



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 3, 2025 – 03:43 PM JST

PDB ID : 8W9S
Title : NAD-dependent dehydrogenase
Authors : Niu, M.; Shen, Y.M.
Deposited on : 2023-09-05
Resolution : 2.09 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)
Xtriage (Phenix) : 1.21
EDS : 3.0
buster-report : 1.1.7 (2018)
Percentile statistics : 20231227.v01 (using entries in the PDB archive December 27th 2023)
CCP4 : 9.0.004 (Gargrove)
Density-Fitness : 1.0.11
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.41.2

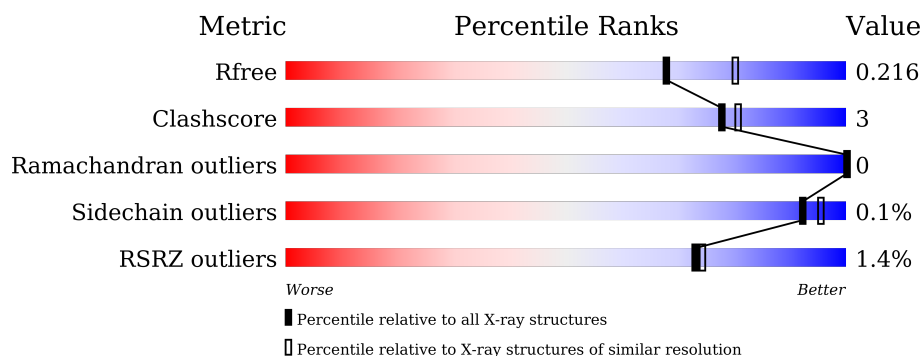
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.09 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}	164625	6234 (2.10-2.10)
Clashscore	180529	6893 (2.10-2.10)
Ramachandran outliers	177936	6839 (2.10-2.10)
Sidechain outliers	177891	6840 (2.10-2.10)
RSRZ outliers	164620	6234 (2.10-2.10)

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	342	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 1%, orange 1%, yellow 1%, green 87%, grey 7%, grey 6%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> % 87% 7% 6% </div> </div>
1	C	342	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 1%, orange 1%, yellow 1%, green 89%, grey 5%, grey 6%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> % 89% 5% 6% </div> </div>
2	B	256	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 1%, orange 1%, yellow 1%, green 96%, grey 3%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> % 96% . . </div> </div>
2	D	256	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 3%, orange 1%, yellow 1%, green 90%, grey 9%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> 3% 90% 9% . </div> </div>

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 9164 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 NADH-dependent dihydrogenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	321	Total	C	N	O	S	0	0	0
			2344	1472	425	443	4			
1	C	322	Total	C	N	O	S	0	0	0
			2355	1478	429	444	4			

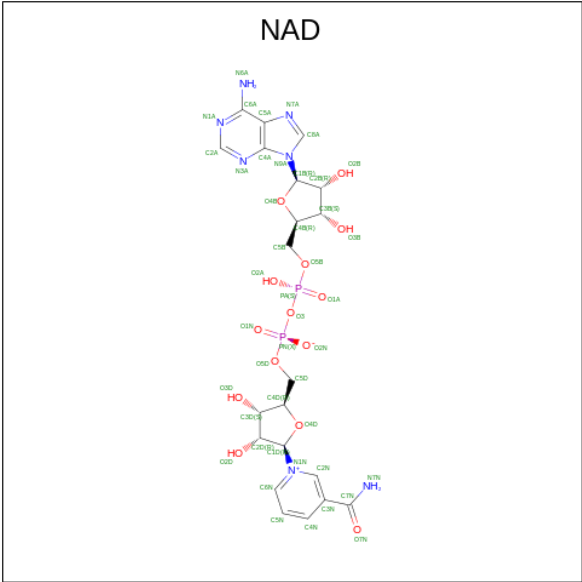
- Molecule 2 is a protein called Rift.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	253	Total	C	N	O	S	0	0	0
			1819	1140	314	360	5			
2	D	253	Total	C	N	O	S	0	0	0
			1825	1143	317	360	5			

There are 4 discrepancies between the modelled and reference sequences:

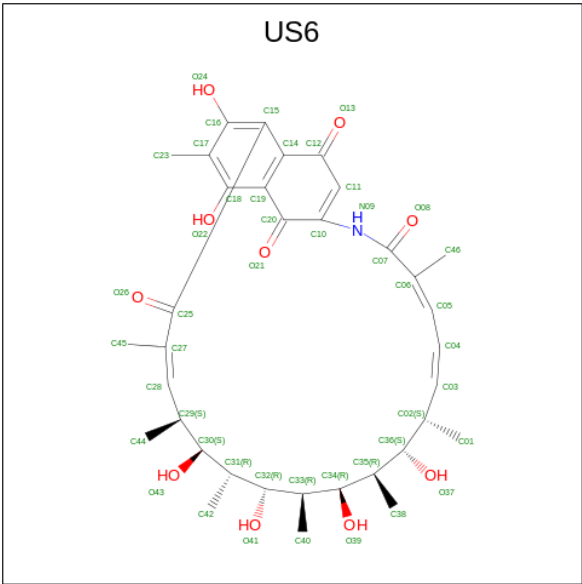
Chain	Residue	Modelled	Actual	Comment	Reference
B	348	MET	-	initiating methionine	UNP O52542
B	349	VAL	-	expression tag	UNP O52542
D	348	MET	-	initiating methionine	UNP O52542
D	349	VAL	-	expression tag	UNP O52542

- Molecule 3 is NICOTINAMIDE-ADENINE-DINUCLEOTIDE (three-letter code: NAD) (formula: C₂₁H₂₇N₇O₁₄P₂) (labeled as "Ligand of Interest" by depositor).



