



Full wwPDB X-ray Structure Validation Report ⓘ

Sep 16, 2023 – 06:32 PM EDT

PDB ID : 4X3Z
Title : Inosine 5'-monophosphate dehydrogenase from *Vibrio cholerae*, deletion mutant, in complex with XMP and NAD
Authors : Osipiuk, J.; MALTSEVA, N.; KIM, Y.; Mulligan, R.; MAKOWSKA-GRZYSKA, M.; Gu, M.; Anderson, W.F.; Joachimiak, A.; Center for Structural Genomics of Infectious Diseases (CSGID)
Deposited on : 2014-12-02
Resolution : 1.62 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](#) ①) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtrriage (Phenix) : 1.13
EDS : 2.35.1
buster-report : 1.1.7 (2018)
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.35.1

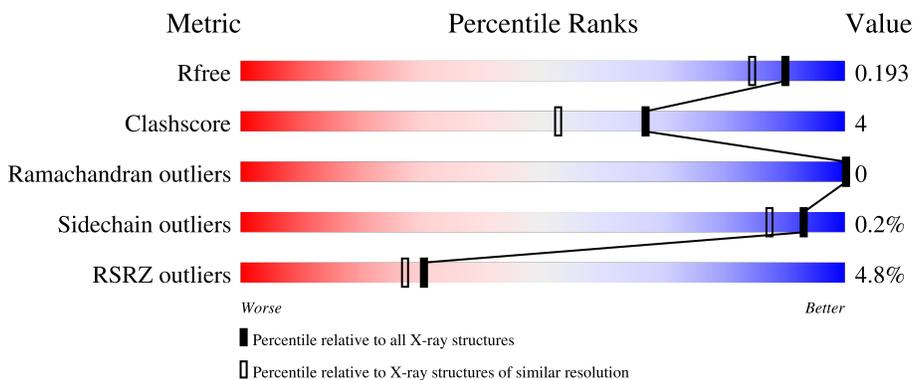
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.62 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}	130704	4693 (1.64-1.60)
Clashscore	141614	5002 (1.64-1.60)
Ramachandran outliers	138981	4888 (1.64-1.60)
Sidechain outliers	138945	4887 (1.64-1.60)
RSRZ outliers	127900	4609 (1.64-1.60)

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	366	
1	B	366	

2 Entry composition i

There are 7 unique types of molecules in this entry. The entry contains 5928 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.

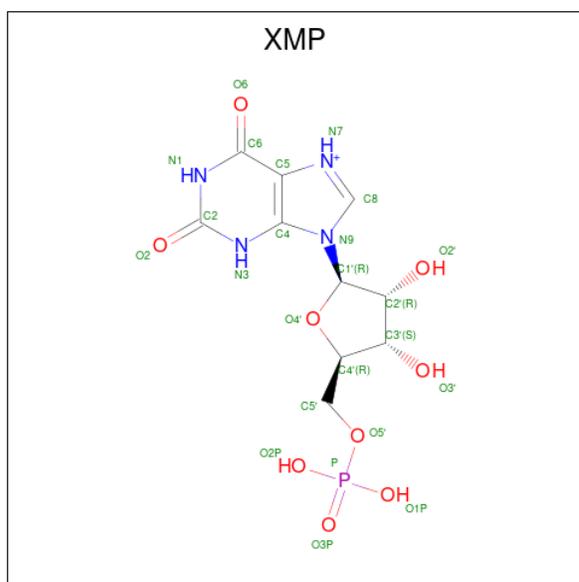
- Molecule 1 is a protein called Inosine-5'-monophosphate dehydrogenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	342	2555	1606	451	481	17	0	15	0
1	B	348	2645	1662	471	494	18	0	20	0

There are 12 discrepancies between the modelled and reference sequences:

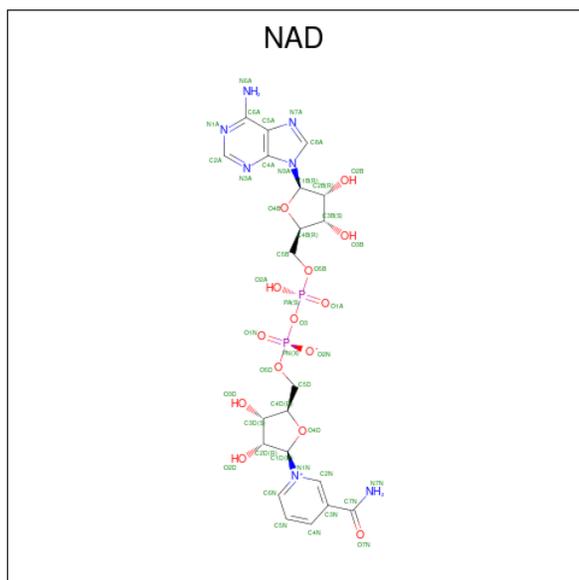
Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	SER	-	expression tag	UNP Q9KTW3
A	-1	ASN	-	expression tag	UNP Q9KTW3
A	0	ALA	-	expression tag	UNP Q9KTW3
A	101	SER	-	linker	UNP Q9KTW3
A	102	GLY	-	linker	UNP Q9KTW3
A	103	GLY	-	linker	UNP Q9KTW3
B	-2	SER	-	expression tag	UNP Q9KTW3
B	-1	ASN	-	expression tag	UNP Q9KTW3
B	0	ALA	-	expression tag	UNP Q9KTW3
B	101	SER	-	linker	UNP Q9KTW3
B	102	GLY	-	linker	UNP Q9KTW3
B	103	GLY	-	linker	UNP Q9KTW3

- Molecule 2 is XANTHOSINE-5'-MONOPHOSPHATE (three-letter code: XMP) (formula: C₁₀H₁₄N₄O₉P).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			24	10	4	9	1		
2	B	1	Total	C	N	O	P	0	0
			24	10	4	9	1		

- Molecule 3 is NICOTINAMIDE-ADENINE-DINUCLEOTIDE (three-letter code: NAD) (formula: $C_{21}H_{27}N_7O_{14}P_2$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	P	0	1
			54	22	7	21	4		

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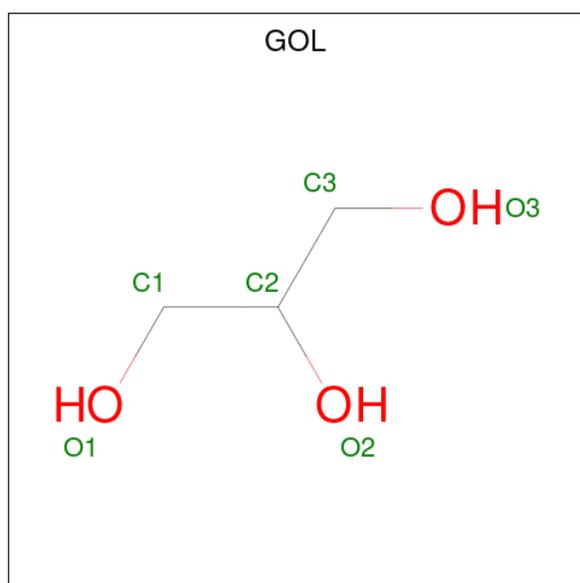
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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	N	O	P		
3	B	1	50	22	7	18	3	0	1

- Molecule 4 is POTASSIUM ION (three-letter code: K) (formula: K).

