 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: NAD, ISE

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	0.69	0/2851	0.90	0/3865
1	B	0.70	0/2851	0.90	2/3865 (0.1%)
1	C	0.69	1/2851 (0.0%)	0.91	4/3865 (0.1%)
1	D	0.67	0/2818	0.89	0/3816
All	All	0.69	1/11371 (0.0%)	0.90	6/15411 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	220	ASP	C-N	5.79	1.40	1.33

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	228	ALA	CA-C-N	-6.93	111.94	122.81
1	C	228	ALA	C-N-CA	-6.93	111.94	122.81
1	B	128	ARG	NE-CZ-NH2	-6.06	113.74	119.20
1	B	128	ARG	CG-CD-NE	-5.72	99.41	112.00
1	C	221	ARG	CA-C-N	5.02	126.11	119.84
1	C	221	ARG	C-N-CA	5.02	126.11	119.84

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	2785	0	2702	18	0
1	B	2785	0	2702	58	0
1	C	2785	0	2702	21	0
1	D	2755	0	2672	23	0
2	A	44	0	26	1	0
2	B	44	0	26	4	0
2	C	44	0	26	1	0
2	D	44	0	26	4	0
3	B	12	0	10	1	0
3	C	12	0	10	1	0
3	D	12	0	10	1	0
4	A	122	0	0	1	0
4	B	124	0	0	5	0
4	C	94	0	0	0	0
4	D	125	0	0	2	0
All	All	11787	0	10912	114	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 5.