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

- Molecule 4 is (7E,9S,10S,11R,12R,13R,14R,15R,16S,17S,18E,20Z)-2,4,10,12,14,16-hexahydroxy-3,7,9,11,13,15,17,21-octamethyl-23-azatricyclo[22.3.1.0^{5,27}]octacos-1,3,5(27),7,18,20,24-heptaene-6,22,26,28-tetrone (three-letter code: US6) (formula: C₃₅H₄₅NO₁₀) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	N	O	0	0
			46	35	1	10		

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	C	1	Total	C	N	O	0	0
			46	35	1	10		

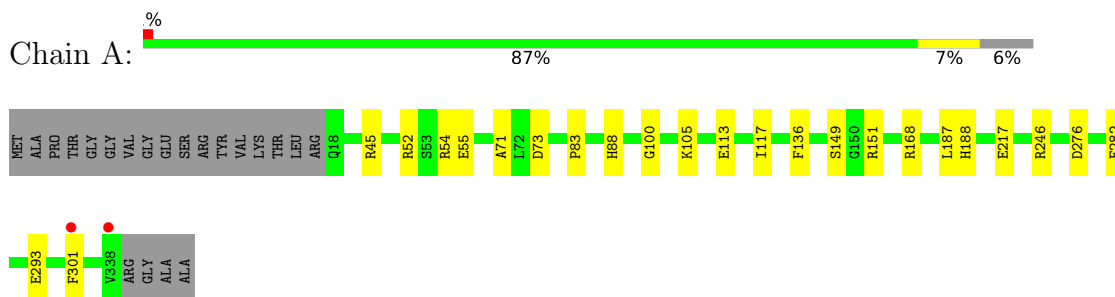
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	181	Total	O	0	0
			181	181		
5	B	171	Total	O	0	0
			171	171		
5	C	198	Total	O	0	0
			198	198		
5	D	135	Total	O	0	0
			135	135		

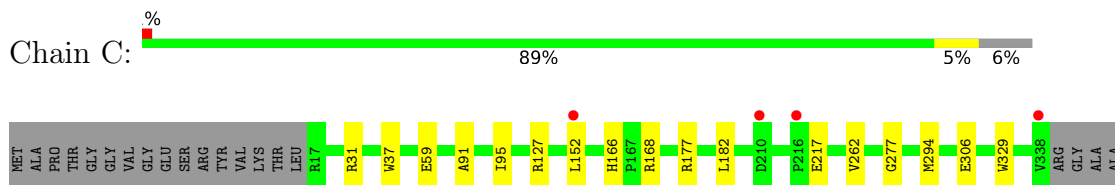
3 Residue-property plots

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.

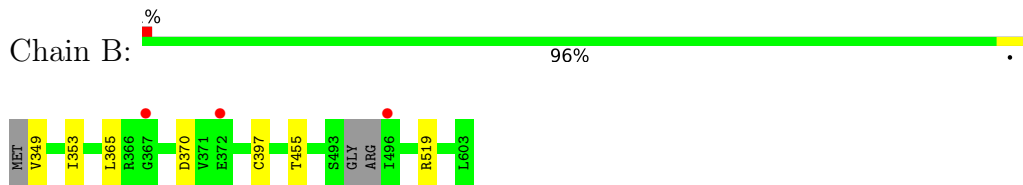
- Molecule 1: NADH-dependent dihydrogenase



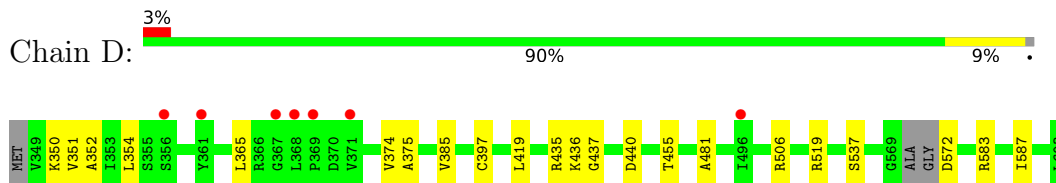
- Molecule 1: NADH-dependent dihydrogenase



- Molecule 2: RiFT



- Molecule 2: RiFT



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	191.54Å 82.11Å 88.56Å 90.00° 99.64° 90.00°	Depositor
Resolution (Å)	44.78 – 2.09 44.78 – 2.09	Depositor EDS
% Data completeness (in resolution range)	91.7 (44.78-2.09) 91.4 (44.78-2.09)	Depositor EDS
R_{merge}	0.03	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.33 (at 2.08Å)	Xtriage
Refinement program	PHENIX (1.18.2_3874: ???)	Depositor
R, R_{free}	0.176 , 0.216 0.178 , 0.216	Depositor DCC
R_{free} test set	78268 reflections (2.58%)	wwPDB-VP
Wilson B-factor (Å ²)	31.2	Xtriage
Anisotropy	0.073	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 37.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	9164	wwPDB-VP
Average B, all atoms (Å ²)	36.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.97% 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: US6, 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	0.42	0/2400	0.58	0/3276
1	C	0.52	0/2411	0.62	0/3290
2	B	0.58	1/1850 (0.1%)	0.66	0/2519
2	D	0.42	0/1856	0.60	0/2526
All	All	0.49	1/8517 (0.0%)	0.61	0/11611

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	397	CYS	CB-SG	-6.42	1.71	1.82

There are no bond angle outliers.

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	2344	0	2269	18	0
1	C	2355	0	2282	10	0
2	B	1819	0	1790	3	0
2	D	1825	0	1798	16	0
3	A	44	0	25	5	0
4	A	46	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	C	46	0	0	0	0
5	A	181	0	0	0	0
5	B	171	0	0	0	0
5	C	198	0	0	0	0
5	D	135	0	0	2	0
All	All	9164	0	8164	45	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 3.