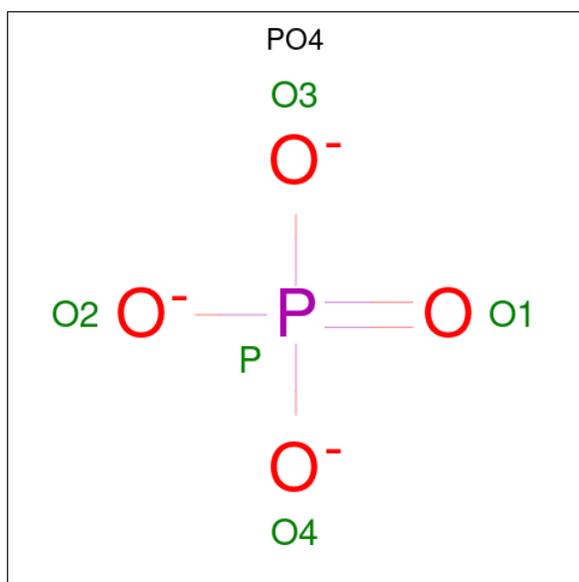
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total K 1 1	0	0
4	B	1	Total K 1 1	0	0

- Molecule 5 is GLYCEROL (three-letter code: GOL) (formula: C₃H₈O₃).



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

- Molecule 6 is PHOSPHATE ION (three-letter code: PO4) (formula: O₄P).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	B	1	Total	O	P	0	0
			5	4	1		

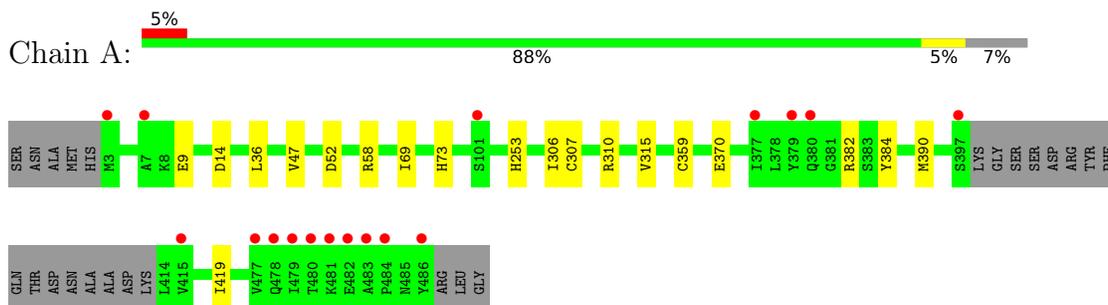
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	268	Total	O	0	0
			268	268		
7	B	295	Total	O	0	1
			295	295		

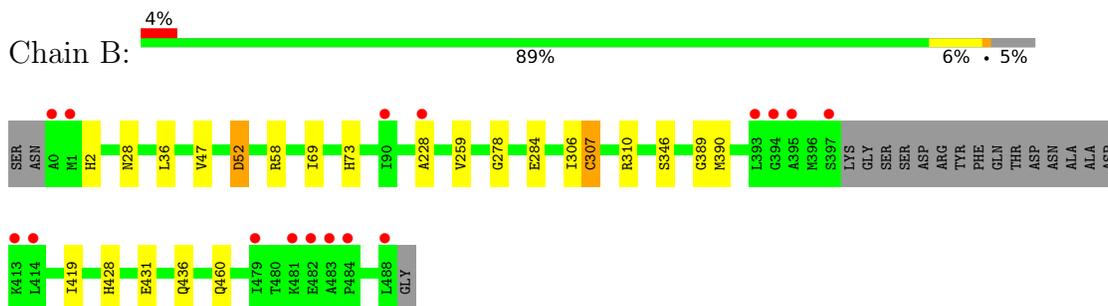
3 Residue-property plots [i](#)

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: Inosine-5'-monophosphate dehydrogenase



- Molecule 1: Inosine-5'-monophosphate dehydrogenase



4 Data and refinement statistics

Property	Value	Source
Space group	P 4 21 2	Depositor
Cell constants a, b, c, α , β , γ	91.26Å 91.26Å 171.21Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	44.20 – 1.62 44.09 – 1.62	Depositor EDS
% Data completeness (in resolution range)	99.9 (44.20-1.62) 99.9 (44.09-1.62)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.55 (at 1.62Å)	Xtrriage
Refinement program	REFMAC 5.8.0073	Depositor
R, R_{free}	0.152 , 0.182 0.166 , 0.193	Depositor DCC
R_{free} test set	4553 reflections (4.93%)	wwPDB-VP
Wilson B-factor (Å ²)	20.1	Xtrriage
Anisotropy	0.106	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 51.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	5928	wwPDB-VP
Average B, all atoms (Å ²)	24.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.58% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: XMP, NAD, K, GOL, PO4

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	0.70	0/2631	0.82	2/3556 (0.1%)
1	B	0.78	1/2740 (0.0%)	0.88	4/3696 (0.1%)
All	All	0.74	1/5371 (0.0%)	0.85	6/7252 (0.1%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	278	GLY	N-CA	8.81	1.59	1.46

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	52	ASP	CB-CG-OD1	7.71	125.24	118.30
1	A	9	GLU	N-CA-CB	6.28	121.90	110.60
1	A	14	ASP	CB-CG-OD1	6.02	123.72	118.30
1	B	58	ARG	NE-CZ-NH1	5.87	123.23	120.30
1	B	58	ARG	NE-CZ-NH2	-5.31	117.64	120.30
1	B	310	ARG	NE-CZ-NH1	-5.27	117.67	120.30

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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	2555	0	2663	17	0
1	B	2645	0	2777	21	0
2	A	24	0	12	3	0
2	B	24	0	12	5	0
3	A	54	0	10	3	0
3	B	50	0	6	1	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0
5	A	6	0	8	0	0
6	B	5	0	0	0	0
7	A	268	0	0	0	0
7	B	295	0	0	5	0
All	All	5928	0	5488	42	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 4.