All (114) 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:B:177:TRP:CE2	1:B:178:ARG:HG3	1.55	1.39
1:B:177:TRP:CZ2	1:B:178:ARG:HG3	1.96	1.01
1:B:177:TRP:CG	1:B:178:ARG:H	1.80	1.00
1:B:177:TRP:HE1	1:B:178:ARG:NH1	1.63	0.96
1:B:177:TRP:CE2	1:B:178:ARG:CG	2.48	0.95
1:B:177:TRP:CZ2	1:B:178:ARG:CG	2.50	0.95
1:B:177:TRP:HE1	1:B:178:ARG:HH11	1.02	0.94
1:C:89:ARG:NH1	1:C:93:GLU:OE2	2.04	0.91
1:B:177:TRP:O	1:B:179:LEU:N	2.08	0.86
1:D:49:ASP:HB2	4:D:620:HOH:O	1.76	0.85
1:A:191:ASP:HB2	4:A:638:HOH:O	1.82	0.80
1:B:177:TRP:NE1	1:B:178:ARG:HG3	1.96	0.80
1:B:175:TRP:CZ3	1:B:223:HIS:CE1	2.71	0.78
1:B:177:TRP:HZ2	1:B:178:ARG:CZ	1.98	0.76
1:B:177:TRP:CG	1:B:178:ARG:N	2.42	0.75
1:B:177:TRP:NE1	1:B:178:ARG:HH11	1.84	0.72
1:B:177:TRP:NE1	1:B:178:ARG:NH1	2.37	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:106:LYS:HD3	1:B:195:HIS:NE2	2.07	0.70
1:B:177:TRP:CD1	1:B:178:ARG:H	2.09	0.70
1:B:168:MET:HE2	1:B:221:ARG:CZ	2.21	0.70
1:A:39:ARG:NH1	1:A:41:GLU:OE2	2.25	0.69
1:B:209:ARG:HG3	4:B:618:HOH:O	1.94	0.68
1:B:101:HIS:HD2	1:B:128:ARG:H	1.39	0.68
2:D:401:NAD:C3N	3:D:402:ISE:O2	2.43	0.67
1:B:177:TRP:CZ2	1:B:178:ARG:CZ	2.80	0.65
1:C:101:HIS:HD2	1:C:128:ARG:H	1.44	0.64
1:C:112:VAL:O	1:C:116:GLN:HG3	1.99	0.63
1:B:177:TRP:CZ2	1:B:178:ARG:HG2	2.34	0.62
1:D:106:LYS:HD3	1:D:194:CYS:SG	2.40	0.61
1:C:173:LEU:O	1:C:221:ARG:HD3	2.02	0.60
1:B:177:TRP:CD2	1:B:178:ARG:N	2.68	0.59
1:D:342:LYS:HE2	4:D:501:HOH:O	2.02	0.58
1:B:175:TRP:CE3	1:B:223:HIS:CE1	2.91	0.58
1:B:231:GLU:C	4:B:501:HOH:O	2.47	0.58
1:B:182:LYS:O	1:B:184:GLY:N	2.37	0.57
1:A:150:GLU:OE1	1:D:293:GLU:OE1	2.22	0.57
1:B:71:VAL:CG1	1:B:95:ALA:HA	2.34	0.57
1:C:168:MET:O	1:C:221:ARG:NH2	2.38	0.56
1:D:168:MET:HE3	1:D:179:LEU:HD21	1.86	0.56
1:B:177:TRP:HZ2	1:B:178:ARG:NE	2.03	0.56
1:B:71:VAL:HG13	1:B:95:ALA:HA	1.89	0.55
1:D:175:TRP:CE3	1:D:221:ARG:HG3	2.41	0.54
1:B:22:HIS:HE1	1:B:105:GLU:OE2	1.90	0.54
1:B:175:TRP:CE3	1:B:175:TRP:O	2.61	0.54
1:A:177:TRP:HD1	1:A:183:ASP:HB2	1.73	0.53
1:B:177:TRP:CZ2	1:B:178:ARG:NE	2.77	0.53
1:D:106:LYS:HD3	1:D:194:CYS:HG	1.73	0.53
1:A:281:MET:HG2	1:B:319:ASN:HD22	1.72	0.53
1:B:179:LEU:HD11	1:B:221:ARG:HD2	1.92	0.52
1:A:22:HIS:HE1	1:A:105:GLU:OE2	1.93	0.51
2:B:401:NAD:C3N	3:B:402:ISE:O2	2.58	0.51
1:A:31:THR:O	1:B:20:LYS:HE3	2.11	0.51
1:C:220:ASP:HA	1:C:228:ALA:O	2.11	0.51
1:D:179:LEU:CD1	1:D:221:ARG:HD2	2.41	0.51
1:A:138:SER:HB3	1:A:325:GLN:HE21	1.75	0.50
1:B:177:TRP:C	1:B:179:LEU:N	2.69	0.50
1:D:18:MET:HE3	2:D:401:NAD:C4N	2.41	0.50
1:C:175:TRP:HH2	1:C:177:TRP:CD1	2.30	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:177:TRP:C	1:B:179:LEU:H	2.21	0.48
1:B:105:GLU:OE1	2:B:401:NAD:H2N	2.13	0.48
1:B:177:TRP:O	1:B:178:ARG:C	2.57	0.48
1:B:22:HIS:HD2	1:B:80:SER:OG	1.96	0.48
1:B:85:ASN:HB2	1:B:183:ASP:O	2.14	0.48
1:B:223:HIS:CE1	4:B:510:HOH:O	2.68	0.47
1:C:223:HIS:CE1	1:C:224:GLN:OE1	2.68	0.47
1:C:281:MET:HG2	1:D:319:ASN:HD22	1.79	0.47
1:A:126:ASP:O	1:A:128:ARG:NH2	2.47	0.47
1:D:106:LYS:HD2	1:D:106:LYS:C	2.39	0.47
1:D:106:LYS:HE2	1:D:195:HIS:NE2	2.30	0.47
1:A:22:HIS:HD2	1:A:80:SER:OG	1.97	0.47
1:A:147:LEU:HD21	1:D:293:GLU:HG2	1.97	0.47
1:B:181:ARG:C	1:B:182:LYS:O	2.58	0.47
1:B:209:ARG:HH22	1:D:213:GLN:HE21	1.63	0.47
1:B:168:MET:HA	1:B:168:MET:HE3	1.97	0.46
1:B:18:MET:HE3	2:B:401:NAD:C4N	2.45	0.46
1:C:223:HIS:O	1:C:224:GLN:HG2	2.16	0.46
1:D:105:GLU:OE1	2:D:401:NAD:H2N	2.16	0.46
1:B:177:TRP:CD2	1:B:178:ARG:HG3	2.36	0.45
1:B:293:GLU:HG2	1:C:147:LEU:HD21	1.98	0.45
1:A:239:LEU:HG	1:C:239:LEU:HG	1.98	0.45
1:C:119:GLU:OE2	1:C:123:ARG:NH1	2.50	0.45
1:B:175:TRP:CE3	1:B:223:HIS:ND1	2.84	0.45
1:D:18:MET:HG3	2:D:401:NAD:C2N	2.48	0.44
1:D:179:LEU:HD11	1:D:221:ARG:HD2	2.00	0.44
1:B:68:ARG:NH2	1:B:90:GLU:OE2	2.51	0.44
1:D:43:LEU:C	1:D:43:LEU:HD23	2.43	0.43
1:B:293:GLU:OE2	1:C:150:GLU:OE1	2.36	0.43
1:C:170:ASP:O	1:C:221:ARG:NH1	2.35	0.43
1:A:180:THR:HA	1:A:231:GLU:HB2	2.01	0.42
1:D:209:ARG:HD2	1:D:209:ARG:C	2.44	0.42
1:B:39:ARG:NH2	1:B:75:ALA:O	2.52	0.42
1:B:180:THR:O	1:B:232:ASN:ND2	2.52	0.42
1:C:223:HIS:C	1:C:224:GLN:HG2	2.44	0.42
1:B:179:LEU:HD13	1:B:230:VAL:HG22	2.02	0.42
1:D:336:ALA:HA	1:D:341:GLY:O	2.19	0.42
1:A:43:LEU:C	1:A:43:LEU:HD23	2.45	0.42
1:A:18:MET:HG3	2:A:500:NAD:C2N	2.50	0.42
1:C:175:TRP:NE1	1:C:229:ARG:O	2.50	0.42
1:D:168:MET:CE	1:D:179:LEU:HD21	2.50	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:169:ALA:HB1	1:D:257:ARG:HD3	2.02	0.42
1:B:18:MET:HG3	2:B:401:NAD:C2N	2.49	0.41
1:B:27:ARG:HD2	1:B:40:LEU:HD11	2.01	0.41
1:C:173:LEU:HA	1:C:174:PRO:HD3	1.94	0.41
1:C:25:ALA:HA	1:C:327:VAL:HG22	2.03	0.41
1:C:130:ILE:HD12	1:C:343:ALA:HB1	2.03	0.41
1:B:293:GLU:HB2	4:B:606:HOH:O	2.19	0.41
1:A:67:TRP:CD1	1:A:91:MET:HE2	2.56	0.41
1:A:112:VAL:O	1:A:116:GLN:HG3	2.21	0.41
2:C:401:NAD:C3N	3:C:402:ISE:O2	2.69	0.40
1:B:209:ARG:C	1:B:209:ARG:HD2	2.45	0.40
1:A:104:LEU:O	1:A:131:ILE:HA	2.21	0.40
1:B:232:ASN:N	4:B:501:HOH:O	2.55	0.40
1:C:27:ARG:HG2	1:C:40:LEU:HD11	2.04	0.40
1:D:169:ALA:CB	1:D:257:ARG:HD3	2.52	0.40

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	364/380 (96%)	357 (98%)	7 (2%)	0	100	100
1	B	364/380 (96%)	349 (96%)	12 (3%)	3 (1%)	16	11
1	C	364/380 (96%)	354 (97%)	9 (2%)	1 (0%)	36	35
1	D	358/380 (94%)	351 (98%)	7 (2%)	0	100	100
All	All	1450/1520 (95%)	1411 (97%)	35 (2%)	4 (0%)	36	35