All (45) 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)
3:A:501:NAD:O4B	3:A:501:NAD:C1B	1.64	1.15
1:A:83:PRO:HD3	3:A:501:NAD:H52A	1.70	0.74
2:B:353:ILE:HD11	2:B:365:LEU:HD12	1.76	0.65
2:D:350:LYS:HD3	2:D:374:VAL:HG11	1.79	0.63
2:D:397:CYS:SG	2:D:419:LEU:HD21	2.39	0.63
2:D:506:ARG:HH11	2:D:506:ARG:HG2	1.67	0.59
1:C:152:LEU:HD21	1:C:262:VAL:HG23	1.87	0.57
2:B:349:VAL:N	2:B:370:ASP:O	2.38	0.57
1:A:188:HIS:CE1	3:A:501:NAD:H72N	2.23	0.55
1:C:37:TRP:CE2	1:C:294:MET:HG3	2.42	0.55
1:A:45:ARG:NH1	1:A:73:ASP:O	2.38	0.54
1:C:91:ALA:O	1:C:95:ILE:HG12	2.09	0.52
1:A:105:LYS:HD2	1:A:187:LEU:HD22	1.92	0.50
1:C:127:ARG:NH1	1:C:306:GLU:OE1	2.44	0.50
2:D:354:LEU:HG	2:D:385:VAL:HG11	1.93	0.50
2:D:506:ARG:HG2	2:D:506:ARG:NH1	2.28	0.49
2:D:572:ASP:N	2:D:572:ASP:OD1	2.47	0.48
1:A:282:PHE:CE2	2:D:437:GLY:HA3	2.50	0.47
1:A:246:ARG:NH2	1:C:277:GLY:O	2.48	0.46
1:A:105:LYS:HE2	3:A:501:NAD:N7N	2.31	0.46
1:C:166:HIS:ND1	1:C:168:ARG:O	2.48	0.46
2:D:583:ARG:O	2:D:587:ILE:HG12	2.16	0.46
1:A:52:ARG:HA	1:A:55:GLU:OE2	2.15	0.45
1:A:54:ARG:NH1	1:A:55:GLU:OE1	2.49	0.45
1:A:149:SER:OG	1:A:151:ARG:HG2	2.17	0.45
1:A:276:ASP:OD2	2:D:537:SER:OG	2.28	0.44
1:A:88:HIS:HA	1:A:117:ILE:HD11	1.99	0.44
1:A:88:HIS:CD2	1:A:113:GLU:HG2	2.53	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:100:GLY:HA3	1:A:301:PHE:HE1	1.82	0.43
1:A:45:ARG:HD2	1:A:71:ALA:O	2.18	0.43
1:C:177:ARG:HD3	1:C:329:TRP:CG	2.54	0.43
2:D:436:LYS:NZ	2:D:440:ASP:OD2	2.52	0.43
1:C:182:LEU:HD23	1:C:182:LEU:HA	1.88	0.42
1:C:31:ARG:NH2	1:C:59:GLU:OE2	2.49	0.42
1:A:105:LYS:HE2	3:A:501:NAD:C7N	2.49	0.42
2:B:455:THR:O	2:B:519:ARG:HA	2.18	0.41
2:D:455:THR:O	2:D:519:ARG:HA	2.20	0.41
2:D:351:VAL:HG11	2:D:365:LEU:HD13	2.02	0.41
2:D:572:ASP:N	5:D:707:HOH:O	2.53	0.41
1:C:177:ARG:HD3	1:C:329:TRP:CD1	2.55	0.41
2:D:481:ALA:HB2	2:D:587:ILE:HD11	2.02	0.41
1:A:168:ARG:C	1:A:217:GLU:HG3	2.41	0.41
2:D:352:ALA:HA	2:D:375:ALA:O	2.21	0.41
1:A:136:PHE:CD1	1:A:293:GLU:HG3	2.56	0.40
2:D:435:ARG:HD3	5:D:814:HOH:O	2.21	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	319/342 (93%)	316 (99%)	3 (1%)	0	100	100
1	C	320/342 (94%)	316 (99%)	4 (1%)	0	100	100
2	B	249/256 (97%)	246 (99%)	3 (1%)	0	100	100
2	D	249/256 (97%)	245 (98%)	4 (2%)	0	100	100
All	All	1137/1196 (95%)	1123 (99%)	14 (1%)	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	222/236 (94%)	222 (100%)	0	100	100
1	C	223/236 (94%)	222 (100%)	1 (0%)	89	93
2	B	179/181 (99%)	179 (100%)	0	100	100
2	D	180/181 (99%)	180 (100%)	0	100	100
All	All	804/834 (96%)	803 (100%)	1 (0%)	92	95

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

Mol	Chain	Res	Type
1	C	217	GLU

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

Mol	Chain	Res	Type
2	B	387	GLN

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 oligosaccharides in this entry.

