



Full wwPDB X-ray Structure Validation Report ⓘ

Sep 18, 2023 – 04:59 PM EDT

PDB ID : 5BJV
Title : X-ray structure of the PglF UDP-N-acetylglucosamine 4,6-dehydratase from *Campylobacter jejuni*, D396N/K397A variant in complex with UDP-N-acetylglucosamine
Authors : Riegert, A.S.; Thoden, J.B.; Holden, H.M.
Deposited on : 2017-09-12
Resolution : 1.80 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 : 4.02b-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

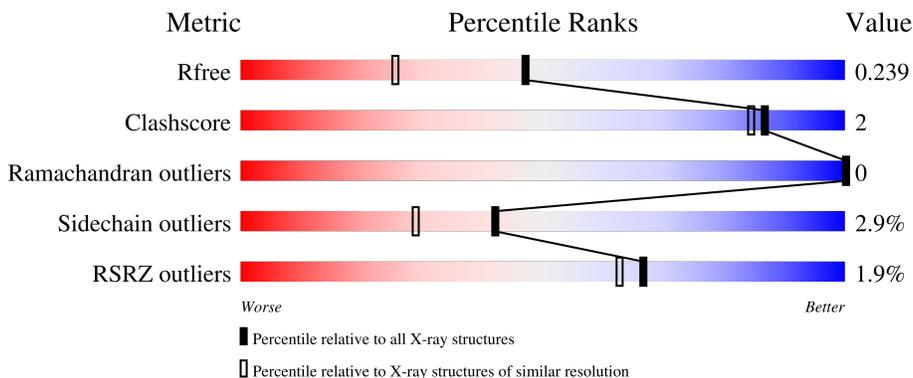
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.80 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	5950 (1.80-1.80)
Clashscore	141614	6793 (1.80-1.80)
Ramachandran outliers	138981	6697 (1.80-1.80)
Sidechain outliers	138945	6696 (1.80-1.80)
RSRZ outliers	127900	5850 (1.80-1.80)

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 6 unique types of molecules in this entry. The entry contains 6112 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 WlaL protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	339	2658	1701	439	503	15	0	3	0
1	B	338	2656	1701	440	500	15	0	4	0

There are 42 discrepancies between the modelled and reference sequences:

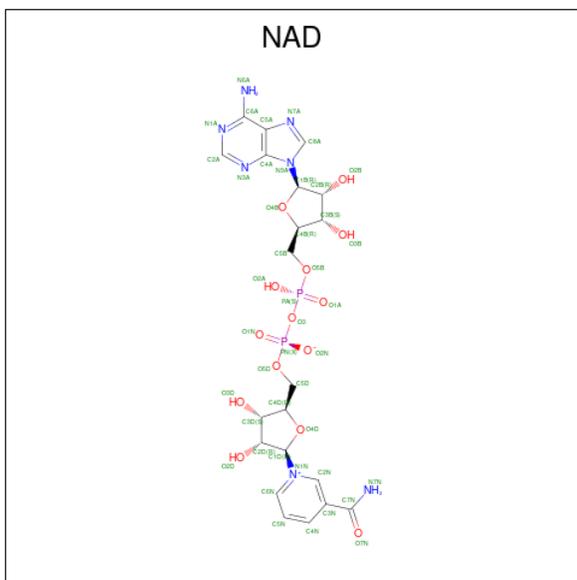
Chain	Residue	Modelled	Actual	Comment	Reference
A	396	ASN	ASP	engineered mutation	UNP O86159
A	397	ALA	LYS	engineered mutation	UNP O86159
A	591	ALA	-	expression tag	UNP O86159
A	592	ALA	-	expression tag	UNP O86159
A	593	GLY	-	expression tag	UNP O86159
A	594	PHE	-	expression tag	UNP O86159
A	595	ASN	-	expression tag	UNP O86159
A	596	ARG	-	expression tag	UNP O86159
A	597	ILE	-	expression tag	UNP O86159
A	598	PRO	-	expression tag	UNP O86159
A	599	ALA	-	expression tag	UNP O86159
A	600	ALA	-	expression tag	UNP O86159
A	601	ALA	-	expression tag	UNP O86159
A	602	LEU	-	expression tag	UNP O86159
A	603	GLU	-	expression tag	UNP O86159
A	604	HIS	-	expression tag	UNP O86159
A	605	HIS	-	expression tag	UNP O86159
A	606	HIS	-	expression tag	UNP O86159
A	607	HIS	-	expression tag	UNP O86159
A	608	HIS	-	expression tag	UNP O86159
A	609	HIS	-	expression tag	UNP O86159
B	396	ASN	ASP	engineered mutation	UNP O86159
B	397	ALA	LYS	engineered mutation	UNP O86159
B	591	ALA	-	expression tag	UNP O86159
B	592	ALA	-	expression tag	UNP O86159

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Chain	Residue	Modelled	Actual	Comment	Reference
B	593	GLY	-	expression tag	UNP O86159
B	594	PHE	-	expression tag	UNP O86159
B	595	ASN	-	expression tag	UNP O86159
B	596	ARG	-	expression tag	UNP O86159
B	597	ILE	-	expression tag	UNP O86159
B	598	PRO	-	expression tag	UNP O86159
B	599	ALA	-	expression tag	UNP O86159
B	600	ALA	-	expression tag	UNP O86159
B	601	ALA	-	expression tag	UNP O86159
B	602	LEU	-	expression tag	UNP O86159
B	603	GLU	-	expression tag	UNP O86159
B	604	HIS	-	expression tag	UNP O86159
B	605	HIS	-	expression tag	UNP O86159
B	606	HIS	-	expression tag	UNP O86159
B	607	HIS	-	expression tag	UNP O86159
B	608	HIS	-	expression tag	UNP O86159
B	609	HIS	-	expression tag	UNP O86159

- Molecule 2 is NICOTINAMIDE-ADENINE-DINUCLEOTIDE (three-letter code: NAD) (formula: C₂₁H₂₇N₇O₁₄P₂).