All (42) 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:307:CYS:HB2	2:B:500:XMP:C2	1.93	0.98
1:B:390:MET:H	2:B:500:XMP:H7	1.25	0.79
1:A:58:ARG:NH2	1:A:370[B]:GLU:OE2	2.16	0.77
1:A:307[B]:CYS:HB2	2:A:500:XMP:C2	2.17	0.74
1:A:47:VAL:HG12	1:A:69[A]:ILE:HG22	1.70	0.73
1:A:390:MET:H	2:A:500:XMP:H7	1.33	0.73
1:A:307[A]:CYS:HB3	2:A:500:XMP:C2	2.21	0.71
1:B:36:LEU:HD22	1:B:69[C]:ILE:HG21	1.79	0.64
1:B:228:ALA:HB1	1:B:259:VAL:HG22	1.79	0.64
1:B:307:CYS:HB2	2:B:500:XMP:N1	2.16	0.60
1:B:47:VAL:HG12	1:B:69[C]:ILE:HG22	1.84	0.57
1:B:47:VAL:HG12	1:B:69[A]:ILE:HG22	1.88	0.56
1:B:307:CYS:CB	2:B:500:XMP:C2	2.78	0.54
1:B:47:VAL:HG12	1:B:69[C]:ILE:CG2	2.38	0.54
1:B:460[A]:GLN:NE2	7:B:603:HOH:O	2.41	0.52
1:B:307:CYS:SG	2:B:500:XMP:O2	2.69	0.51
1:A:310:ARG:HG2	1:A:315:VAL:O	2.11	0.51
1:A:36:LEU:HD22	1:A:69[A]:ILE:HG21	1.91	0.51
1:A:58:ARG:NH2	1:A:370[B]:GLU:CD	2.64	0.51
1:B:428:HIS:HB2	1:B:431[B]:GLU:OE1	2.12	0.50
1:A:306[A]:ILE:CD1	1:A:419:ILE:HD11	2.41	0.49
1:B:47:VAL:CG1	1:B:69[A]:ILE:HG22	2.43	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:2:HIS:ND1	7:B:878:HOH:O	2.35	0.48
1:B:284[B]:GLU:HG2	7:B:804:HOH:O	2.13	0.48
1:B:36:LEU:HD13	1:B:47:VAL:HG11	1.94	0.48
1:A:47:VAL:CG1	1:A:69[A]:ILE:HG22	2.41	0.47
1:A:58:ARG:NH2	1:A:370[B]:GLU:OE1	2.49	0.46
1:A:382:ARG:NE	1:A:384:TYR:OH	2.41	0.46
1:B:28[B]:ASN:ND2	7:B:601:HOH:O	2.32	0.45
1:B:419:ILE:C	1:B:419:ILE:HD12	2.37	0.44
1:A:306[A]:ILE:HD12	1:A:419:ILE:HD11	2.00	0.43
1:B:52:ASP:HA	1:B:73:HIS:CD2	2.55	0.42
1:A:253:HIS:HD2	3:A:501[B]:NAD:O2A	2.03	0.42
1:A:52:ASP:HA	1:A:73:HIS:CD2	2.55	0.41
1:A:47:VAL:HG13	1:A:359[B]:CYS:SG	2.60	0.41
3:A:501[B]:NAD:O2N	3:A:501[B]:NAD:C5B	2.68	0.41
1:A:306[A]:ILE:HD11	1:A:419:ILE:CD1	2.51	0.41
1:B:346:SER:HB3	1:B:436[A]:GLN:NE2	2.37	0.40
1:B:306:ILE:HD13	1:B:389:GLY:HA2	2.02	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	353/366 (96%)	342 (97%)	11 (3%)	0	100	100
1	B	365/366 (100%)	354 (97%)	11 (3%)	0	100	100
All	All	718/732 (98%)	696 (97%)	22 (3%)	0	100	100

There are no Ramachandran outliers to report.

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	270/274 (98%)	270 (100%)	0	100	100
1	B	281/274 (103%)	280 (100%)	1 (0%)	91	84
All	All	551/548 (100%)	550 (100%)	1 (0%)	93	87

All (1) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	307	CYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	2	HIS

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no monosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 10 ligands modelled in this entry, 2 are monoatomic - leaving 8 for Mogul analysis.

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
6	PO4	B	503	-	4,4,4	0.77	0	6,6,6	1.77	2 (33%)
2	XMP	A	500	-	20,26,26	1.48	3 (15%)	27,40,40	2.31	5 (18%)
3	NAD	B	501[A]	-	42,48,48	1.18	3 (7%)	50,73,73	1.59	9 (18%)
2	XMP	B	500	-	20,26,26	1.36	4 (20%)	27,40,40	2.01	7 (25%)
3	NAD	A	501[A]	-	42,48,48	0.95	3 (7%)	50,73,73	1.41	8 (16%)
5	GOL	A	503	-	5,5,5	0.35	0	5,5,5	0.21	0
3	NAD	B	501[B]	-	42,48,48	1.21	3 (7%)	50,73,73	1.64	8 (16%)
3	NAD	A	501[B]	-	42,48,48	0.96	3 (7%)	50,73,73	1.49	8 (16%)

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	XMP	A	500	-	-	0/6/26/26	0/3/3/3
3	NAD	B	501[A]	-	-	2/26/62/62	0/5/5/5
2	XMP	B	500	-	-	0/6/26/26	0/3/3/3
3	NAD	A	501[A]	-	-	2/26/62/62	0/5/5/5
5	GOL	A	503	-	-	2/4/4/4	-
3	NAD	B	501[B]	-	-	4/26/62/62	0/5/5/5
3	NAD	A	501[B]	-	-	4/26/62/62	0/5/5/5

All (19) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	500	XMP	O4'-C1'	4.71	1.47	1.41
3	B	501[A]	NAD	C5A-C4A	3.26	1.49	1.40
3	B	501[B]	NAD	C5A-C4A	3.26	1.49	1.40
3	B	501[A]	NAD	O4B-C1B	3.23	1.45	1.41
3	B	501[B]	NAD	O4B-C1B	3.23	1.45	1.41
2	B	500	XMP	O4'-C1'	3.17	1.45	1.41

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	501[A]	NAD	O4B-C1B	2.81	1.45	1.41
3	A	501[B]	NAD	O4B-C1B	2.81	1.45	1.41
2	A	500	XMP	C6-N1	-2.68	1.33	1.37
2	B	500	XMP	C2'-C1'	-2.60	1.49	1.53
3	B	501[A]	NAD	C2A-N1A	2.47	1.38	1.33
3	B	501[B]	NAD	C2A-N1A	2.47	1.38	1.33
3	A	501[A]	NAD	C2A-N3A	2.29	1.35	1.32
3	A	501[B]	NAD	C2A-N3A	2.29	1.35	1.32
2	A	500	XMP	C2-N3	-2.26	1.32	1.36
2	B	500	XMP	O2'-C2'	2.17	1.48	1.43
2	B	500	XMP	O6-C6	2.11	1.27	1.23
3	A	501[A]	NAD	C5A-C4A	2.00	1.46	1.40
3	A	501[B]	NAD	C5A-C4A	2.00	1.46	1.40

All (47) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	500	XMP	N3-C2-N1	8.23	120.68	115.90
2	B	500	XMP	C4-N3-C2	-5.45	119.59	128.11
2	B	500	XMP	N3-C2-N1	5.44	119.06	115.90
2	A	500	XMP	C4-N3-C2	-5.13	120.09	128.11
3	B	501[A]	NAD	O2N-PN-O1N	4.30	133.49	112.24
3	B	501[B]	NAD	O2N-PN-O1N	4.30	133.49	112.24
3	B	501[A]	NAD	C4A-C5A-N7A	-4.15	105.07	109.40
3	B	501[B]	NAD	C4A-C5A-N7A	-4.15	105.07	109.40
3	B	501[A]	NAD	C2A-N1A-C6A	3.92	125.47	118.75
3	B	501[B]	NAD	C2A-N1A-C6A	3.92	125.47	118.75
3	A	501[A]	NAD	N3A-C2A-N1A	-3.91	122.56	128.68
3	A	501[B]	NAD	N3A-C2A-N1A	-3.91	122.56	128.68
3	A	501[B]	NAD	O5B-C5B-C4B	-3.91	95.53	108.99
3	B	501[A]	NAD	N3A-C2A-N1A	-3.80	122.73	128.68
3	B	501[B]	NAD	N3A-C2A-N1A	-3.80	122.73	128.68
2	A	500	XMP	C6-N1-C2	-3.35	124.24	126.88
3	A	501[A]	NAD	C3D-C2D-C1D	3.33	106.00	100.98
3	A	501[B]	NAD	C3D-C2D-C1D	3.33	106.00	100.98
3	B	501[B]	NAD	PN-O3-PA	-3.31	121.46	132.83
3	B	501[A]	NAD	C1B-N9A-C4A	-3.27	120.90	126.64
3	B	501[B]	NAD	C1B-N9A-C4A	-3.27	120.90	126.64
2	B	500	XMP	P-O5'-C5'	3.09	126.82	118.30
6	B	503	PO4	O3-P-O2	3.08	117.85	107.97
3	A	501[B]	NAD	O5B-PA-O1A	3.07	121.07	109.07
2	A	500	XMP	C8-N7-C5	2.96	108.62	102.99