5.6 Ligand geometry

3 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	$\# Z > 2$
4	US6	C	401	-	48,48,48	2.31	12 (25%)	69,72,72	1.67	10 (14%)
3	NAD	A	501	-	42,48,48	4.47	13 (30%)	50,73,73	1.18	4 (8%)
4	US6	A	502	-	48,48,48	2.24	14 (29%)	69,72,72	1.62	13 (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
4	US6	C	401	-	-	31/58/74/74	0/2/3/3
3	NAD	A	501	-	-	6/26/62/62	0/5/5/5
4	US6	A	502	-	-	20/58/74/74	0/2/3/3

All (39) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	501	NAD	O4B-C1B	16.57	1.64	1.41
3	A	501	NAD	C2B-C1B	-12.47	1.34	1.53
3	A	501	NAD	O4D-C1D	9.52	1.54	1.41
3	A	501	NAD	C7N-N7N	8.55	1.49	1.33
3	A	501	NAD	C2D-C3D	-8.39	1.30	1.53
4	C	401	US6	C11-C10	6.74	1.50	1.36
4	A	502	US6	C11-C10	6.48	1.49	1.36
3	A	501	NAD	O4B-C4B	-6.09	1.31	1.45
3	A	501	NAD	O4D-C4D	-6.01	1.31	1.45
4	C	401	US6	C07-N09	5.22	1.47	1.37
3	A	501	NAD	C3N-C7N	4.90	1.57	1.50
4	C	401	US6	C14-C12	4.66	1.58	1.46
4	A	502	US6	C07-N09	4.55	1.45	1.37

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	502	US6	C04-C05	4.31	1.56	1.43
4	A	502	US6	C19-C20	4.28	1.57	1.46
4	C	401	US6	C04-C05	4.16	1.56	1.43
4	C	401	US6	C19-C20	4.16	1.57	1.46
4	A	502	US6	C28-C27	4.11	1.38	1.33
4	C	401	US6	C28-C27	4.07	1.38	1.33
4	A	502	US6	C27-C25	4.01	1.58	1.48
4	C	401	US6	C27-C25	3.92	1.57	1.48
4	A	502	US6	C14-C12	3.88	1.56	1.46
3	A	501	NAD	C3D-C4D	3.76	1.62	1.53
4	C	401	US6	C11-C12	3.59	1.52	1.44
4	A	502	US6	C15-C25	3.50	1.58	1.51
4	C	401	US6	C15-C25	3.26	1.58	1.51
4	A	502	US6	C11-C12	3.24	1.51	1.44
3	A	501	NAD	O7N-C7N	-2.95	1.18	1.24
3	A	501	NAD	O3D-C3D	-2.82	1.36	1.43
4	C	401	US6	O13-C12	-2.60	1.18	1.24
4	C	401	US6	C10-N09	2.59	1.47	1.41
4	A	502	US6	O13-C12	-2.56	1.18	1.24
3	A	501	NAD	C6A-N6A	2.50	1.43	1.34
4	A	502	US6	C10-N09	2.39	1.47	1.41
3	A	501	NAD	C3B-C4B	2.35	1.59	1.53
4	A	502	US6	O08-C07	-2.28	1.19	1.23
4	A	502	US6	O41-C32	-2.09	1.38	1.43
4	A	502	US6	O26-C25	-2.06	1.18	1.23
4	C	401	US6	O24-C16	2.05	1.41	1.37

All (27) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	502	US6	C20-C10-N09	6.46	117.91	112.35
4	C	401	US6	C20-C10-N09	6.44	117.89	112.35
4	C	401	US6	C35-C34-C33	-4.23	108.34	115.43
3	A	501	NAD	N3A-C2A-N1A	-4.20	122.12	128.68
4	C	401	US6	C33-C32-C31	-3.69	109.24	115.43
4	A	502	US6	C01-C02-C36	-3.58	103.81	111.31
4	A	502	US6	C33-C32-C31	-3.53	109.51	115.43
4	A	502	US6	C11-C10-N09	-3.27	120.86	126.57
4	C	401	US6	C44-C29-C30	-3.13	104.74	111.31
4	C	401	US6	C02-C03-C04	-3.09	119.65	126.16
4	C	401	US6	C11-C10-N09	-3.07	121.20	126.57
3	A	501	NAD	PN-O3-PA	-2.97	122.65	132.83

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	502	US6	C45-C27-C25	2.96	120.96	115.53
4	C	401	US6	O13-C12-C11	-2.94	116.87	121.79
4	C	401	US6	C02-C36-C35	-2.92	109.00	114.96
4	A	502	US6	C44-C29-C30	-2.88	105.27	111.31
4	A	502	US6	C35-C34-C33	-2.82	110.71	115.43
3	A	501	NAD	C4A-C5A-N7A	-2.81	106.47	109.40
4	A	502	US6	O13-C12-C11	-2.63	117.39	121.79
4	C	401	US6	C45-C27-C25	2.49	120.10	115.53
4	A	502	US6	C45-C27-C28	-2.48	120.27	124.30
4	C	401	US6	C10-N09-C07	-2.39	115.50	125.01
4	A	502	US6	C36-C35-C34	-2.31	107.90	112.54
4	A	502	US6	C02-C03-C04	-2.20	121.53	126.16
4	A	502	US6	C05-C04-C03	-2.20	119.19	124.53
4	A	502	US6	C29-C30-C31	-2.16	110.55	114.96
3	A	501	NAD	O4D-C1D-C2D	-2.13	103.81	106.93

There are no chirality outliers.

All (57) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	501	NAD	C5B-O5B-PA-O1A
4	A	502	US6	C28-C29-C30-C31
4	A	502	US6	C44-C29-C30-C31
4	A	502	US6	C44-C29-C30-O43
4	C	401	US6	C11-C10-N09-C07
4	C	401	US6	C20-C10-N09-C07
4	C	401	US6	C28-C29-C30-C31
4	C	401	US6	C28-C29-C30-O43
4	C	401	US6	C44-C29-C30-C31
4	C	401	US6	O43-C30-C31-C32
4	C	401	US6	O43-C30-C31-C42
4	C	401	US6	C32-C33-C34-C35
4	C	401	US6	C32-C33-C34-O39
4	C	401	US6	C40-C33-C34-C35
4	C	401	US6	C40-C33-C34-O39
4	C	401	US6	C34-C35-C36-C02
4	C	401	US6	C34-C35-C36-O37
4	C	401	US6	C38-C35-C36-C02
4	C	401	US6	C38-C35-C36-O37
4	A	502	US6	C38-C35-C36-O37
4	C	401	US6	C29-C30-C31-C42
4	C	401	US6	C29-C30-C31-C32