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

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total C O 4 2 2	0	0
4	A	1	Total C O 8 4 4	0	1
4	B	1	Total C O 4 2 2	0	0

- Molecule 5 is SODIUM ION (three-letter code: NA) (formula: Na).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	2	Total Na 2 2	0	0
5	B	2	Total Na 2 2	0	0

- Molecule 6 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	A	302	Total O 302 302	0	0
6	B	310	Total O 310 310	0	0

4 Data and refinement statistics i

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	69.82Å 108.47Å 108.96Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.00 – 1.80 29.39 – 1.80	Depositor EDS
% Data completeness (in resolution range)	96.4 (30.00-1.80) 96.5 (29.39-1.80)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	0.10	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.09 (at 1.80Å)	Xtrriage
Refinement program	REFMAC 5.8.0124	Depositor
R, R_{free}	0.192 , 0.232 0.203 , 0.239	Depositor DCC
R_{free} test set	3754 reflections (5.03%)	wwPDB-VP
Wilson B-factor (Å ²)	11.4	Xtrriage
Anisotropy	0.254	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 48.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.31$	Xtrriage
Estimated twinning fraction	0.000 for -h,l,k	Xtrriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	6112	wwPDB-VP
Average B, all atoms (Å ²)	18.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 50.81 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 6.1365e-05. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹Intensities estimated from amplitudes.

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

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: NAD, EDO, NA, UD1

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	0.59	0/2705	0.84	0/3651
1	B	0.57	0/2704	0.84	0/3650
All	All	0.58	0/5409	0.84	0/7301

There are no bond length outliers.

There are no bond angle outliers.

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	2658	0	2766	11	0
1	B	2656	0	2762	11	0
2	A	44	0	26	1	0
2	B	44	0	26	0	0
3	A	39	0	25	0	0
3	B	39	0	25	1	0
4	A	12	0	17	0	0
4	B	4	0	6	0	0
5	A	2	0	0	0	0
5	B	2	0	0	0	0
6	A	302	0	0	4	2

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
6	B	310	0	0	4	2
All	All	6112	0	5653	23	2

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 2.

All (23) 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:542[A]:GLN:OE1	6:B:801:HOH:O	1.87	0.91
1:A:249:ILE:HG12	1:A:506:MET:HG2	1.65	0.78
1:A:317:ASN:ND2	6:A:802:HOH:O	2.21	0.67
1:A:255:ARG:C	1:A:256:LYS:HG2	2.16	0.67
1:B:253:LEU:HD11	1:B:502:LEU:HD13	1.84	0.59
1:A:340:LEU:HD12	1:A:381:SER:OG	2.10	0.52
1:B:552:GLU:H	1:B:552:GLU:CD	2.12	0.52
1:A:394:SER:O	2:A:701:NAD:H6N	2.10	0.51
1:B:287:LEU:HD23	1:B:473:VAL:HG13	1.92	0.51
1:B:259:ASP:HB3	6:B:1046:HOH:O	2.09	0.51
1:A:255:ARG:O	1:A:255:ARG:HD2	2.12	0.50
1:A:346:GLU:OE1	6:A:801:HOH:O	2.20	0.47
1:A:322:LYS:NZ	6:A:806:HOH:O	2.37	0.47
1:B:357:VAL:HB	1:B:358:PRO:HD3	1.99	0.44
1:B:510:SER:O	1:B:512:ARG:HG2	2.18	0.44
1:A:287:LEU:HD23	1:A:473:VAL:HG13	2.00	0.44
1:A:255:ARG:NH2	6:A:808:HOH:O	2.41	0.43
1:B:444:LYS:NZ	6:B:824:HOH:O	2.52	0.43
1:B:583:PHE:CZ	1:B:585:HIS:HB2	2.55	0.42
1:B:463:VAL:HG12	1:B:498:LYS:HA	2.01	0.42
3:B:702:UD1:H8'3	6:B:860:HOH:O	2.20	0.41
1:B:258:LYS:HE3	1:B:258:LYS:HB2	1.89	0.41
1:A:567:GLN:O	1:A:568:ILE:HD13	2.21	0.40

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

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:A:1025:HOH:O	6:B:996:HOH:O[2_454]	2.09	0.11
6:A:858:HOH:O	6:B:890:HOH:O[2_454]	2.18	0.02