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	178	ARG
1	B	183	ASP
1	C	224	GLN
1	B	182	LYS

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	275/288 (96%)	267 (97%)	8 (3%)	37	40
1	B	275/288 (96%)	267 (97%)	8 (3%)	37	40
1	C	275/288 (96%)	269 (98%)	6 (2%)	45	50
1	D	271/288 (94%)	265 (98%)	6 (2%)	45	50
All	All	1096/1152 (95%)	1068 (97%)	28 (3%)	40	44

All (28) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	46	MET
1	A	49	ASP
1	A	106	LYS
1	A	128	ARG
1	A	134	ASN
1	A	177	TRP
1	A	237	GLN
1	A	342	LYS
1	B	28	ASN
1	B	69	GLU
1	B	106	LYS
1	B	134	ASN
1	B	173	LEU
1	B	176	SER
1	B	182	LYS
1	B	263	LEU
1	C	28	ASN
1	C	106	LYS

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Mol	Chain	Res	Type
1	C	134	ASN
1	C	178	ARG
1	C	183	ASP
1	C	263	LEU
1	D	28	ASN
1	D	62	ARG
1	D	106	LYS
1	D	134	ASN
1	D	176	SER
1	D	221	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (24) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	22	HIS
1	A	28	ASN
1	A	235	GLN
1	A	237	GLN
1	A	278	GLN
1	A	325	GLN
1	A	358	HIS
1	B	22	HIS
1	B	28	ASN
1	B	101	HIS
1	B	223	HIS
1	B	319	ASN
1	B	358	HIS
1	B	368	GLN
1	B	371	ASN
1	C	28	ASN
1	C	101	HIS
1	C	223	HIS
1	C	335	GLN
1	C	358	HIS
1	D	28	ASN
1	D	213	GLN
1	D	319	ASN
1	D	371	ASN

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 ⓘ

7 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	500	-	46,48,48	1.47	7 (15%)	64,73,73	1.80	12 (18%)
2	NAD	B	401	-	46,48,48	1.45	7 (15%)	64,73,73	1.76	14 (21%)
2	NAD	C	401	-	46,48,48	1.42	6 (13%)	64,73,73	1.71	13 (20%)
2	NAD	D	401	-	46,48,48	1.35	7 (15%)	64,73,73	1.84	11 (17%)
3	ISE	D	402	-	12,12,12	1.03	0	15,18,18	1.49	3 (20%)
3	ISE	B	402	-	12,12,12	0.96	0	15,18,18	1.71	2 (13%)
3	ISE	C	402	-	12,12,12	0.87	0	15,18,18	1.58	2 (13%)

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	500	-	-	3/30/62/62	0/5/5/5

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	NAD	B	401	-	-	2/30/62/62	0/5/5/5
2	NAD	C	401	-	-	3/30/62/62	0/5/5/5
2	NAD	D	401	-	-	4/30/62/62	0/5/5/5
3	ISE	D	402	-	-	-	0/1/1/1
3	ISE	B	402	-	-	-	0/1/1/1
3	ISE	C	402	-	-	-	0/1/1/1

All (27) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	401	NAD	C5A-C4A	5.08	1.48	1.39
2	C	401	NAD	C5A-C4A	4.87	1.47	1.39
2	D	401	NAD	C5A-C4A	4.71	1.47	1.39
2	A	500	NAD	C5A-C4A	4.41	1.46	1.39
2	A	500	NAD	PN-O3	4.25	1.64	1.59
2	C	401	NAD	O4D-C1D	3.51	1.45	1.40
2	B	401	NAD	O4D-C1D	3.42	1.45	1.40
2	C	401	NAD	PN-O3	3.15	1.62	1.59
2	D	401	NAD	C5A-C6A	3.13	1.49	1.41
2	A	500	NAD	O4D-C1D	3.04	1.44	1.40
2	C	401	NAD	C5A-C6A	2.94	1.49	1.41
2	B	401	NAD	PN-O3	2.85	1.62	1.59
2	A	500	NAD	C5A-C6A	2.84	1.48	1.41
2	B	401	NAD	C5A-C6A	2.80	1.48	1.41
2	C	401	NAD	C8A-N7A	2.76	1.37	1.31
2	A	500	NAD	C8A-N7A	2.69	1.36	1.31
2	B	401	NAD	C8A-N7A	2.62	1.36	1.31
2	A	500	NAD	PA-O3	2.59	1.62	1.59
2	D	401	NAD	C8A-N7A	2.54	1.36	1.31
2	D	401	NAD	PN-O3	2.47	1.62	1.59
2	D	401	NAD	C7N-N7N	2.28	1.37	1.33
2	D	401	NAD	O4D-C1D	2.26	1.43	1.40
2	B	401	NAD	C7N-N7N	2.25	1.37	1.33
2	A	500	NAD	C7N-N7N	2.22	1.37	1.33
2	C	401	NAD	PA-O3	2.20	1.61	1.59
2	B	401	NAD	PA-O3	2.10	1.61	1.59
2	D	401	NAD	C4N-C3N	2.10	1.42	1.39

