Full wwPDB X-ray Structure Validation Report (i)

Jun 12, 2024 – 07:46 PM EDT

PDB ID	:	3VPS
Title	:	Structure of a novel NAD dependent-NDP-hexosamine 5,6-dehydratase, TunA, involved in tunicamycin biosynthesis
Authors	:	Wyszynski, F.J.; Lee, S.S.; Yabe, T.; Wang, H.; Gomez-Escribano, J.P.; Bibb, M.J.; Lee, S.J.; Davies, G.J.; Davis, B.G.
Deposited on	:	2012-03-12
Resolution	:	1.90 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 (i) 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 (i)) were used in the production of this report:

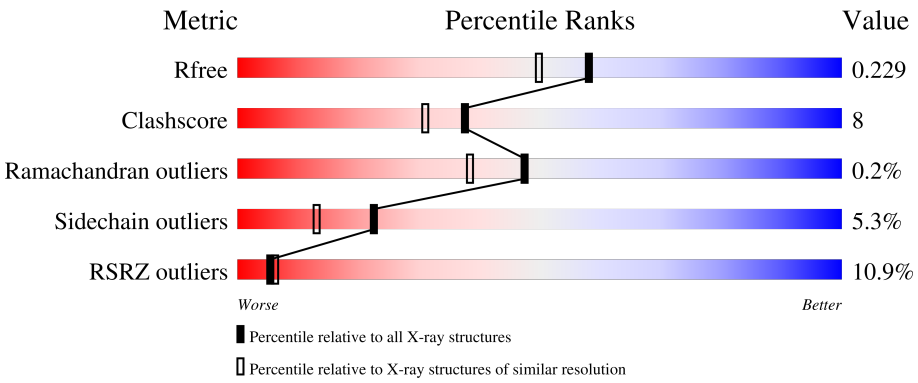
MolProbity	:	4.02b-467
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	1.20.1
EDS	:	2.36.2
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36.2

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 1.90 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	6207 (1.90-1.90)
Clashscore	141614	6847 (1.90-1.90)
Ramachandran outliers	138981	6760 (1.90-1.90)
Sidechain outliers	138945	6760 (1.90-1.90)
RSRZ outliers	127900	6082 (1.90-1.90)

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 <=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	321	10% 78% 14% • 6%
1	B	321	10% 77% 14% • 6%

2 Entry composition (i)

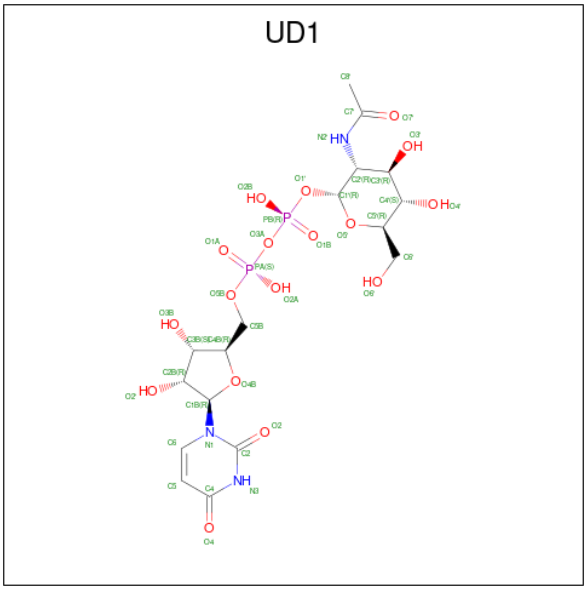
There are 4 unique types of molecules in this entry. The entry contains 5003 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 NAD-dependent epimerase/dehydratase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	303	2315	1453	413	444	5	0	0	0
			Total	C	N	O	S			
1	B	301	2303	1447	410	441	5	0	0	0

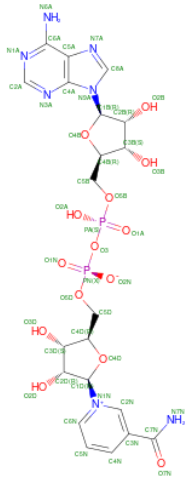
- Molecule 2 is URIDINE-DIPHOSPHATE-N-ACETYLGLUCOSAMINE (three-letter code: UD1) (formula: $C_{17}H_{27}N_3O_{17}P_2$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	N	O	P		
2	A	1	39	17	3	17	2	0	0
			Total	C	N	O	P		
2	B	1	39	17	3	17	2	0	0

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

NAD



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

- Molecule 4 is water.