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	340/366 (93%)	332 (98%)	8 (2%)	0	100	100
1	B	340/366 (93%)	334 (98%)	6 (2%)	0	100	100
All	All	680/732 (93%)	666 (98%)	14 (2%)	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	300/316 (95%)	291 (97%)	9 (3%)	41	27
1	B	299/316 (95%)	290 (97%)	9 (3%)	41	27
All	All	599/632 (95%)	581 (97%)	18 (3%)	42	27

All (18) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	249	ILE
1	A	256	LYS
1	A	304	HIS
1	A	337	ASP
1	A	484	LYS
1	A	542[A]	GLN
1	A	542[B]	GLN
1	A	561	LYS

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Mol	Chain	Res	Type
1	A	578	LYS
1	B	304	HIS
1	B	317	ASN
1	B	320	LYS
1	B	346	GLU
1	B	392	MET
1	B	505	LYS
1	B	511	ASN
1	B	518	LYS
1	B	552	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	304	HIS
1	A	396	ASN
1	A	537	ASN
1	B	313	ASN
1	B	396	ASN
1	B	511	ASN

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 12 ligands modelled in this entry, 4 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
2	NAD	B	701	-	42,48,48	0.82	0	50,73,73	2.05	13 (26%)
4	EDO	A	703	-	3,3,3	0.60	0	2,2,2	0.40	0
4	EDO	A	704[A]	-	3,3,3	0.65	0	2,2,2	0.43	0
2	NAD	A	701	-	42,48,48	0.70	1 (2%)	50,73,73	1.64	12 (24%)
3	UD1	B	702	-	38,41,41	1.20	6 (15%)	57,62,62	1.43	8 (14%)
4	EDO	A	704[B]	5	3,3,3	0.52	0	2,2,2	0.28	0
4	EDO	B	703	-	3,3,3	0.59	0	2,2,2	0.47	0
3	UD1	A	702	-	38,41,41	0.99	3 (7%)	57,62,62	1.10	6 (10%)

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	701	-	-	1/26/62/62	0/5/5/5
4	EDO	A	703	-	-	0/1/1/1	-
4	EDO	A	704[A]	-	-	0/1/1/1	-
2	NAD	A	701	-	-	3/26/62/62	0/5/5/5
3	UD1	B	702	-	-	7/26/63/63	0/3/3/3
4	EDO	A	704[B]	5	-	1/1/1/1	-
4	EDO	B	703	-	-	1/1/1/1	-
3	UD1	A	702	-	-	3/26/63/63	0/3/3/3

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	702	UD1	O7'-C7'	3.42	1.31	1.23
3	A	702	UD1	C4-N3	-2.48	1.34	1.38
3	A	702	UD1	C6-C5	2.45	1.40	1.35
3	A	702	UD1	C2-N1	2.44	1.42	1.38
3	B	702	UD1	C2-N1	2.41	1.42	1.38
3	B	702	UD1	C4-N3	-2.41	1.34	1.38
3	B	702	UD1	C8'-C7'	-2.30	1.45	1.50
3	B	702	UD1	C6-C5	2.25	1.40	1.35

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	702	UD1	C2-N3	-2.20	1.34	1.38
2	A	701	NAD	C5A-C4A	2.17	1.46	1.40

All (39) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	701	NAD	O3D-C3D-C2D	-6.89	89.53	111.82
2	B	701	NAD	O3D-C3D-C4D	6.69	130.40	111.05
2	A	701	NAD	O2D-C2D-C3D	-5.20	94.99	111.82
2	B	701	NAD	O4D-C1D-C2D	-3.99	101.10	106.93
3	B	702	UD1	N3-C2-N1	3.86	120.02	114.89
2	A	701	NAD	N3A-C2A-N1A	-3.82	122.72	128.68
2	A	701	NAD	C6N-N1N-C2N	-3.63	118.66	121.97
2	B	701	NAD	C6N-N1N-C2N	-3.03	119.21	121.97
3	B	702	UD1	C8'-C7'-N2'	-3.03	110.97	116.10
3	B	702	UD1	C4-N3-C2	-2.99	122.64	126.58
2	A	701	NAD	O3D-C3D-C4D	-2.78	103.02	111.05
3	B	702	UD1	O7'-C7'-N2'	2.76	127.02	121.95
2	A	701	NAD	O4D-C1D-C2D	-2.76	102.90	106.93
2	B	701	NAD	N6A-C6A-N1A	2.74	124.26	118.57
3	B	702	UD1	C2'-N2'-C7'	2.72	129.79	123.18
2	B	701	NAD	N3A-C2A-N1A	-2.70	124.46	128.68
3	B	702	UD1	C5-C4-N3	2.65	118.81	114.84
2	B	701	NAD	C3D-C2D-C1D	2.53	104.79	100.98
2	B	701	NAD	C1B-N9A-C4A	-2.51	122.22	126.64
3	B	702	UD1	C3'-C2'-N2'	2.48	115.29	110.62
2	B	701	NAD	C5N-C4N-C3N	-2.42	117.48	120.34
3	A	702	UD1	C5-C4-N3	2.42	118.45	114.84
3	A	702	UD1	N3-C2-N1	2.41	118.09	114.89
2	A	701	NAD	O7N-C7N-C3N	-2.40	116.76	119.63
2	A	701	NAD	O2D-C2D-C1D	2.35	119.53	110.85
3	A	702	UD1	C1'-C2'-N2'	-2.32	107.02	111.00
3	A	702	UD1	O3A-PB-O1'	-2.31	97.82	102.48
3	B	702	UD1	C1'-C2'-N2'	-2.29	107.06	111.00
2	A	701	NAD	C2A-N1A-C6A	2.16	122.45	118.75
2	B	701	NAD	O2D-C2D-C3D	-2.15	104.86	111.82
2	B	701	NAD	O4D-C4D-C3D	2.07	109.20	105.11
2	A	701	NAD	N6A-C6A-N1A	2.06	122.84	118.57
2	A	701	NAD	C3N-C7N-N7N	2.05	120.22	117.75
2	B	701	NAD	O2A-PA-O1A	2.04	122.31	112.24
2	B	701	NAD	C5A-C6A-N6A	-2.03	117.27	120.35
3	A	702	UD1	C4-N3-C2	-2.01	123.93	126.58