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	500	XMP	C8-N7-C5	2.88	108.47	102.99
3	B	501[A]	NAD	C5A-C6A-N1A	-2.79	114.02	120.35
3	B	501[B]	NAD	C5A-C6A-N1A	-2.79	114.02	120.35
6	B	503	PO4	O3-P-O1	-2.70	101.01	110.89
3	A	501[A]	NAD	O2N-PN-O1N	2.61	125.13	112.24
3	B	501[A]	NAD	N6A-C6A-N1A	2.42	123.60	118.57
3	B	501[B]	NAD	N6A-C6A-N1A	2.42	123.60	118.57
3	A	501[A]	NAD	C3N-C7N-N7N	2.41	120.64	117.75
3	A	501[B]	NAD	C3N-C7N-N7N	2.41	120.64	117.75
3	A	501[A]	NAD	PN-O3-PA	-2.40	124.60	132.83
2	B	500	XMP	O6-C6-C5	-2.39	119.70	124.37
3	A	501[A]	NAD	C3B-C2B-C1B	2.26	104.38	100.98
3	A	501[B]	NAD	C3B-C2B-C1B	2.26	104.38	100.98
2	B	500	XMP	O6-C6-N1	2.23	123.28	120.65
3	A	501[A]	NAD	C2N-C3N-C4N	-2.19	115.77	118.26
3	A	501[B]	NAD	C2N-C3N-C4N	-2.19	115.77	118.26
3	A	501[A]	NAD	N6A-C6A-N1A	2.16	123.05	118.57
3	A	501[B]	NAD	N6A-C6A-N1A	2.16	123.05	118.57
2	A	500	XMP	O2'-C2'-C1'	-2.13	102.97	110.85
2	B	500	XMP	O2'-C2'-C1'	-2.08	103.16	110.85
3	B	501[A]	NAD	C5B-C4B-C3B	-2.05	107.49	115.18
3	B	501[A]	NAD	O4B-C4B-C5B	2.01	115.97	109.37

There are no chirality outliers.

All (14) torsion outliers are listed below:

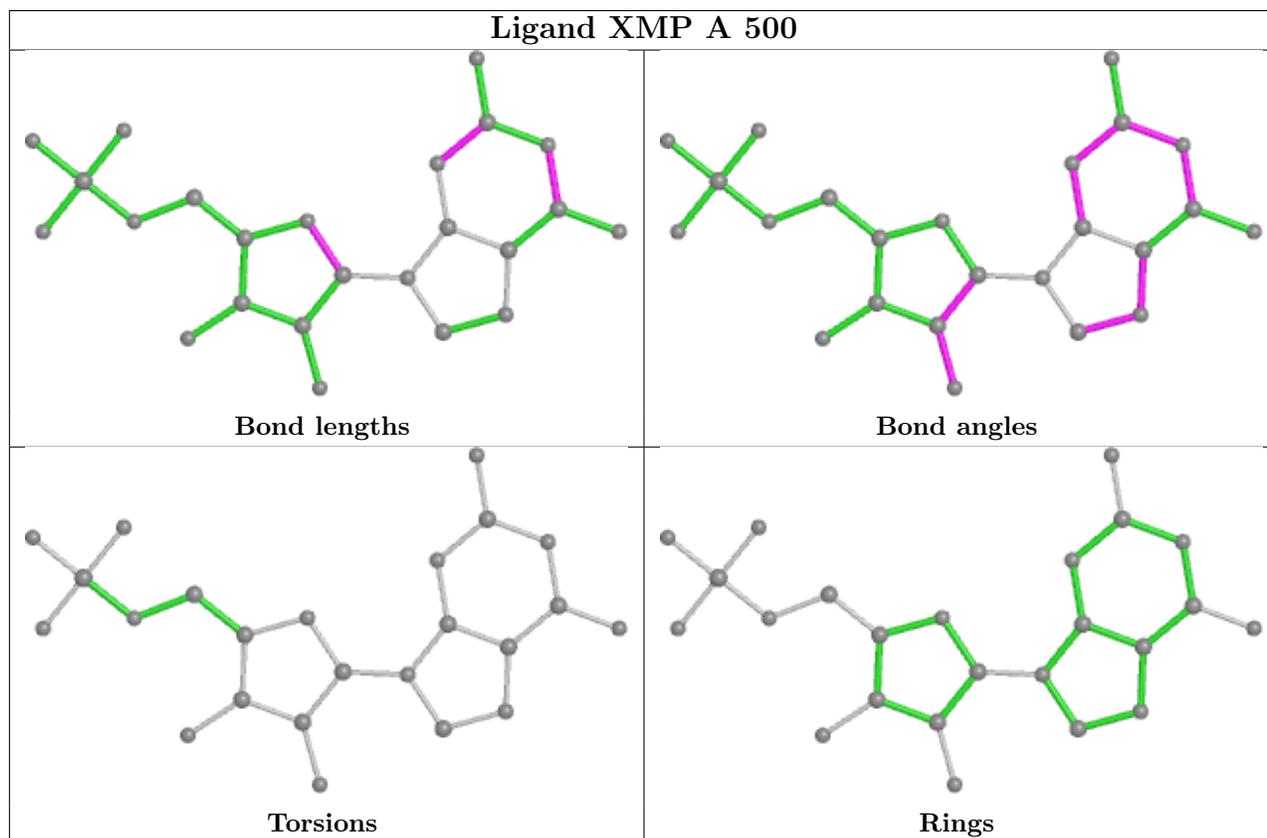
Mol	Chain	Res	Type	Atoms
3	A	501[B]	NAD	O4B-C4B-C5B-O5B
3	A	501[B]	NAD	C3B-C4B-C5B-O5B
3	B	501[A]	NAD	PN-O3-PA-O5B
3	B	501[B]	NAD	PN-O3-PA-O5B
3	A	501[B]	NAD	C4B-C5B-O5B-PA
3	A	501[A]	NAD	PN-O3-PA-O5B
3	A	501[B]	NAD	PN-O3-PA-O5B
3	B	501[B]	NAD	C5D-O5D-PN-O3
3	B	501[B]	NAD	C4B-C5B-O5B-PA
3	A	501[A]	NAD	PN-O3-PA-O1A
5	A	503	GOL	C1-C2-C3-O3
5	A	503	GOL	O2-C2-C3-O3
3	B	501[A]	NAD	C5D-O5D-PN-O2N
3	B	501[B]	NAD	C5D-O5D-PN-O2N