All (57) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	401	NAD	C5A-C4A-N3A	-5.77	118.77	126.72
2	C	401	NAD	C5A-C4A-N3A	-5.09	119.71	126.72
2	A	500	NAD	C4A-N9A-C8A	4.93	110.91	105.74
3	B	402	ISE	O1-C1-C6	-4.89	100.37	110.53
2	B	401	NAD	C5A-C4A-N3A	-4.86	120.02	126.72
2	D	401	NAD	N3A-C4A-N9A	4.84	135.39	127.17
2	C	401	NAD	N3A-C4A-N9A	4.80	135.33	127.17
2	A	500	NAD	C5A-C4A-N3A	-4.71	120.23	126.72
2	A	500	NAD	N3A-C4A-N9A	4.70	135.17	127.17
2	A	500	NAD	N3A-C2A-N1A	-4.65	121.54	128.58
2	C	401	NAD	N3A-C2A-N1A	-4.29	122.09	128.58
2	C	401	NAD	C4A-N9A-C8A	4.25	110.20	105.74
2	B	401	NAD	N3A-C4A-N9A	4.25	134.39	127.17
2	D	401	NAD	N3A-C2A-N1A	-4.24	122.17	128.58
2	D	401	NAD	C2A-N3A-C4A	4.17	122.01	111.83
3	C	402	ISE	O1-C1-C6	-4.07	102.07	110.53
2	D	401	NAD	C4A-N9A-C8A	4.05	109.99	105.74
2	A	500	NAD	C2A-N3A-C4A	3.95	121.48	111.83
2	B	401	NAD	N3A-C2A-N1A	-3.95	122.61	128.58
2	D	401	NAD	C4A-C5A-N7A	-3.89	106.13	110.58
2	C	401	NAD	C2A-N3A-C4A	3.88	121.31	111.83
2	B	401	NAD	C4A-N9A-C8A	3.84	109.77	105.74
2	B	401	NAD	C2A-N3A-C4A	3.45	120.25	111.83
2	B	401	NAD	C4A-C5A-N7A	-3.30	106.81	110.58
3	D	402	ISE	O1-C1-C6	-3.29	103.70	110.53
2	D	401	NAD	C5A-N7A-C8A	3.26	108.57	103.45
2	A	500	NAD	N9A-C8A-N7A	-3.22	109.37	113.94
2	D	401	NAD	N9A-C8A-N7A	-3.15	109.47	113.94
2	D	401	NAD	C6A-C5A-N7A	3.06	137.99	132.09
2	A	500	NAD	O2A-PA-O1A	2.85	125.72	112.44
2	B	401	NAD	C5A-N7A-C8A	2.85	107.93	103.45
2	B	401	NAD	N9A-C8A-N7A	-2.84	109.91	113.94
2	C	401	NAD	N9A-C8A-N7A	-2.77	110.00	113.94
2	A	500	NAD	C6A-C5A-N7A	2.59	137.09	132.09
2	C	401	NAD	C4A-C5A-N7A	-2.55	107.67	110.58
2	C	401	NAD	O2A-PA-O1A	2.54	124.27	112.44
2	B	401	NAD	C2A-N1A-C6A	2.48	122.81	118.73
2	A	500	NAD	C5A-N7A-C8A	2.45	107.29	103.45
2	B	401	NAD	C6A-C5A-N7A	2.43	136.77	132.09
2	C	401	NAD	C6A-C5A-N7A	2.42	136.76	132.09
2	A	500	NAD	C4A-C5A-N7A	-2.42	107.81	110.58
2	C	401	NAD	C5A-N7A-C8A	2.37	107.17	103.45
3	B	402	ISE	C6-C5-C4	-2.31	106.77	110.83

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	401	NAD	O2A-PA-O1A	2.31	123.21	112.44
2	C	401	NAD	C2A-N1A-C6A	2.29	122.50	118.73
2	B	401	NAD	C4A-N9A-C1B	-2.27	121.31	126.63
2	B	401	NAD	C3B-C2B-C1B	-2.26	97.19	101.46
2	A	500	NAD	C2A-N1A-C6A	2.25	122.43	118.73
2	B	401	NAD	O2N-PN-O1N	2.22	122.77	112.44
3	D	402	ISE	O3-C3-C2	2.14	116.04	110.11
2	A	500	NAD	C4A-N9A-C1B	-2.14	121.62	126.63
2	D	401	NAD	O3-PA-O1A	-2.12	104.33	110.70
3	C	402	ISE	C5-C4-C3	2.09	113.52	110.22
2	C	401	NAD	O2N-PN-O3	2.09	112.91	107.27
2	C	401	NAD	O4B-C1B-N9A	2.03	111.99	108.09
2	D	401	NAD	C3B-C2B-C1B	-2.02	97.65	101.46
3	D	402	ISE	O2-C2-C1	-2.01	118.53	121.73

There are no chirality outliers.

All (12) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	500	NAD	O4D-C1D-N1N-C2N
2	A	500	NAD	O4D-C1D-N1N-C6N
2	B	401	NAD	O4D-C1D-N1N-C2N
2	B	401	NAD	O4D-C1D-N1N-C6N
2	C	401	NAD	O4D-C1D-N1N-C2N
2	C	401	NAD	O4D-C1D-N1N-C6N
2	D	401	NAD	O4D-C1D-N1N-C2N
2	D	401	NAD	O4D-C1D-N1N-C6N
2	D	401	NAD	C2D-C1D-N1N-C2N
2	A	500	NAD	C2D-C1D-N1N-C2N
2	C	401	NAD	C2D-C1D-N1N-C2N
2	D	401	NAD	O4B-C4B-C5B-O5B

There are no ring outliers.

7 monomers are involved in 10 short contacts:

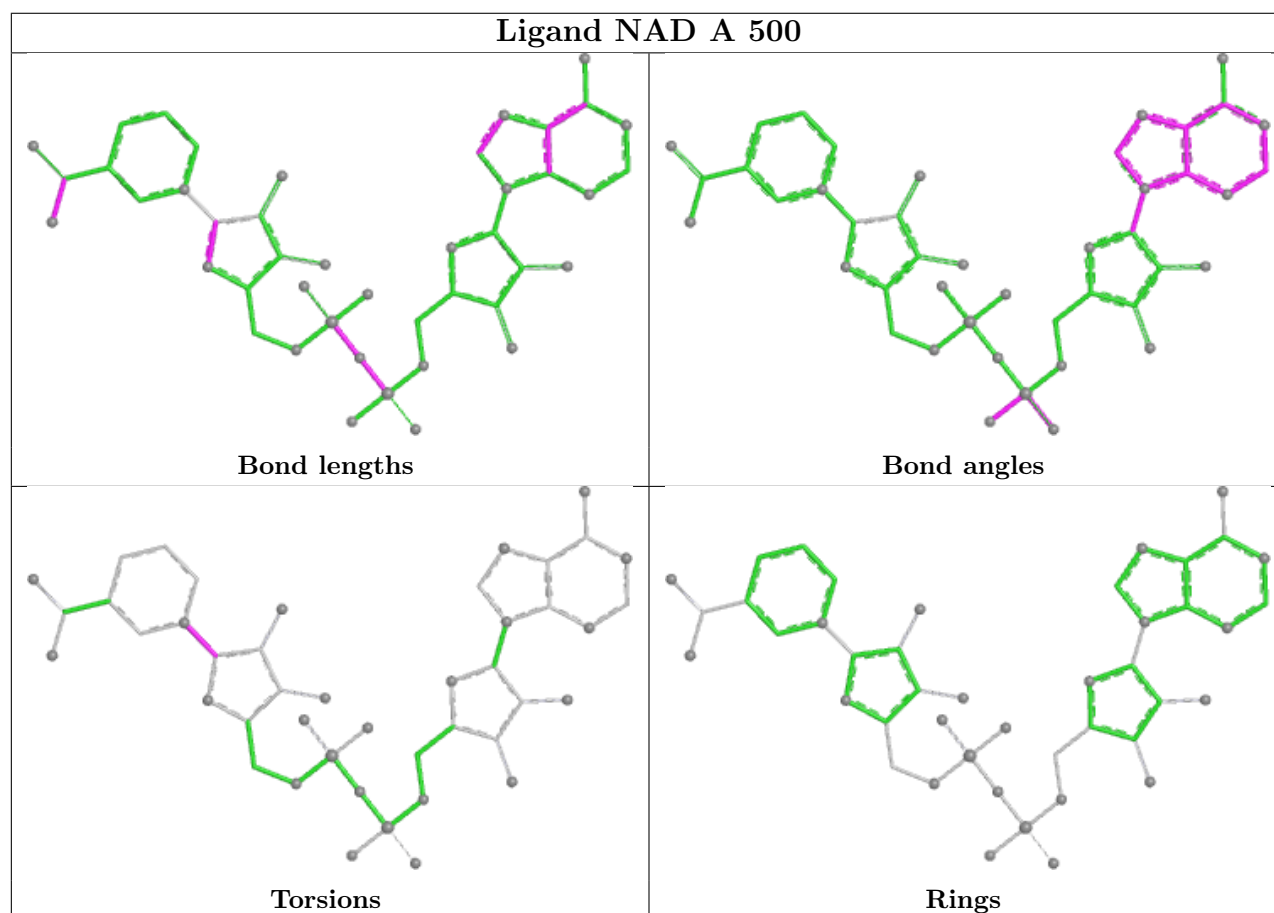
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	500	NAD	1	0
2	B	401	NAD	4	0
2	C	401	NAD	1	0
2	D	401	NAD	4	0
3	D	402	ISE	1	0