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	701	NAD	C1B-N9A-C4A	-2.01	123.12	126.64
3	A	702	UD1	O2A-PA-O1A	2.01	122.15	112.24
2	A	701	NAD	C5N-C4N-C3N	-2.00	117.97	120.34

There are no chirality outliers.

All (16) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	702	UD1	C1'-O1'-PB-O3A
3	B	702	UD1	C1'-O1'-PB-O3A
3	B	702	UD1	C8'-C7'-N2'-C2'
3	B	702	UD1	O7'-C7'-N2'-C2'
3	B	702	UD1	C4'-C5'-C6'-O6'
4	A	704[B]	EDO	O1-C1-C2-O2
3	A	702	UD1	PB-O3A-PA-O5B
3	B	702	UD1	PB-O3A-PA-O5B
3	A	702	UD1	C4'-C5'-C6'-O6'
3	B	702	UD1	C1'-O1'-PB-O1B
3	B	702	UD1	O5'-C5'-C6'-O6'
2	B	701	NAD	O4B-C4B-C5B-O5B
4	B	703	EDO	O1-C1-C2-O2
2	A	701	NAD	C5D-O5D-PN-O3
2	A	701	NAD	O4B-C4B-C5B-O5B
2	A	701	NAD	C5D-O5D-PN-O1N

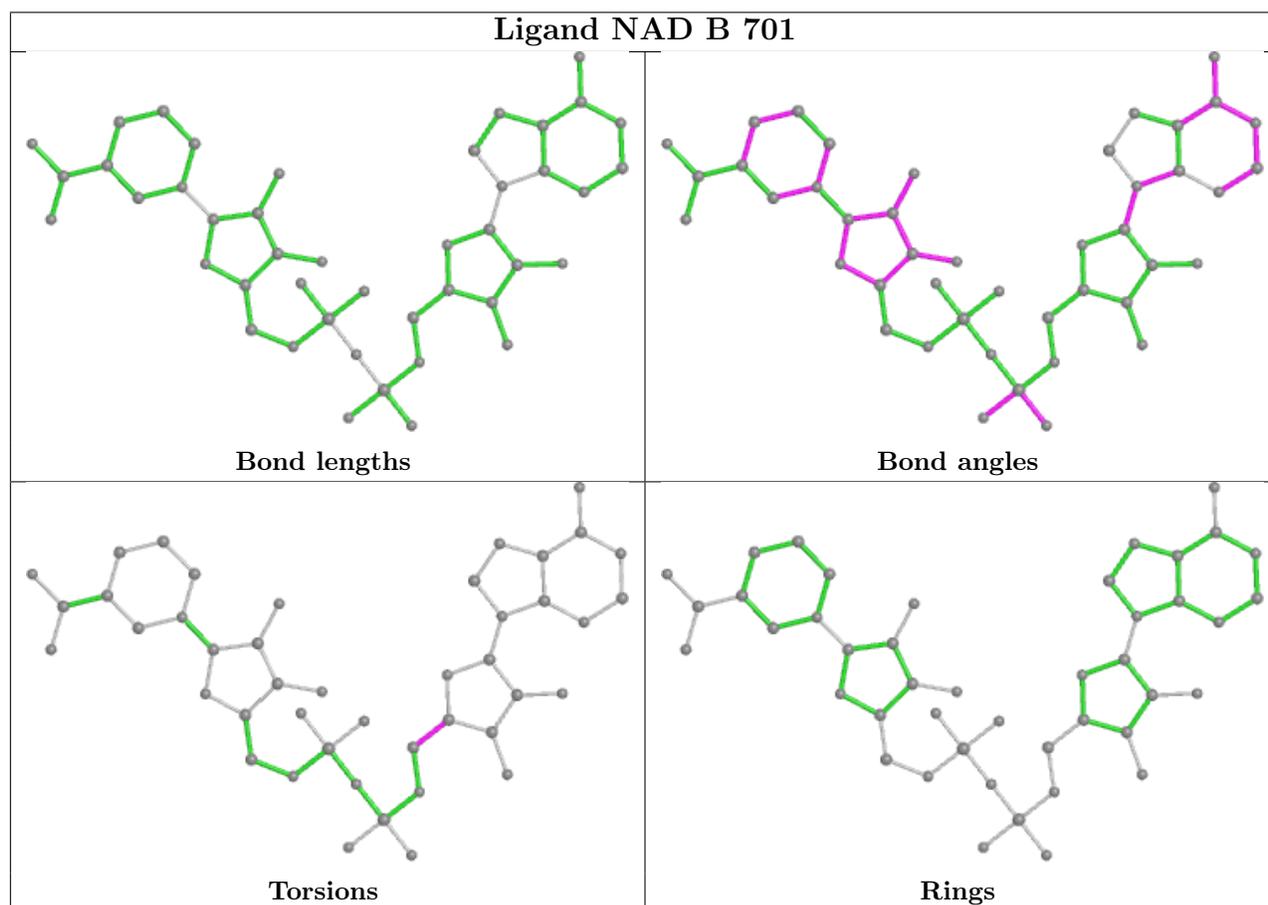
There are no ring outliers.