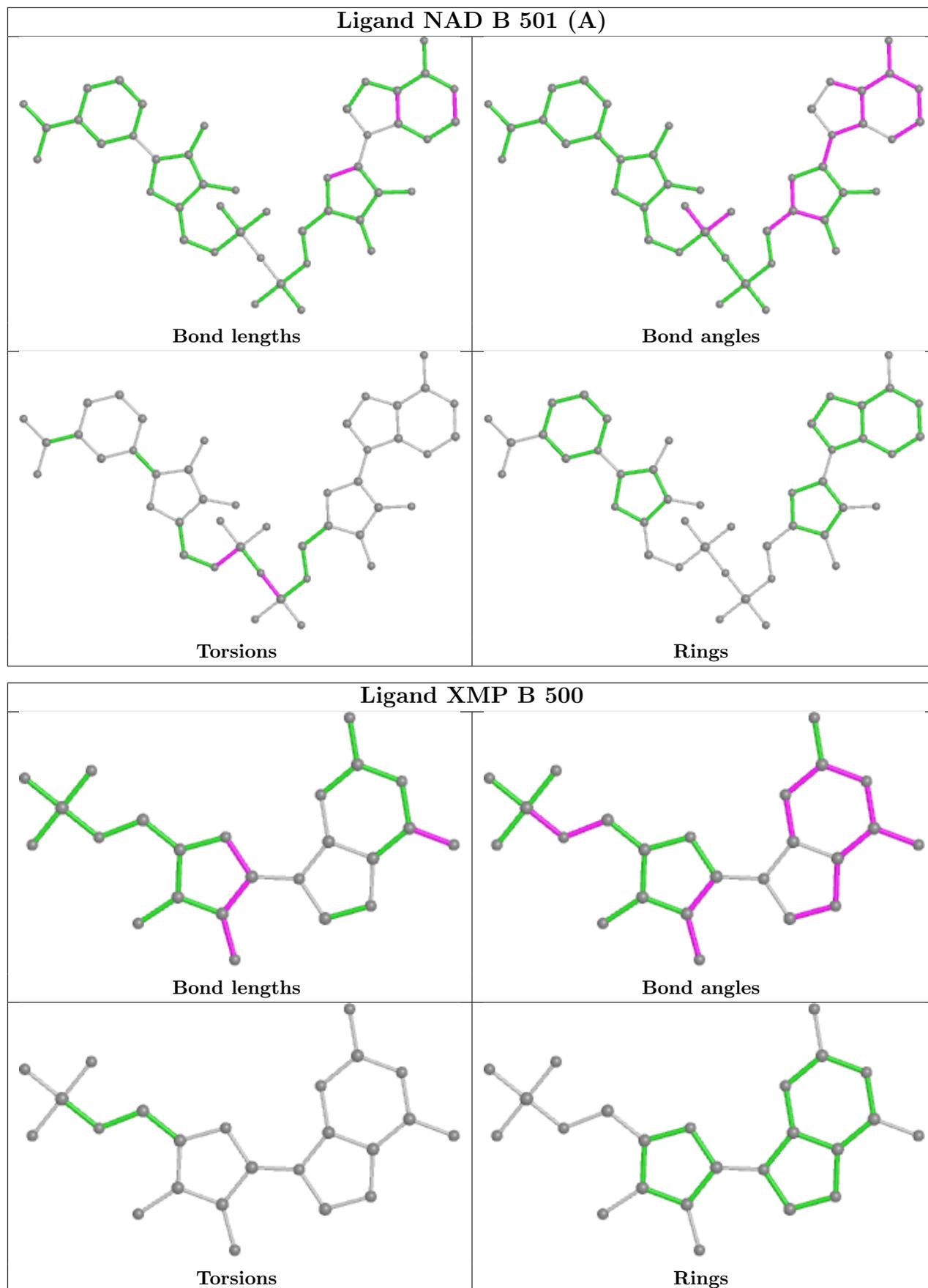
There are no ring outliers.

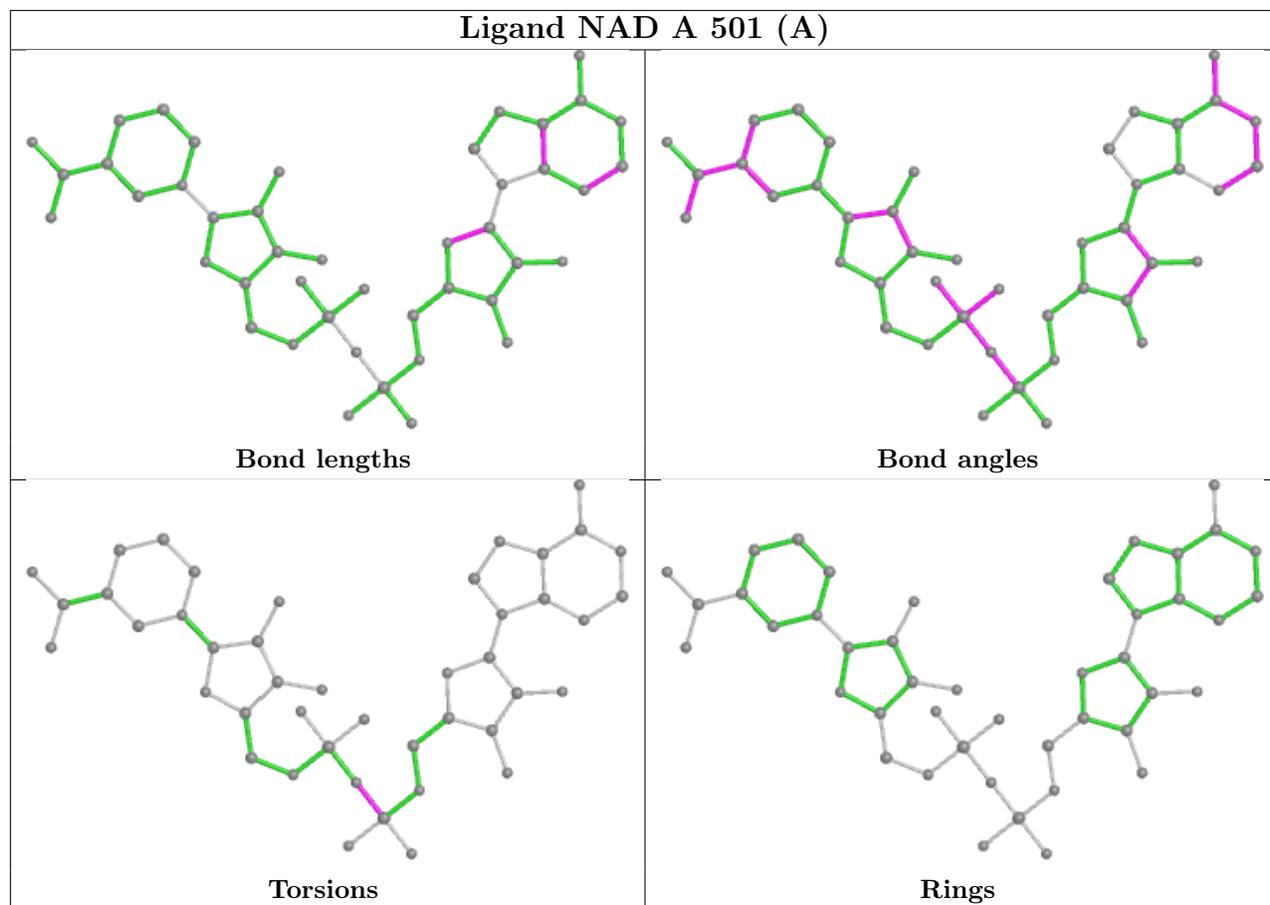
4 monomers are involved in 12 short contacts:

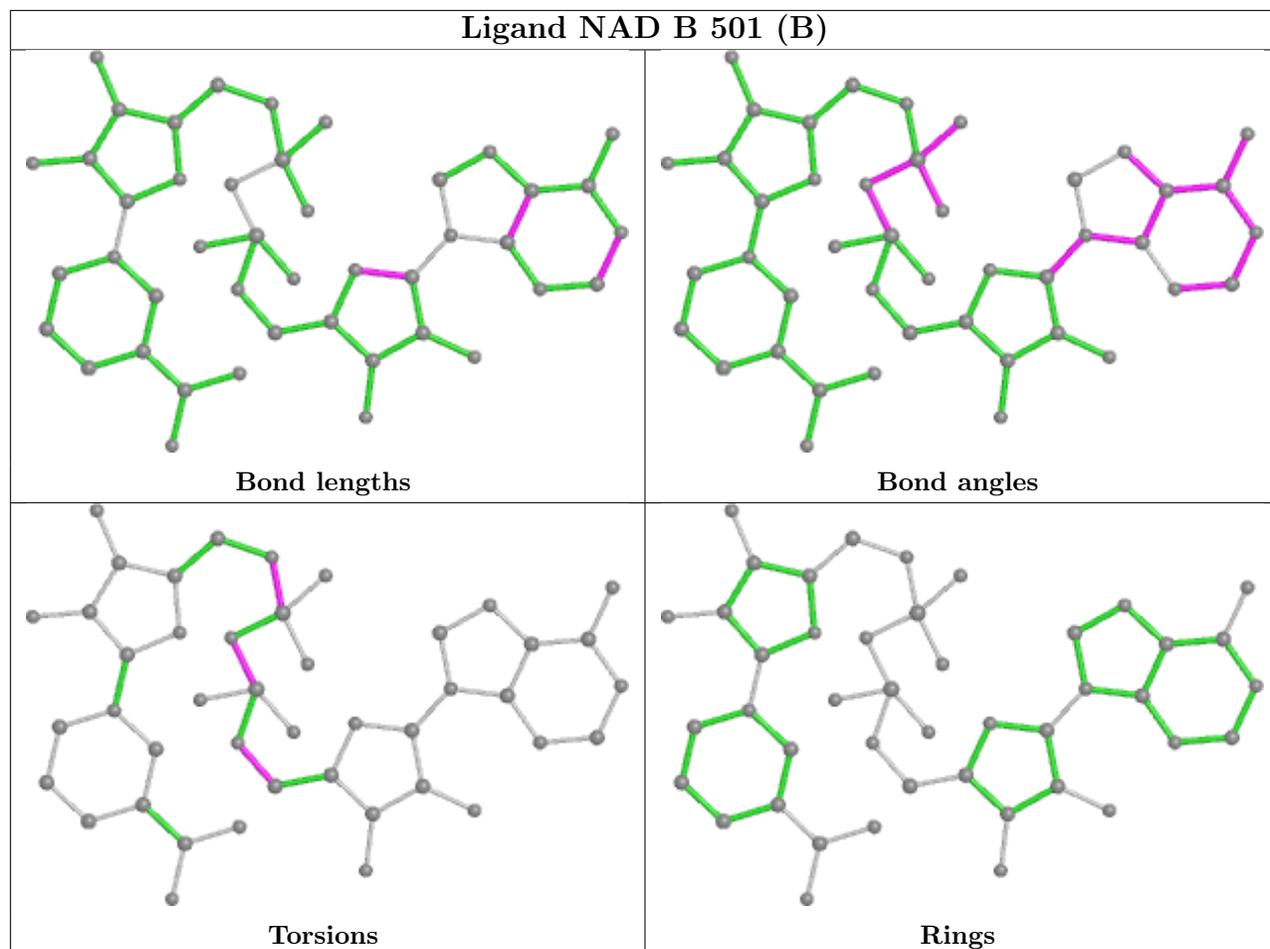
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	500	XMP	3	0
3	B	501[A]	NAD	1	0
2	B	500	XMP	5	0
3	A	501[B]	NAD	3	0

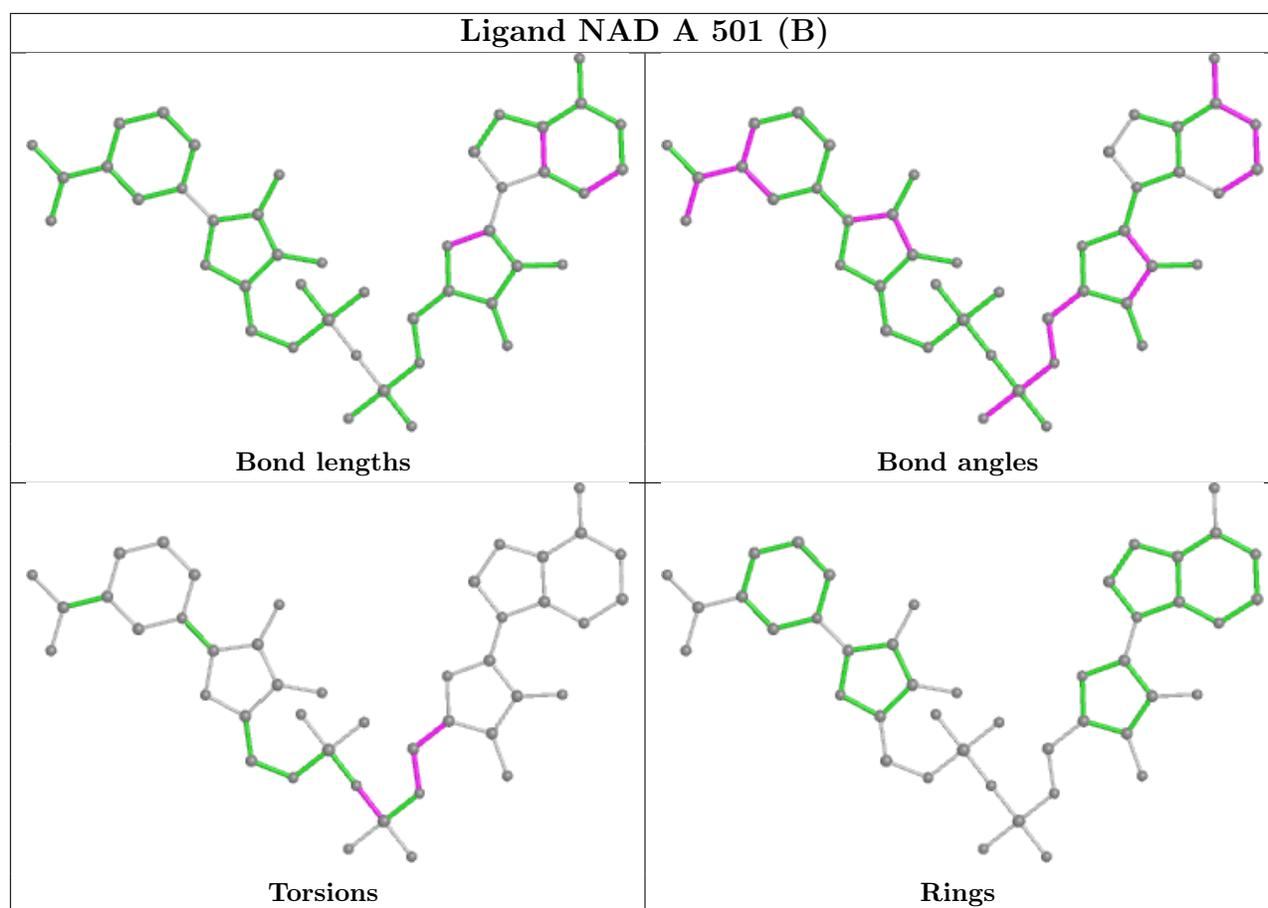
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. 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	342/366 (93%)	-0.02	17 (4%) 28 26	14, 21, 40, 69	2 (0%)
1	B	348/366 (95%)	0.12	16 (4%) 32 29	12, 20, 39, 60	2 (0%)
All	All	690/732 (94%)	0.05	33 (4%) 30 27	12, 21, 40, 69	4 (0%)

All (33) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	397	SER	5.7
1	A	379	TYR	5.6
1	B	0	ALA	4.6
1	B	228	ALA	4.3
1	B	413	LYS	4.1
1	B	1	MET	4.1
1	A	478	GLN	3.8
1	A	380	GLN	3.8
1	B	488	LEU	3.7
1	A	479	ILE	3.7
1	B	395	ALA	3.7
1	B	414	LEU	3.6
1	A	481	LYS	3.5
1	A	484	PRO	3.5
1	A	415	VAL	3.5
1	A	482	GLU	3.5
1	A	397	SER	3.2
1	B	481	LYS	3.1
1	A	486	TYR	3.0
1	A	483	ALA	2.8
1	B	484	PRO	2.6
1	B	393	LEU	2.6
1	A	480	THR	2.6
1	B	482	GLU	2.5

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Mol	Chain	Res	Type	RSRZ
1	A	3	MET	2.5
1	A	7	ALA	2.4
1	A	101[A]	SER	2.4
1	A	477	VAL	2.4
1	B	90	ILE	2.3
1	A	377	ILE	2.3
1	B	394	GLY	2.1
1	B	483	ALA	2.1
1	B	479	ILE	2.1

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

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

6.3 Carbohydrates [i](#)

There are no monosaccharides in this entry.