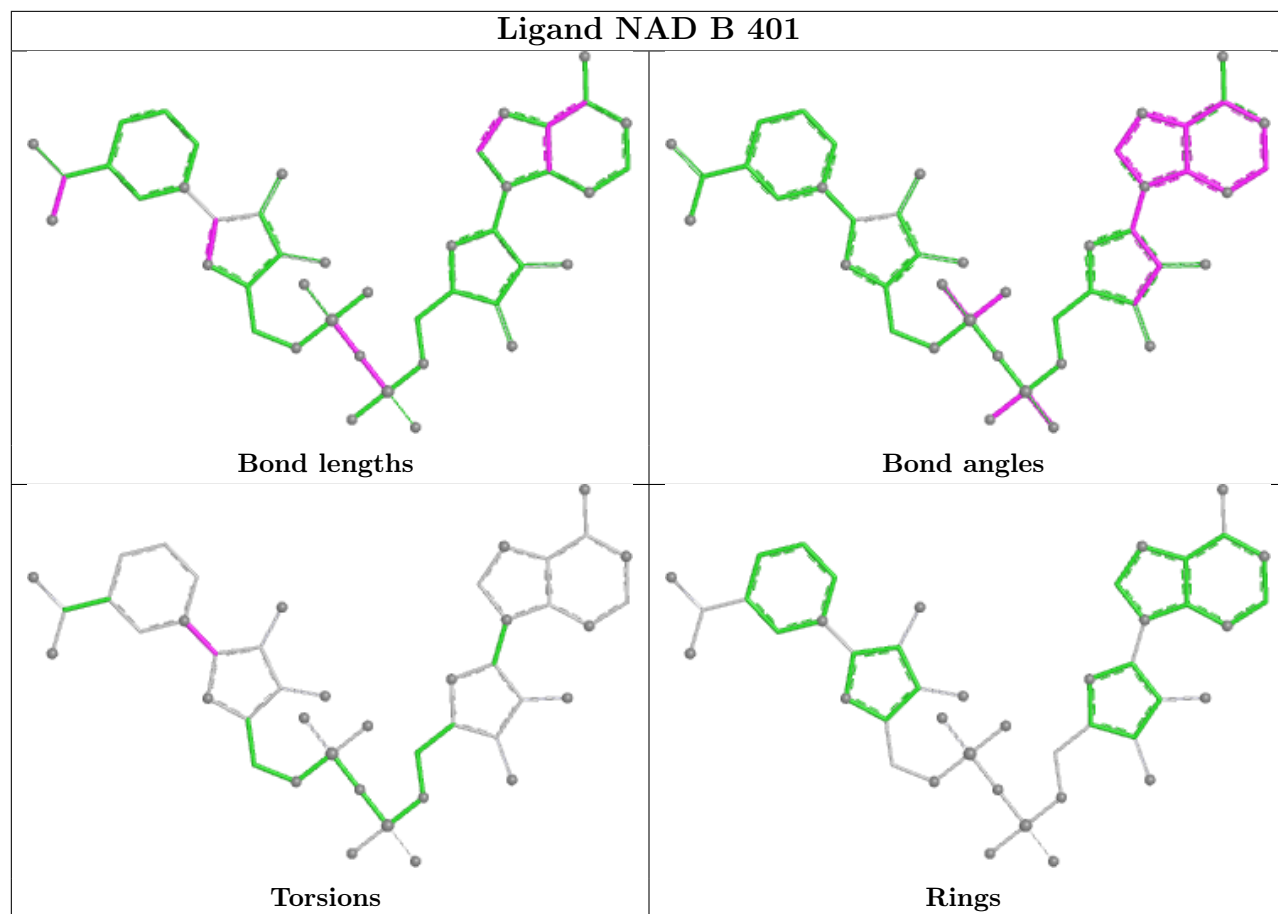
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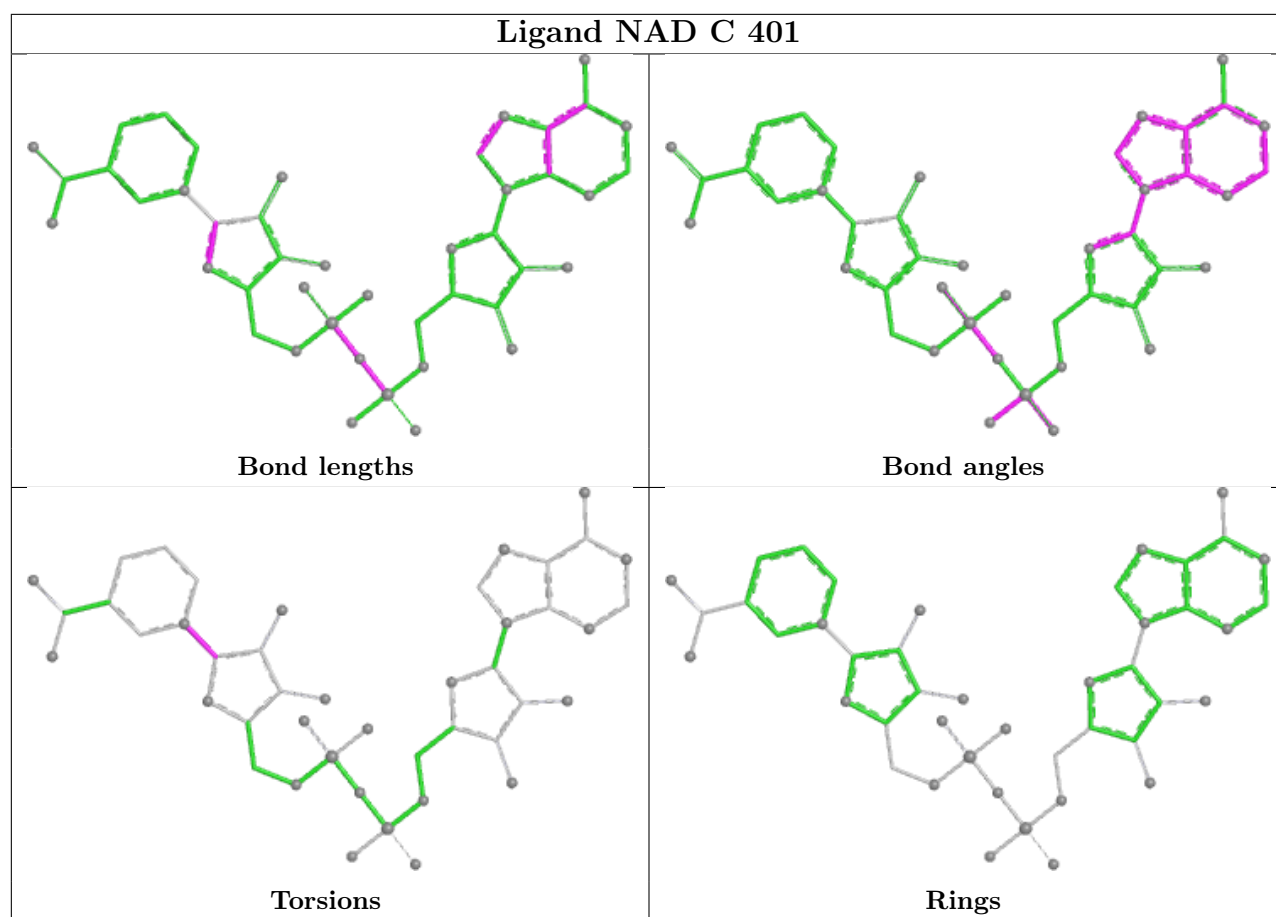
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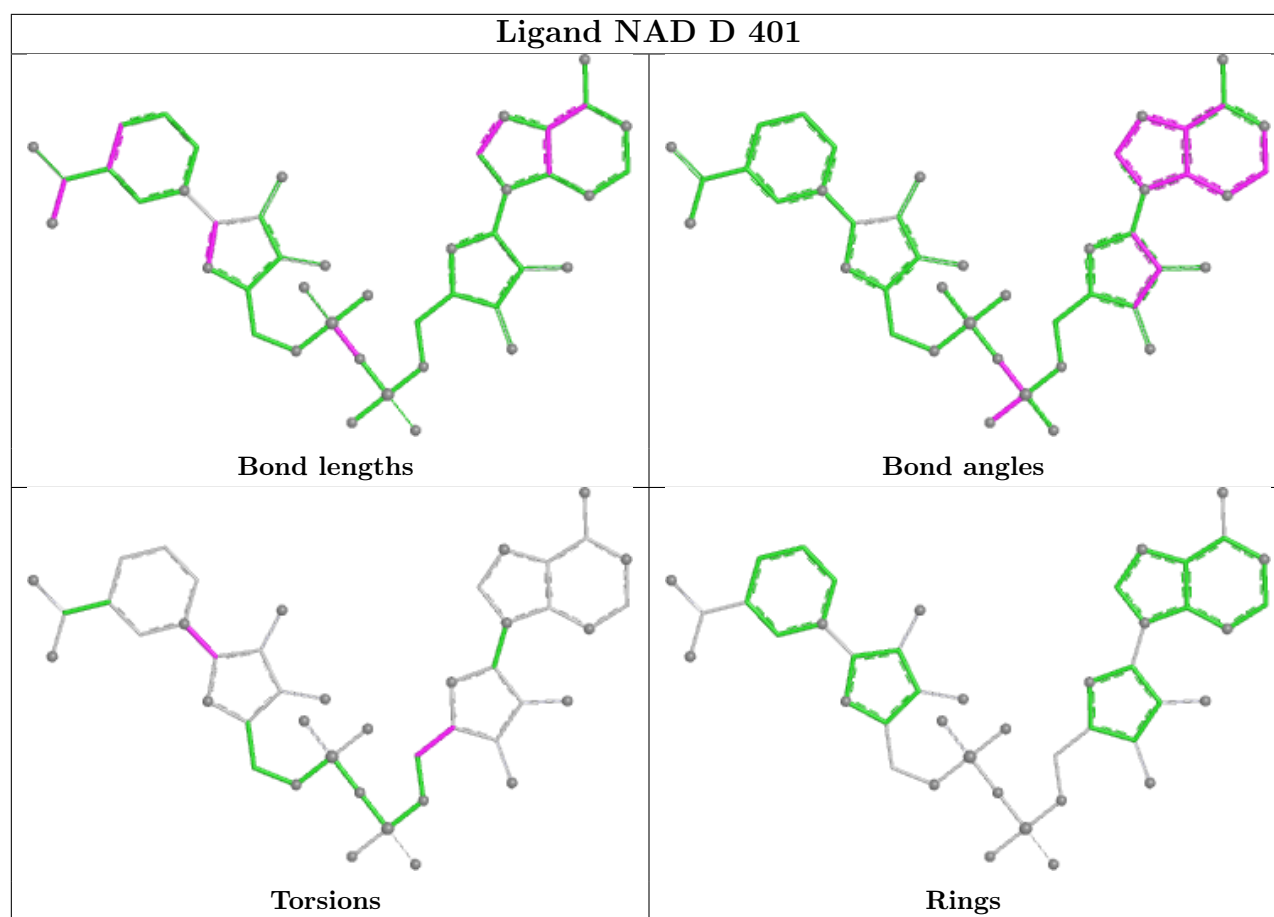
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	402	ISE	1	0
3	C	402	ISE	1	0