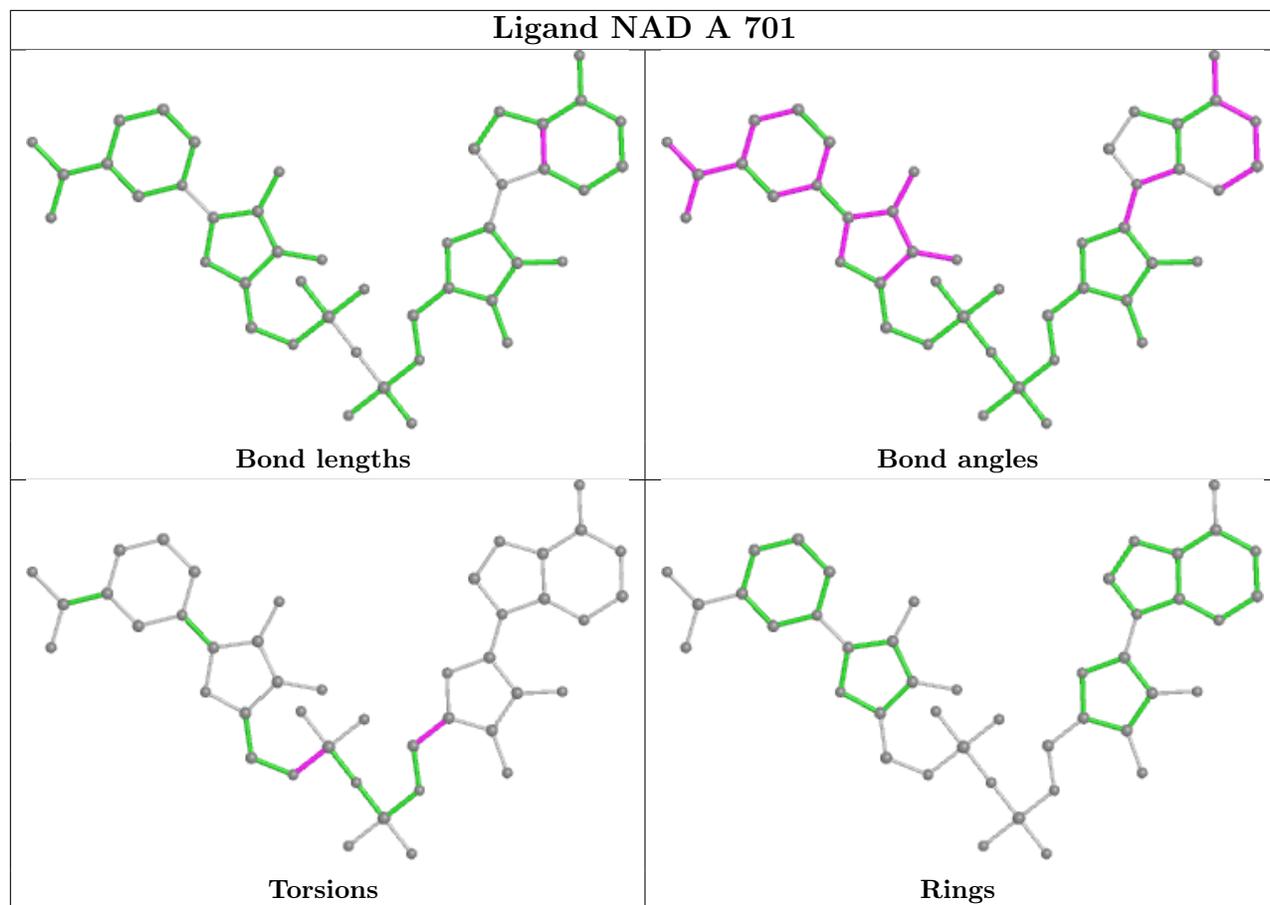
2 monomers are involved in 2 short contacts:

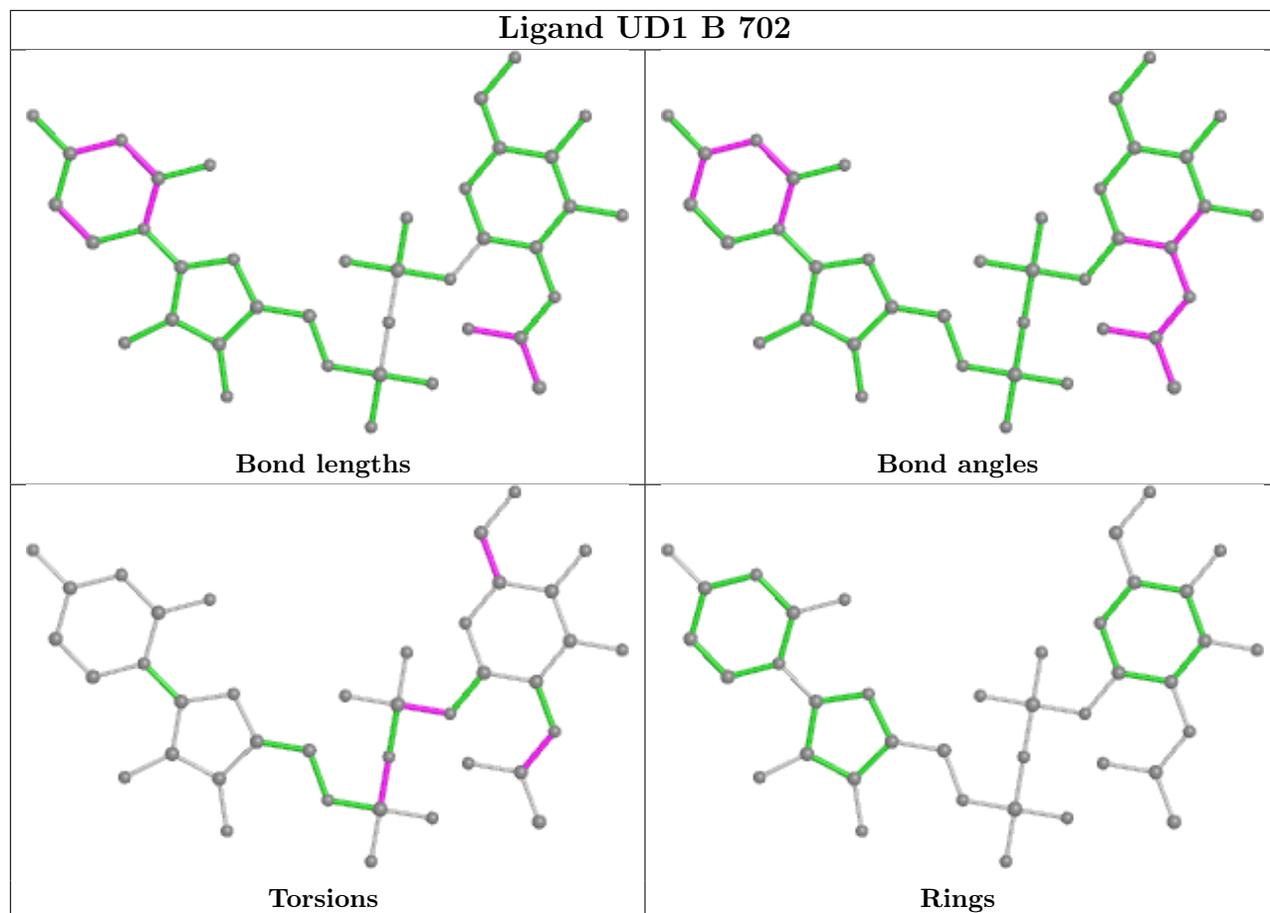
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	701	NAD	1	0
3	B	702	UD1	1	0