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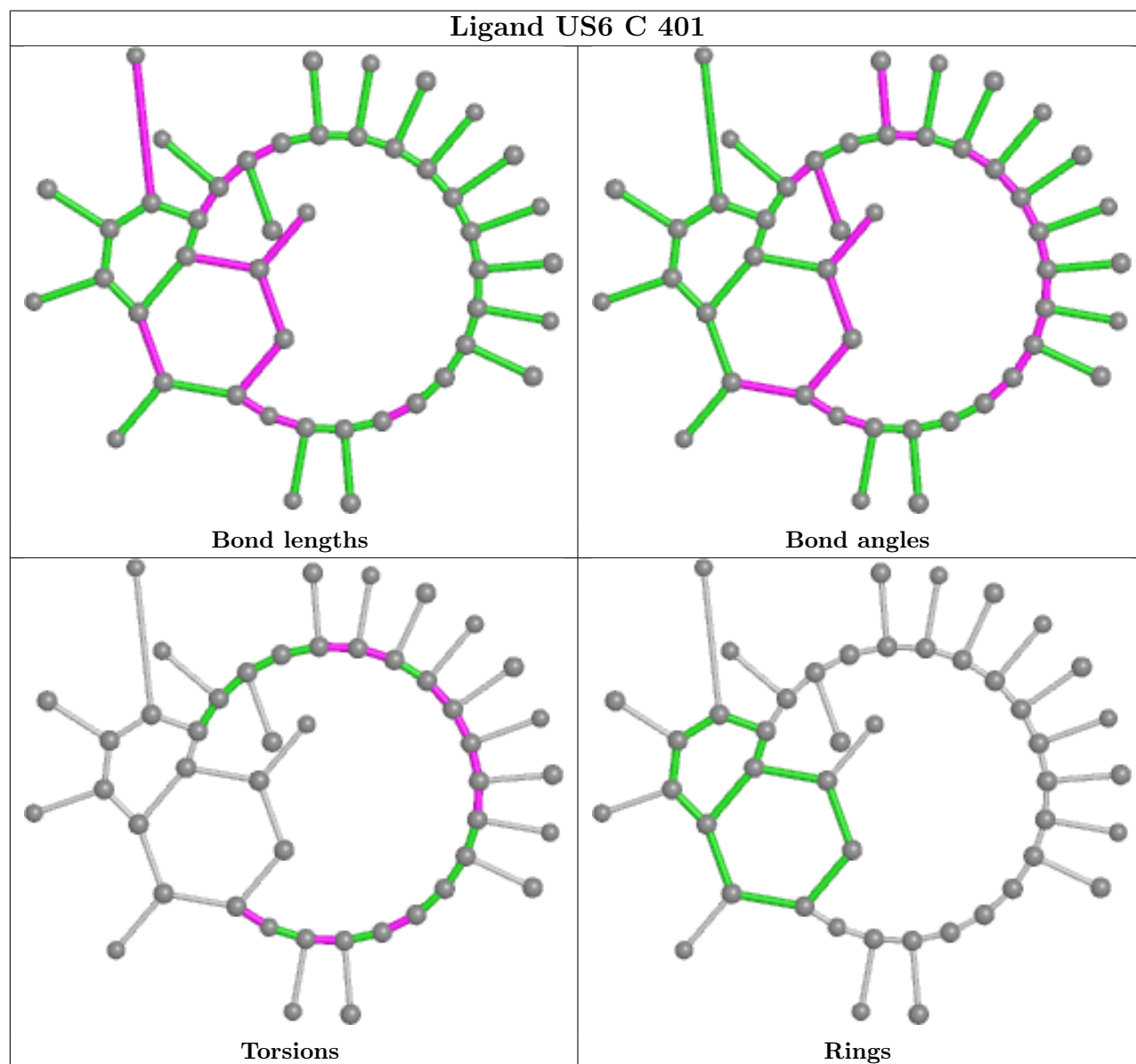
Mol	Chain	Res	Type	Atoms
3	A	501	NAD	O4B-C4B-C5B-O5B
3	A	501	NAD	C3B-C4B-C5B-O5B
4	A	502	US6	C34-C35-C36-O37
4	A	502	US6	C28-C29-C30-O43
4	C	401	US6	C44-C29-C30-O43
4	A	502	US6	C38-C35-C36-C02
4	A	502	US6	C34-C35-C36-C02
4	C	401	US6	O41-C32-C33-C40
4	C	401	US6	O39-C34-C35-C36
4	A	502	US6	C11-C10-N09-C07
4	C	401	US6	O41-C32-C33-C34
4	C	401	US6	O39-C34-C35-C38
4	A	502	US6	C05-C06-C07-N09
4	C	401	US6	C05-C06-C07-N09
4	A	502	US6	C20-C10-N09-C07
4	A	502	US6	C01-C02-C03-C04
4	A	502	US6	C42-C31-C32-C33
4	C	401	US6	C31-C32-C33-C40
4	C	401	US6	C33-C34-C35-C38
4	C	401	US6	C31-C32-C33-C34
3	A	501	NAD	C5B-O5B-PA-O3
4	C	401	US6	C05-C06-C07-O08
3	A	501	NAD	C5B-O5B-PA-O2A
4	A	502	US6	C30-C31-C32-C33
4	C	401	US6	C33-C34-C35-C36
4	A	502	US6	C30-C31-C32-O41
4	A	502	US6	C46-C06-C07-N09
4	C	401	US6	C46-C06-C07-N09
4	A	502	US6	C05-C06-C07-O08
4	A	502	US6	C03-C04-C05-C06
4	C	401	US6	C03-C04-C05-C06
4	A	502	US6	C16-C15-C25-O26
3	A	501	NAD	PA-O3-PN-O1N
4	A	502	US6	C46-C06-C07-O08
4	C	401	US6	C46-C06-C07-O08