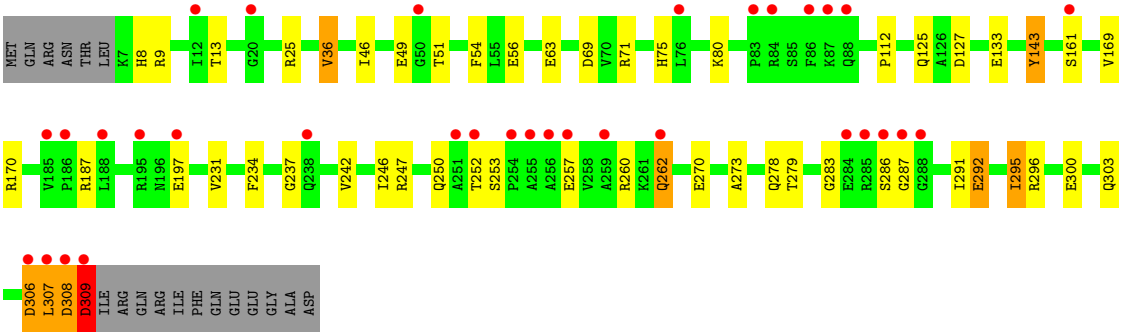
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
4	A	125	125	125	0	0
			Total	O		
4	B	94	94	94	0	0

3 Residue-property plots (i)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density (RSRZ > 2). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

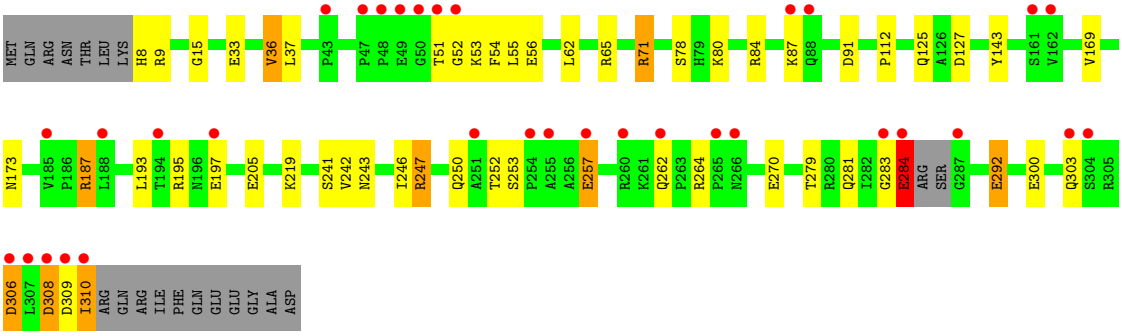
• Molecule 1: NAD-dependent epimerase/dehydratase

Chain A:



• Molecule 1: NAD-dependent epimerase/dehydratase

Chain B:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α, β, γ	45.45Å 51.06Å 67.80Å 98.13° 106.65° 94.36°	Depositor
Resolution (Å)	22.26 – 1.90 22.26 – 1.90	Depositor EDS
% Data completeness (in resolution range)	99.4 (22.26-1.90) 99.4 (22.26-1.90)	Depositor EDS
R_{merge}	0.19	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.99 (at 1.90Å)	Xtriage
Refinement program	CNS, REFMAC 5.5.0109	Depositor
R, R_{free}	0.186 , 0.235 0.183 , 0.229	Depositor DCC
R_{free} test set	2283 reflections (5.08%)	wwPDB-VP
Wilson B-factor (Å^2)	27.5	Xtriage
Anisotropy	0.084	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.42 , 65.5	EDS
L-test for twinning[2]	$< \|L\| > = 0.51$, $< L^2 > = 0.35$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.95	EDS
Total number of atoms	5003	wwPDB-VP
Average B, all atoms (Å^2)	41.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.92% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ 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: NAD, 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	1.19	10/2361 (0.4%)	0.98	8/3215 (0.2%)
1	B	1.05	7/2348 (0.3%)	1.02	13/3197 (0.4%)
All	All	1.12	17/4709 (0.4%)	1.00	21/6412 (0.3%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand.A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	B	0	1