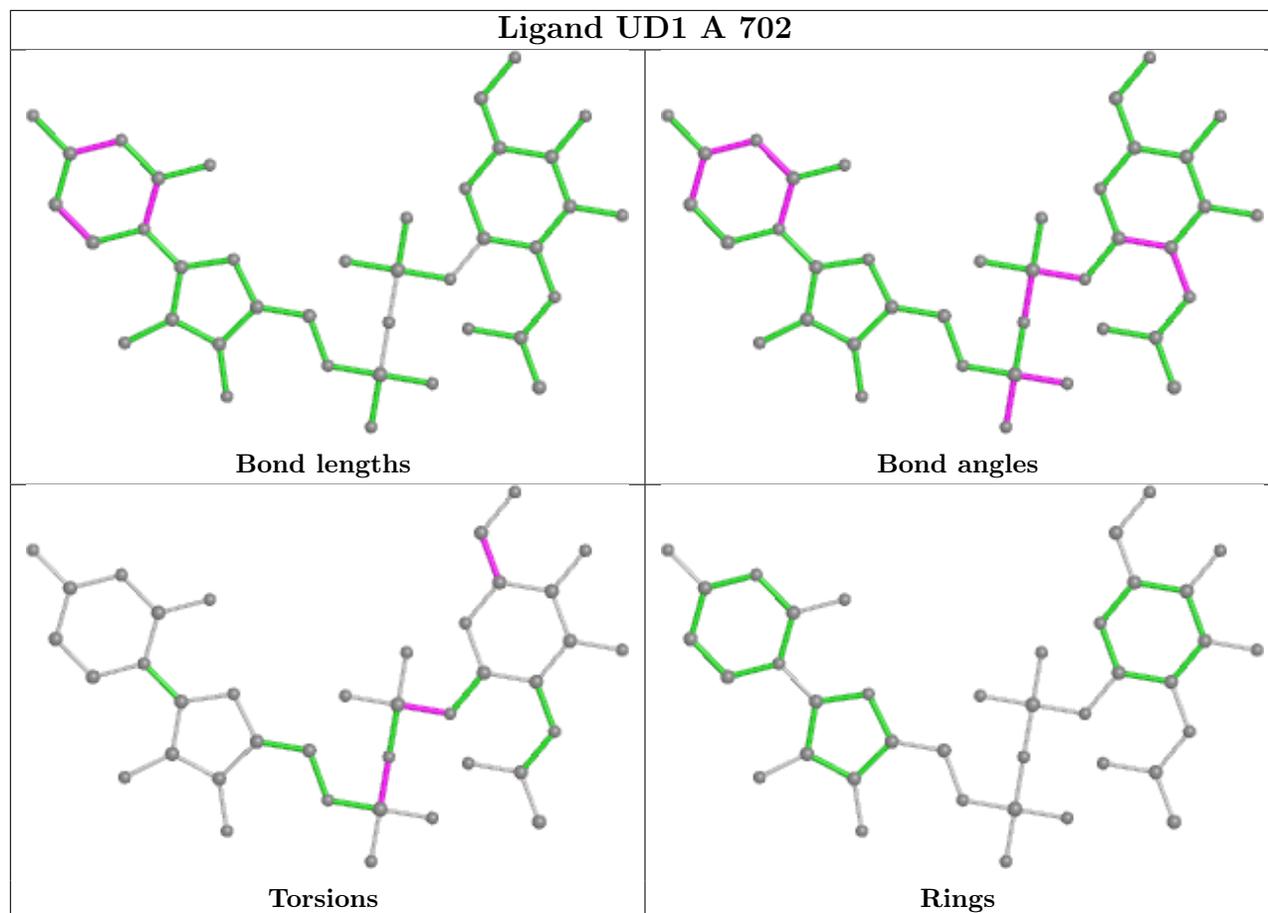
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. 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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	339/366 (92%)	-0.37	10 (2%) 51 46	6, 14, 41, 59	0
1	B	338/366 (92%)	-0.42	3 (0%) 84 82	6, 15, 39, 60	0
All	All	677/732 (92%)	-0.39	13 (1%) 66 63	6, 15, 40, 60	0

All (13) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	249	ILE	4.4
1	A	256	LYS	4.2
1	A	254	ALA	3.7
1	A	258	LYS	3.3
1	A	422	GLU	3.1
1	A	251	ASP	3.0
1	B	513	ASN	2.7
1	B	422	GLU	2.7
1	B	254	ALA	2.3
1	A	537	ASN	2.3
1	A	586	ASN	2.3
1	A	252	LEU	2.2
1	A	508	LEU	2.0

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

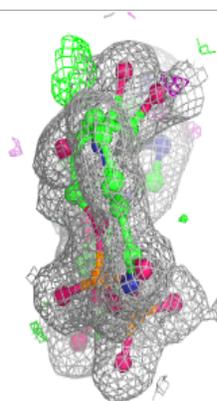
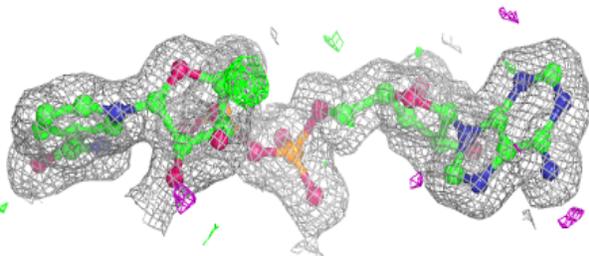
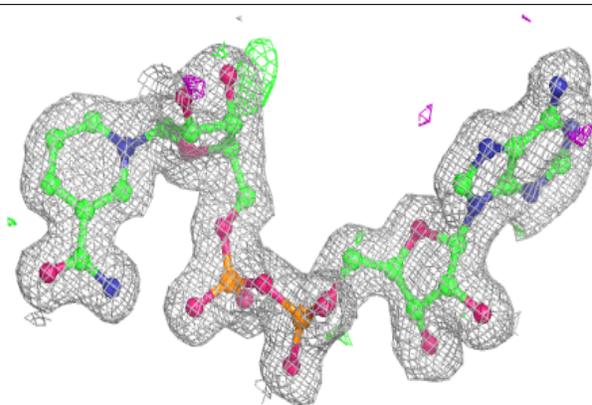
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
4	EDO	A	704[A]	4/4	0.90	0.17	15,17,18,18	4
4	EDO	A	704[B]	4/4	0.90	0.17	37,39,40,42	4
4	EDO	A	703	4/4	0.97	0.09	15,16,20,21	0
5	NA	A	705	1/1	0.97	0.13	23,23,23,23	0
5	NA	A	706	1/1	0.97	0.20	43,43,43,43	0
5	NA	B	704	1/1	0.97	0.13	30,30,30,30	0
5	NA	B	705	1/1	0.97	0.19	31,31,31,31	0
2	NAD	B	701	44/44	0.98	0.08	5,7,18,22	0
3	UD1	A	702	39/39	0.98	0.07	6,8,10,10	0
3	UD1	B	702	39/39	0.98	0.08	6,10,14,15	0
4	EDO	B	703	4/4	0.98	0.06	15,15,15,15	0
2	NAD	A	701	44/44	0.99	0.09	5,7,13,15	0