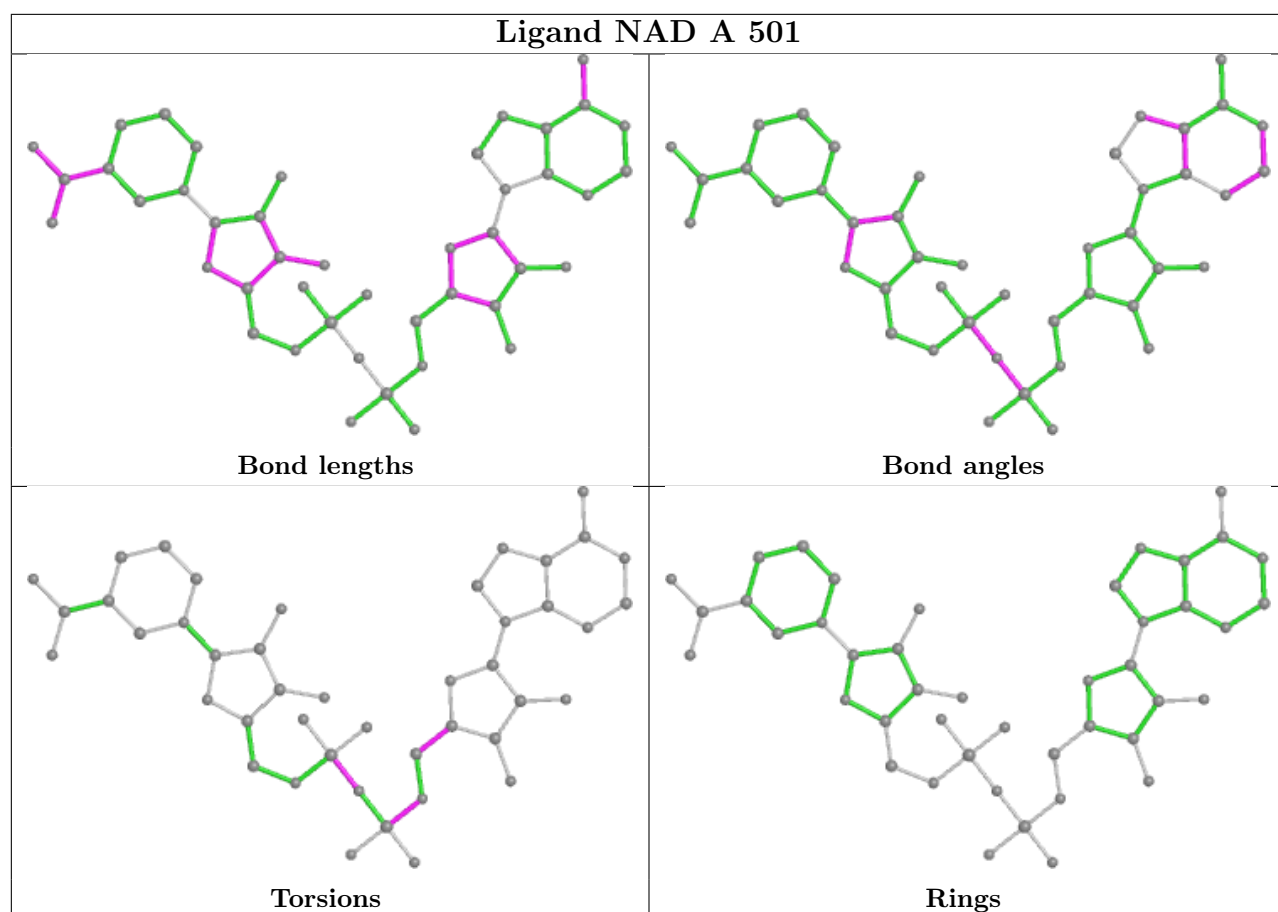
There are no ring outliers.