6.4 Ligands [i](#)

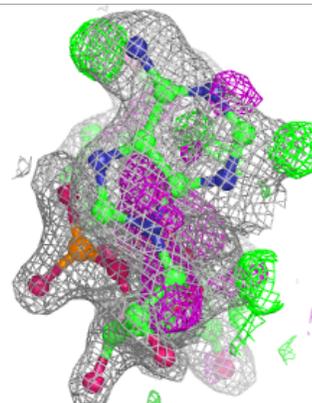
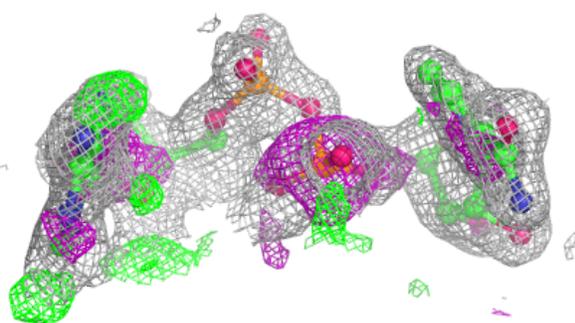
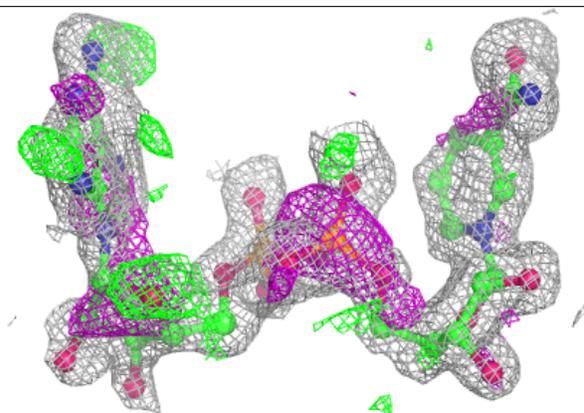
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	NAD	B	501[A]	44/44	0.75	0.25	25,32,45,51	6
3	NAD	B	501[B]	44/44	0.75	0.25	25,34,45,51	6
5	GOL	A	503	6/6	0.80	0.15	42,43,43,45	6
6	PO4	B	503	5/5	0.90	0.15	28,34,40,41	0
3	NAD	A	501[A]	44/44	0.92	0.12	20,24,47,53	10
3	NAD	A	501[B]	44/44	0.92	0.12	20,23,47,53	10
2	XMP	B	500	24/24	0.97	0.09	14,16,23,27	0
2	XMP	A	500	24/24	0.98	0.06	15,17,19,22	0
4	K	A	502	1/1	1.00	0.06	15,15,15,15	1
4	K	B	502	1/1	1.00	0.11	12,12,12,12	1

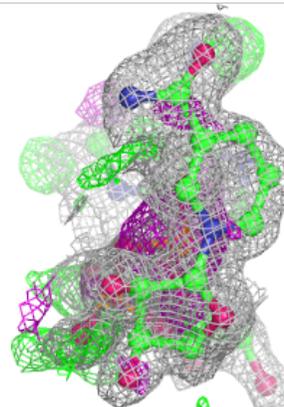
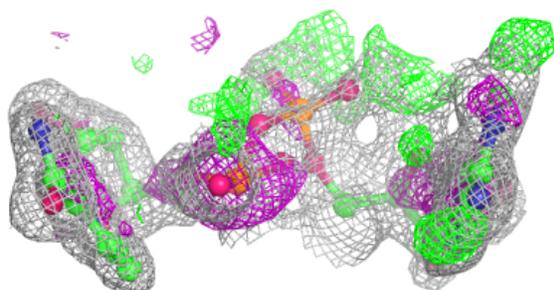
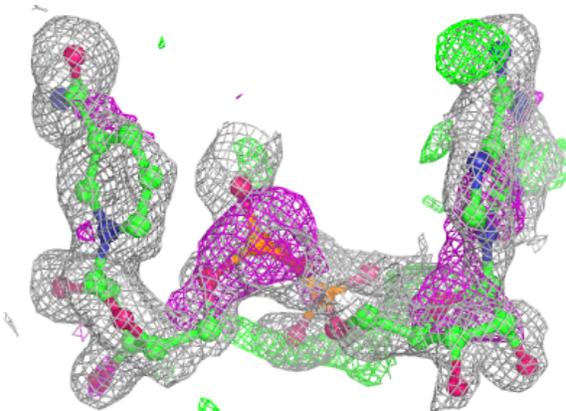
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around NAD B 501 (A):

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

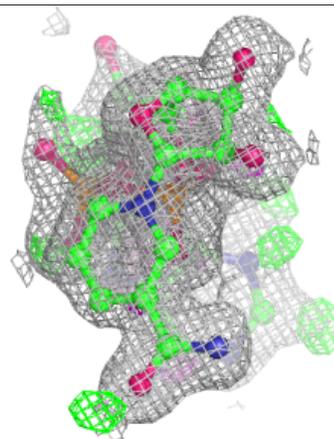
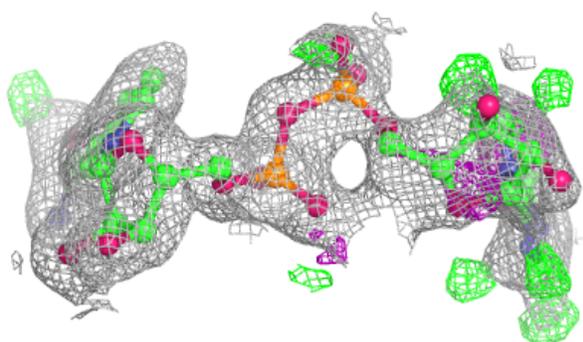
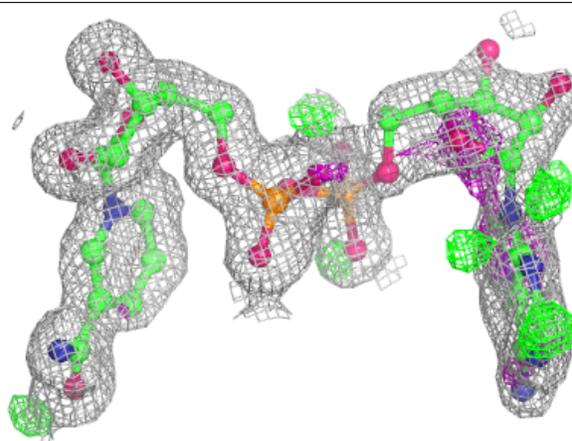
**Electron density around NAD B 501 (B):**