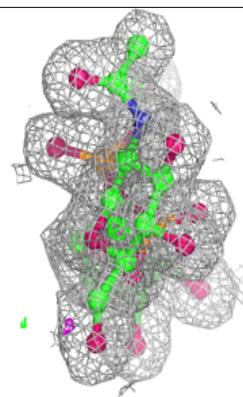
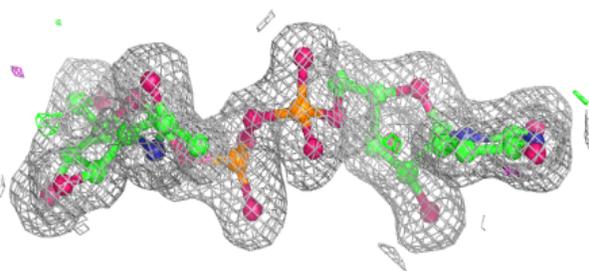
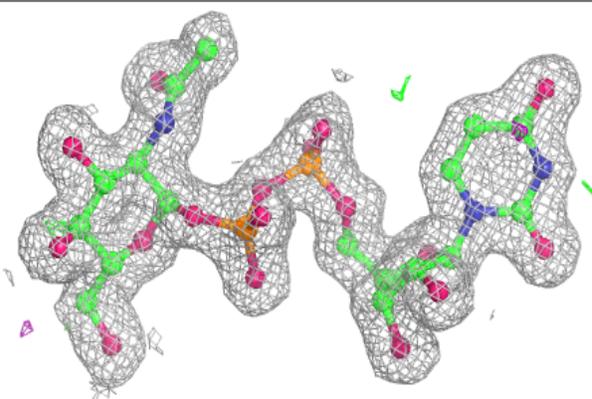
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around NAD B 701:

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

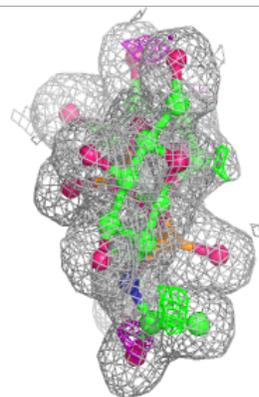
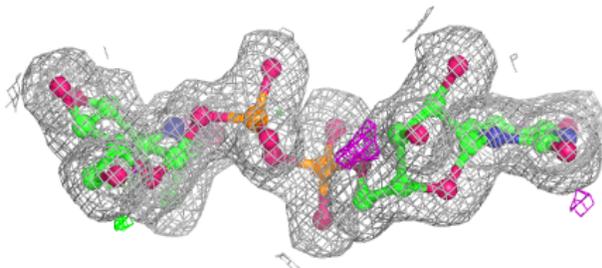
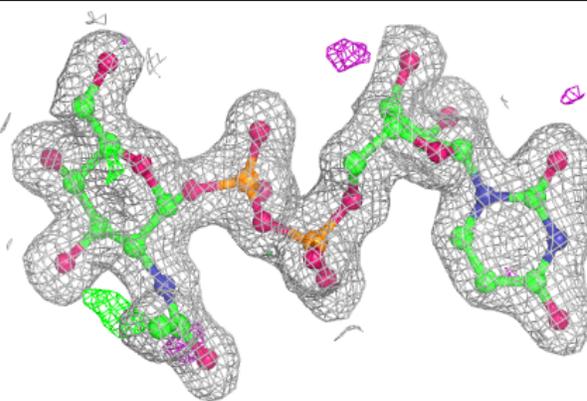
**Electron density around UD1 A 702:**

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

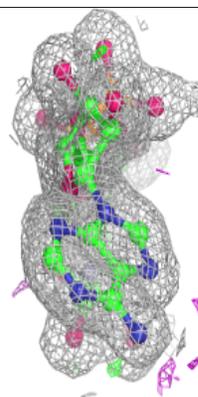
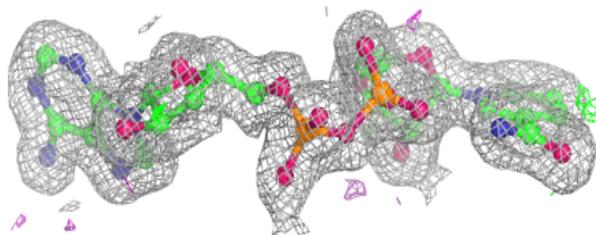
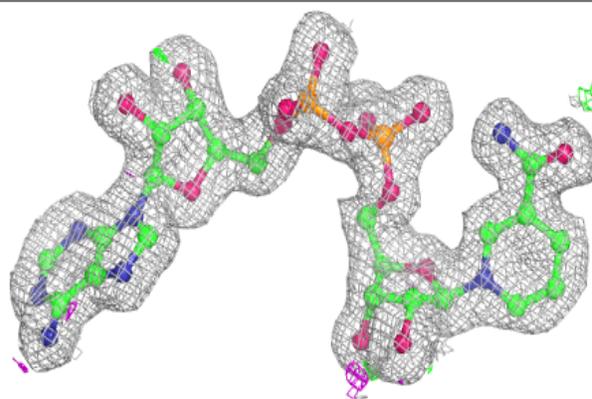


Electron density around UD1 B 702:

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

**Electron density around NAD A 701:**

$2mF_o-DF_c$ (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.