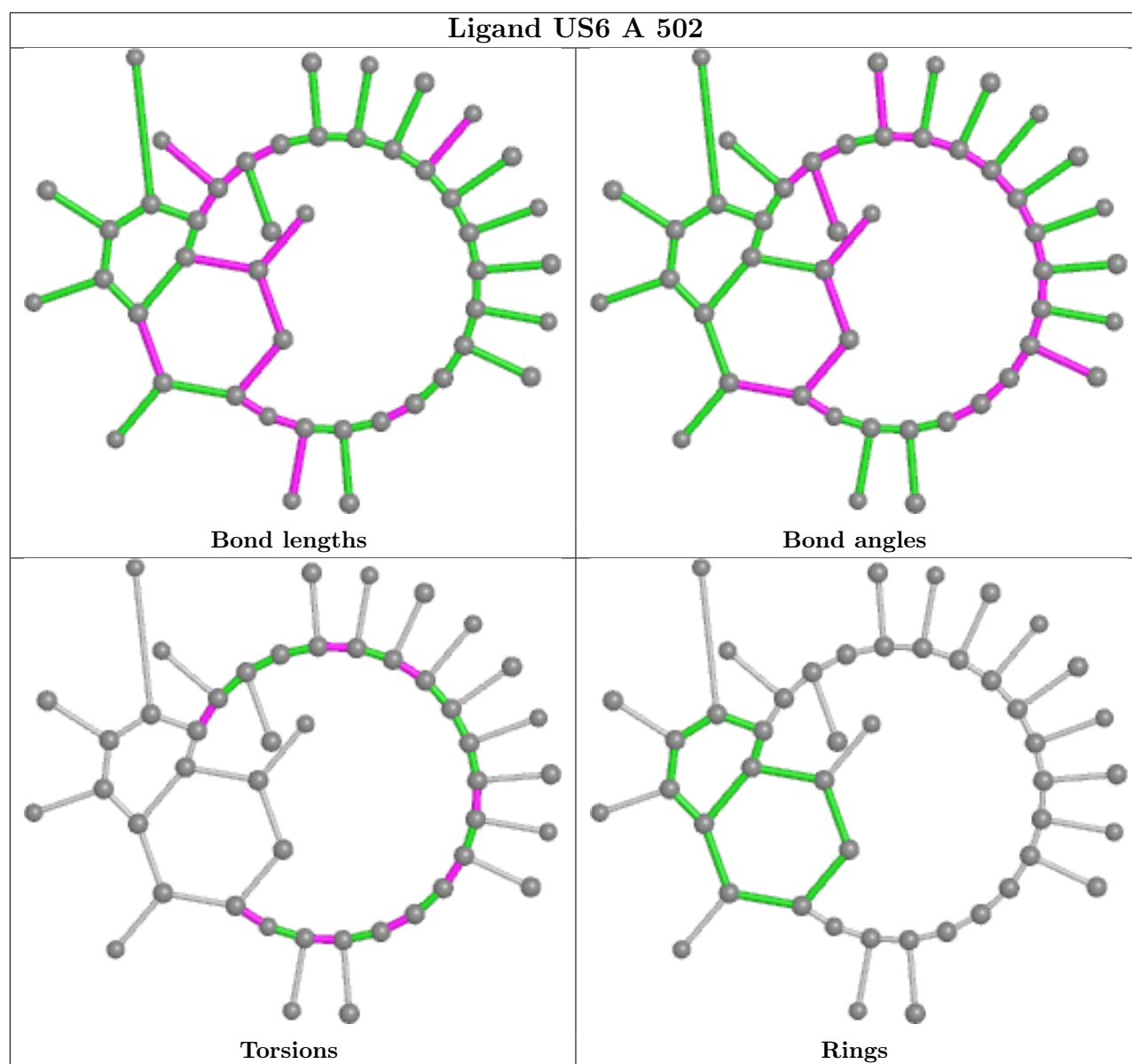
1 monomer is involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	501	NAD	5	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	321/342 (93%)	-0.12	2 (0%) 85 86	23, 36, 53, 74	0
1	C	322/342 (94%)	-0.20	4 (1%) 76 77	23, 34, 47, 66	0
2	B	253/256 (98%)	-0.25	3 (1%) 76 77	23, 31, 48, 81	0
2	D	253/256 (98%)	0.01	7 (2%) 55 57	26, 36, 67, 90	0
All	All	1149/1196 (96%)	-0.14	16 (1%) 73 74	23, 34, 54, 90	0

All (16) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	496	ILE	3.1
2	D	356	SER	2.9
2	D	368	LEU	2.8
2	D	371	VAL	2.6
2	B	367	GLY	2.6
1	C	216	PRO	2.3
1	A	338	VAL	2.3
2	D	361	TYR	2.2
1	C	338	VAL	2.2
2	D	496	ILE	2.1
1	C	152	LEU	2.1
2	D	367	GLY	2.1
2	B	372	GLU	2.0
1	A	301	PHE	2.0
2	D	369	PRO	2.0
1	C	210	ASP	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](#)

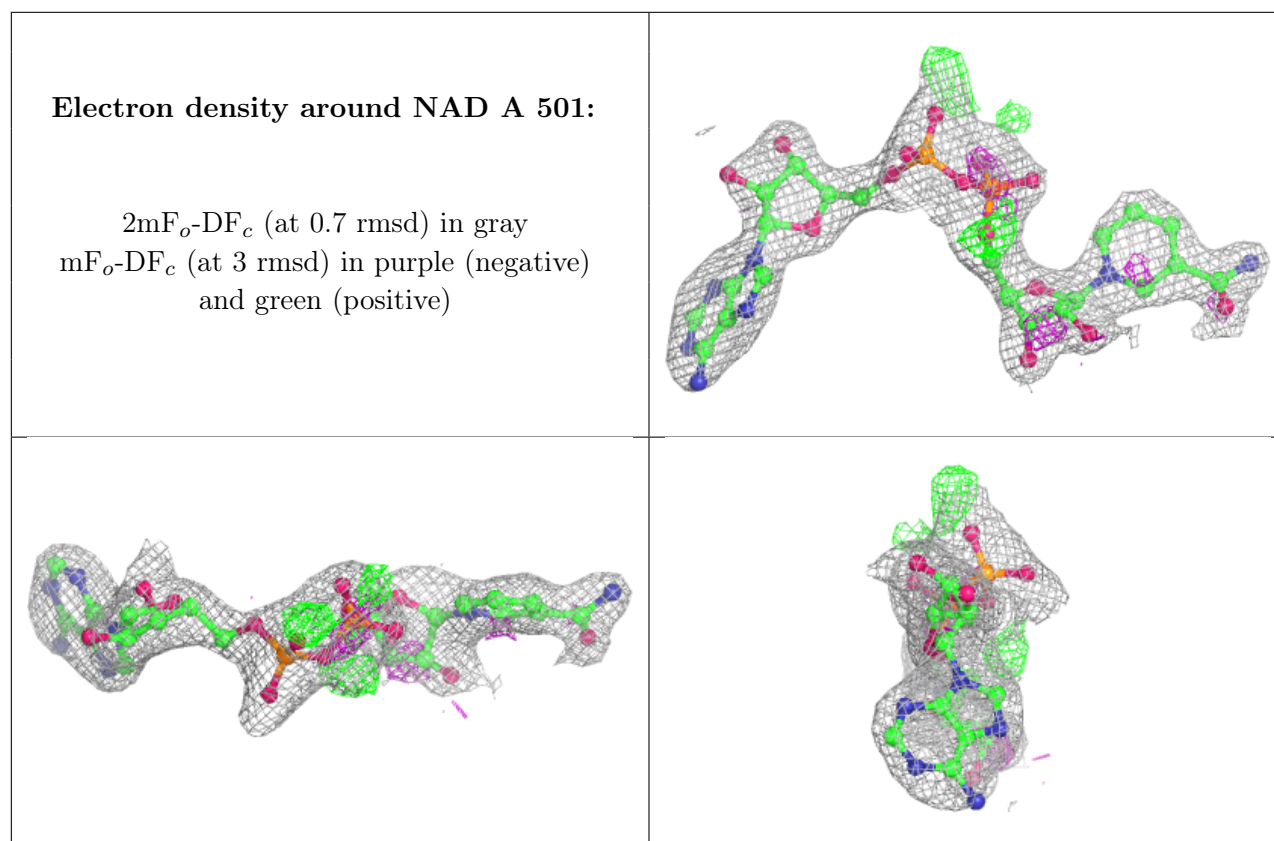
There are no monosaccharides in this entry.