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



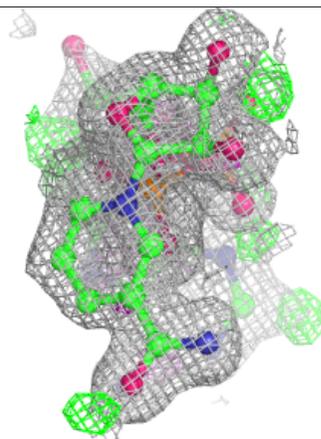
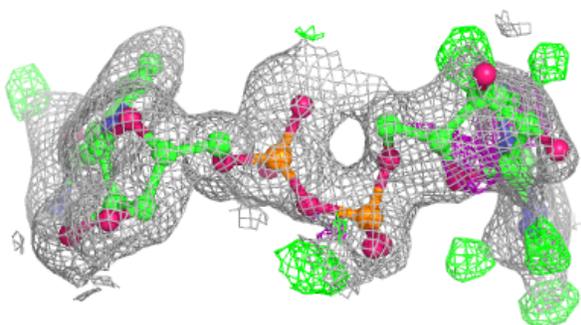
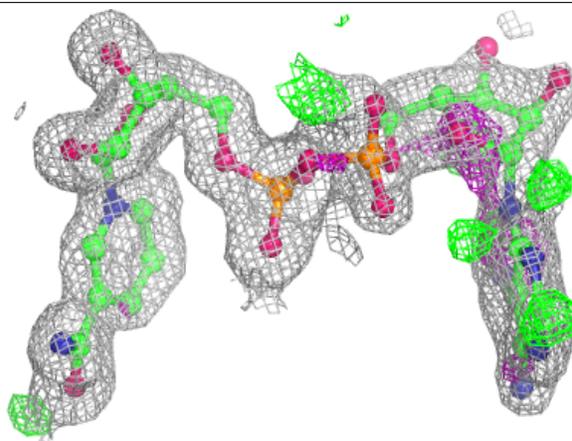
Electron density around NAD A 501 (A):

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

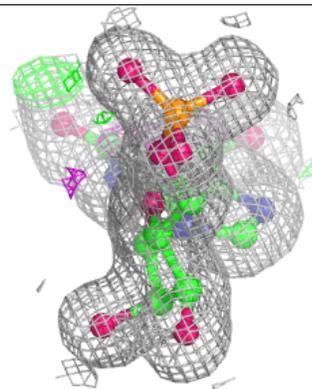
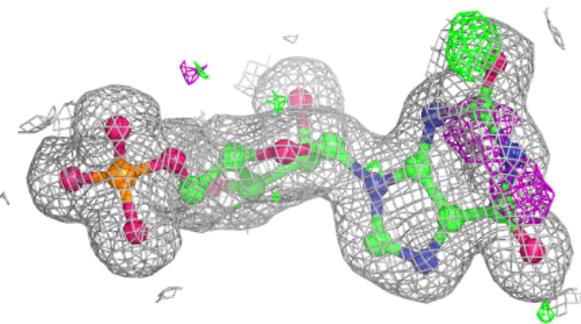
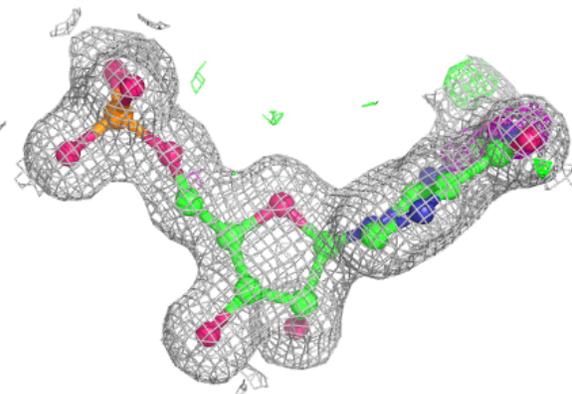


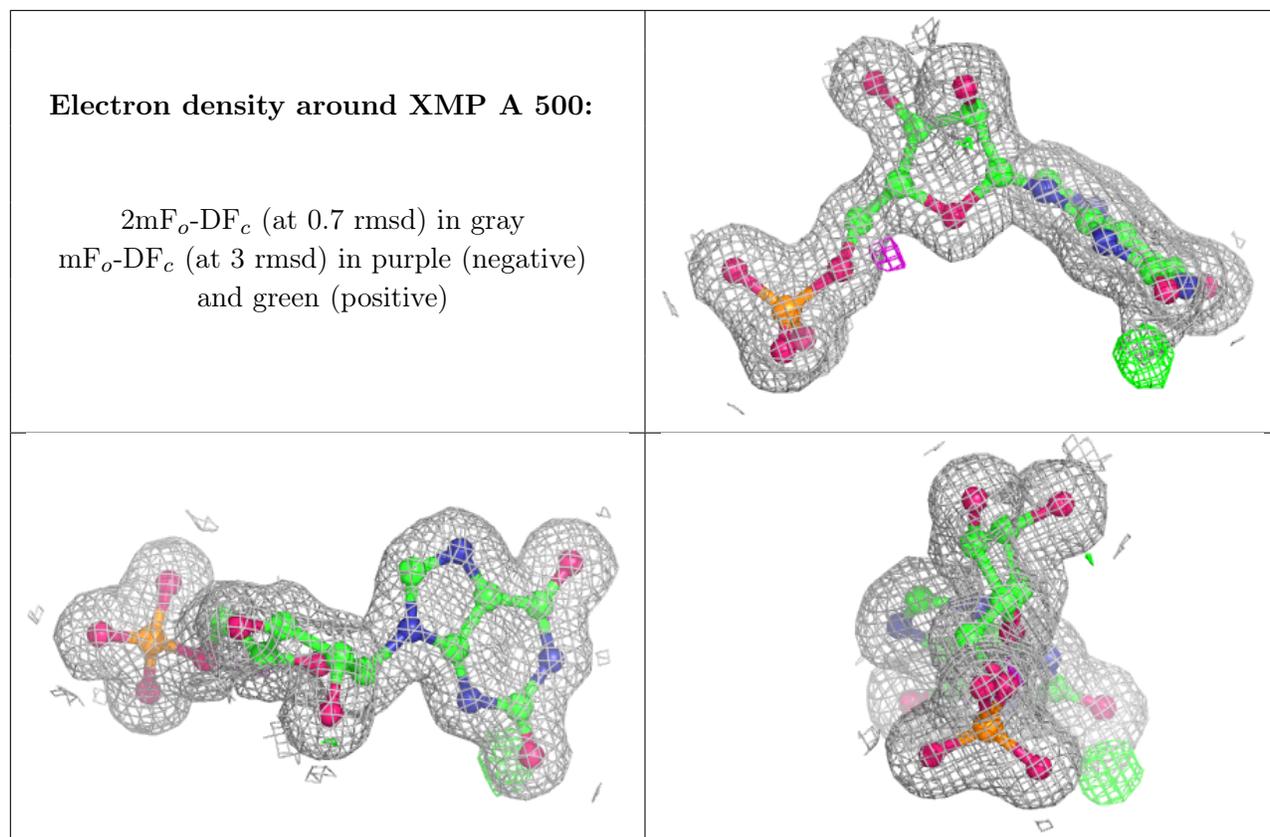
Electron density around NAD A 501 (B):

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

**Electron density around XMP B 500:**

$2mF_o-DF_c$ (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.