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](#)

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	366/380 (96%)	-0.02	13 (3%) 46 45	15, 20, 35, 52	0
1	B	366/380 (96%)	0.07	15 (4%) 41 40	15, 21, 37, 51	0
1	C	366/380 (96%)	0.17	21 (5%) 29 28	15, 22, 41, 58	0
1	D	362/380 (95%)	-0.02	10 (2%) 55 54	15, 21, 33, 50	0
All	All	1460/1520 (96%)	0.05	59 (4%) 42 41	15, 21, 37, 58	0

All (59) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	180	THR	11.0
1	B	177	TRP	9.2
1	B	175	TRP	8.8
1	B	176	SER	6.3
1	C	177	TRP	6.2
1	A	177	TRP	5.2
1	D	225	GLY	4.8
1	C	228	ALA	4.4
1	C	223	HIS	4.0
1	D	223	HIS	4.0
1	B	179	LEU	3.9
1	B	183	ASP	3.7
1	C	224	GLN	3.7
1	C	182	LYS	3.7
1	D	170	ASP	3.6
1	B	181	ARG	3.4
1	D	175	TRP	3.4
1	A	342	LYS	3.3
1	B	182	LYS	3.2
1	A	126	ASP	3.1
1	C	226	GLY	3.0

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Mol	Chain	Res	Type	RSRZ
1	A	181	ARG	3.0
1	B	231	GLU	2.9
1	C	180	THR	2.9
1	A	340	ARG	2.9
1	D	228	ALA	2.8
1	B	184	GLY	2.8
1	C	124	ALA	2.7
1	B	226	GLY	2.7
1	C	181	ARG	2.7
1	B	225	GLY	2.5
1	C	184	GLY	2.5
1	A	182	LYS	2.5
1	B	223	HIS	2.5
1	B	342	LYS	2.5
1	A	225	GLY	2.4
1	C	97	ALA	2.4
1	C	227	THR	2.4
1	A	223	HIS	2.4
1	C	183	ASP	2.4
1	A	183	ASP	2.3
1	C	229	ARG	2.3
1	D	227	THR	2.3
1	A	224	GLN	2.3
1	A	365	GLN	2.3
1	D	340	ARG	2.2
1	C	107	PRO	2.2
1	D	222	PRO	2.2
1	A	339	GLY	2.2
1	C	125	SER	2.2
1	C	175	TRP	2.2
1	C	69	GLU	2.2
1	B	49	ASP	2.1
1	D	220	ASP	2.1
1	C	225	GLY	2.1
1	D	226	GLY	2.1
1	A	174	PRO	2.1
1	C	126	ASP	2.1
1	C	179	LEU	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 [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