6.4 Ligands [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.

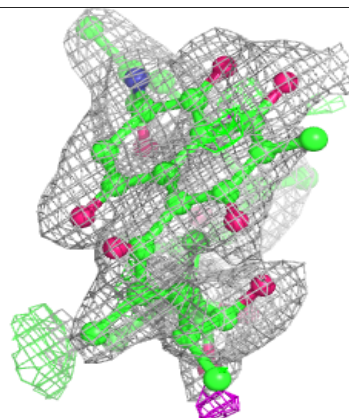
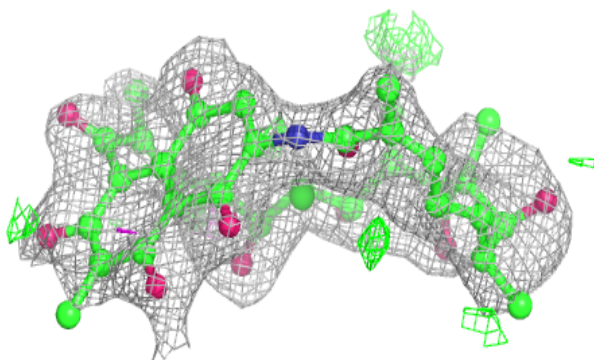
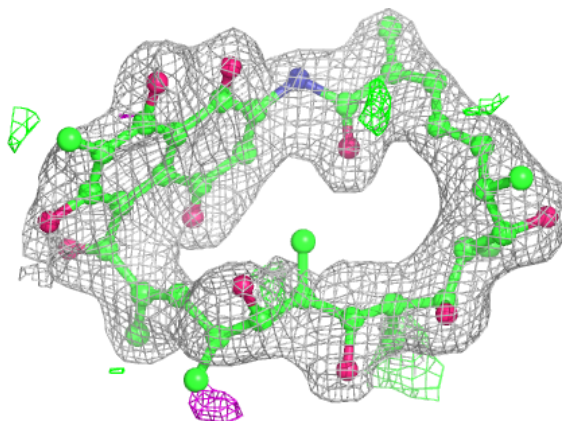
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
3	NAD	A	501	44/44	0.85	0.11	44,52,65,74	0
4	US6	C	401	46/46	0.85	0.14	32,52,62,63	0
4	US6	A	502	46/46	0.89	0.12	29,41,50,56	0

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.



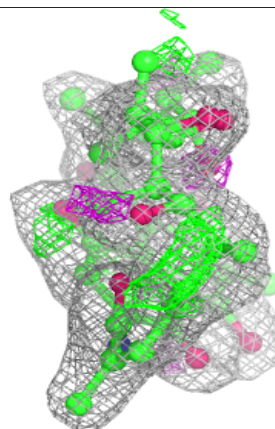
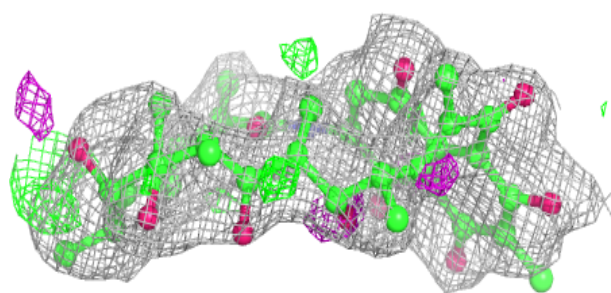
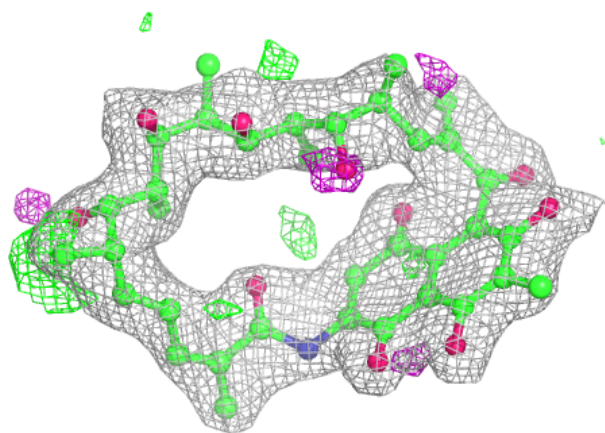
Electron density around US6 C 401:

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



Electron density around US6 A 502:

$2mF_o-DF_c$ (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.