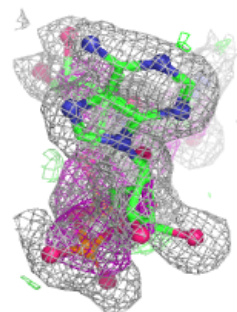
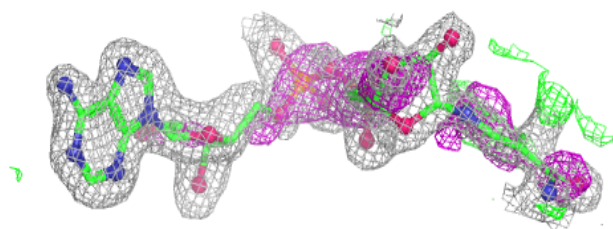
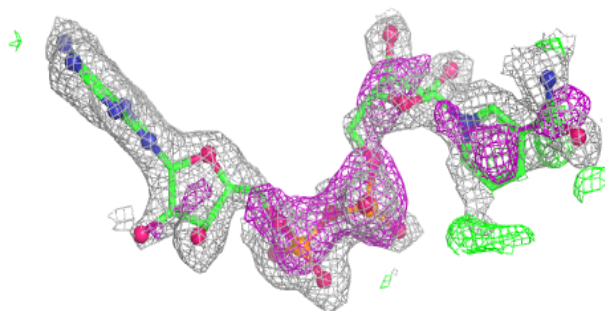
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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
2	NAD	A	500	44/44	0.76	0.14	31,34,40,44	0
2	NAD	C	401	44/44	0.77	0.13	33,37,44,47	0
2	NAD	B	401	44/44	0.85	0.11	25,33,35,37	0
3	ISE	C	402	12/12	0.89	0.09	22,22,24,24	0
3	ISE	B	402	12/12	0.90	0.09	21,22,23,23	0
2	NAD	D	401	44/44	0.91	0.10	25,29,33,35	0
3	ISE	D	402	12/12	0.92	0.07	21,22,23,23	0

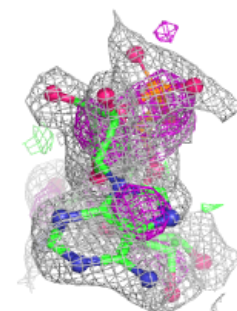
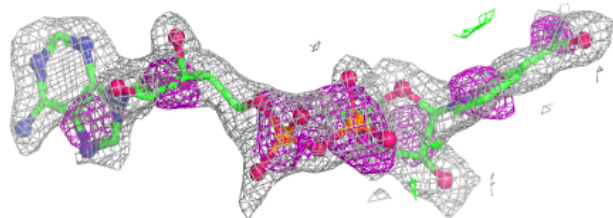
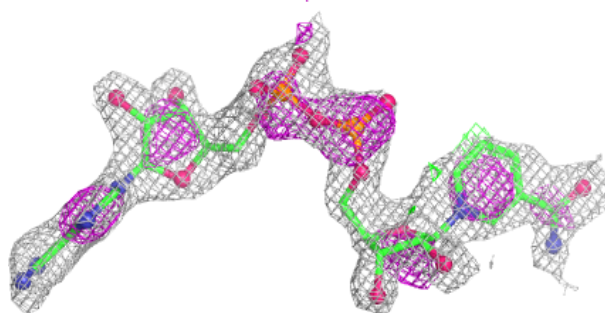
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around NAD A 500:

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

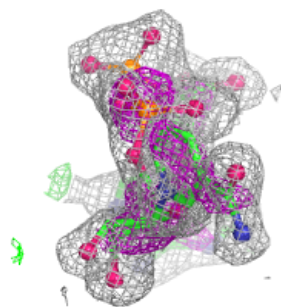
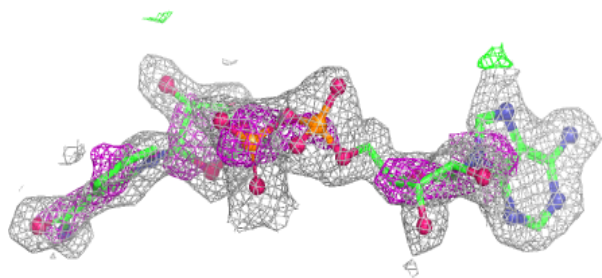
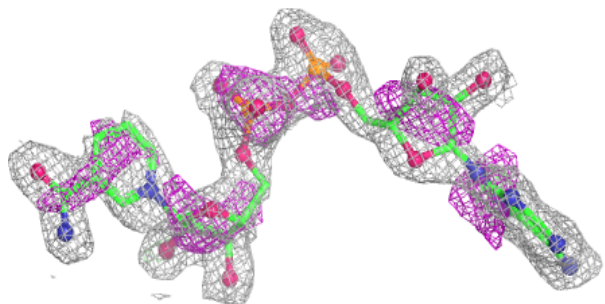
**Electron density around NAD C 401:**

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

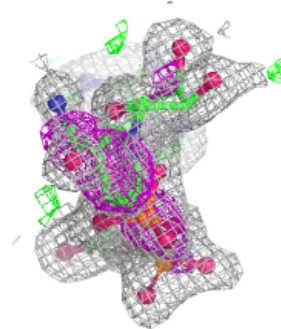
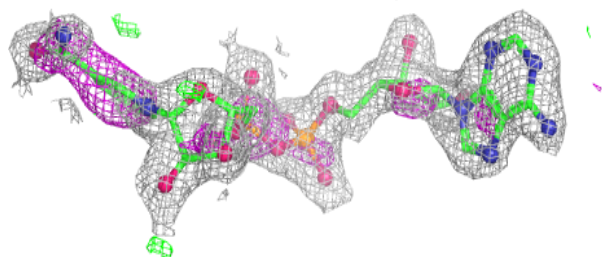
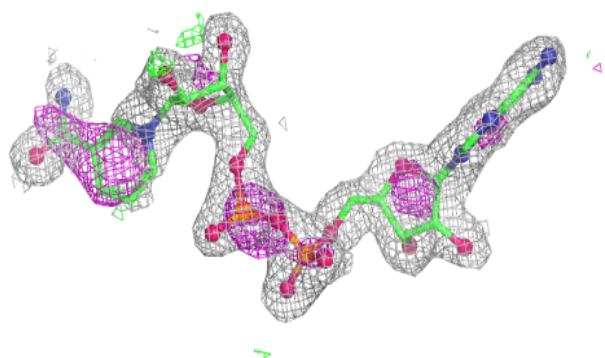


Electron density around NAD B 401:

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

**Electron density around NAD D 401:**

$2mF_o-DF_c$ (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.