All (17) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	284	GLU	CG-CD	-9.57	1.37	1.51
1	A	262	GLN	CG-CD	-8.29	1.31	1.51
1	A	309	ASP	CB-CG	-7.26	1.36	1.51
1	A	36	VAL	CB-CG2	-7.15	1.37	1.52
1	A	262	GLN	CB-CG	-7.15	1.33	1.52
1	A	63	GLU	CD-OE1	-6.99	1.18	1.25
1	B	36	VAL	CB-CG1	-6.55	1.39	1.52
1	A	63	GLU	CB-CG	-6.24	1.40	1.52
1	A	49	GLU	CB-CG	-6.18	1.40	1.52
1	B	71	ARG	NE-CZ	-6.17	1.25	1.33
1	B	257	GLU	CG-CD	-6.12	1.42	1.51
1	B	187	ARG	CZ-NH1	-5.96	1.25	1.33
1	A	308	ASP	CB-CG	-5.62	1.40	1.51
1	A	187	ARG	CZ-NH1	-5.37	1.26	1.33
1	B	71	ARG	CB-CG	-5.32	1.38	1.52
1	A	143	TYR	CD1-CE1	5.18	1.47	1.39

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	292	GLU	CG-CD	5.17	1.59	1.51

All (21) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	187	ARG	NE-CZ-NH2	14.22	127.41	120.30
1	B	187	ARG	NE-CZ-NH1	-12.52	114.04	120.30
1	A	187	ARG	NE-CZ-NH2	11.69	126.15	120.30
1	A	187	ARG	NE-CZ-NH1	-10.08	115.26	120.30
1	A	9	ARG	NE-CZ-NH2	-7.95	116.33	120.30
1	A	36	VAL	CG1-CB-CG2	-7.37	99.11	110.90
1	A	262	GLN	N-CA-CB	-7.13	97.76	110.60
1	B	264	ARG	NE-CZ-NH1	-6.82	116.89	120.30
1	B	71	ARG	CG-CD-NE	-6.72	97.68	111.80
1	B	169	VAL	CG1-CB-CG2	-6.46	100.56	110.90
1	B	71	ARG	NE-CZ-NH1	-6.13	117.23	120.30
1	B	71	ARG	CD-NE-CZ	-6.13	115.02	123.60
1	A	169	VAL	CG1-CB-CG2	-5.72	101.75	110.90
1	B	247	ARG	NE-CZ-NH1	5.61	123.10	120.30
1	B	247	ARG	CG-CD-NE	5.57	123.49	111.80
1	A	262	GLN	CB-CG-CD	-5.56	97.13	111.60
1	B	284	GLU	OE1-CD-OE2	5.54	129.95	123.30
1	A	127	ASP	CB-CG-OD1	5.37	123.14	118.30
1	B	127	ASP	CB-CG-OD1	5.29	123.06	118.30
1	B	91	ASP	CB-CG-OD1	5.27	123.04	118.30
1	B	247	ARG	NE-CZ-NH2	-5.01	117.80	120.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	78	SER	Peptide

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	2315	0	2318	37	0
1	B	2303	0	2308	32	0
2	A	39	0	25	2	0
2	B	39	0	25	2	0
3	A	44	0	26	2	0
3	B	44	0	26	2	0
4	A	125	0	0	7	2
4	B	94	0	0	4	1
All	All	5003	0	4728	73	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 8.

All (73) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:308:ASP:OD1	1:A:309:ASP:N	1.83	1.10
1:A:51:THR:HG21	4:A:610:HOH:O	1.57	1.02
1:A:309:ASP:O	1:A:309:ASP:OD1	1.77	1.00
1:B:262:GLN:OE1	1:B:262:GLN:N	2.00	0.94
1:B:309:ASP:O	1:B:310:ILE:HG13	1.70	0.91
1:B:125:GLN:HE22	1:B:270:GLU:H	1.23	0.81
1:B:309:ASP:O	1:B:310:ILE:CG1	2.30	0.79
1:A:125:GLN:HE22	1:A:270:GLU:H	1.27	0.78
1:B:279:THR:O	1:B:283:GLY:HA2	1.84	0.76
1:B:84:ARG:O	1:B:87:LYS:HG2	1.90	0.71
1:B:195:ARG:HB3	1:B:197:GLU:OE1	1.92	0.69
1:B:308:ASP:OD1	1:B:308:ASP:C	2.30	0.69
2:B:400:UD1:H4'	3:B:401:NAD:C5N	2.26	0.65
2:A:400:UD1:H4'	3:A:401:NAD:C5N	2.28	0.64
1:A:36:VAL:HG23	1:A:54:PHE:HD1	1.63	0.63
1:A:309:ASP:O	1:A:309:ASP:CG	2.38	0.62
1:A:306:ASP:OD1	1:A:306:ASP:C	2.40	0.60
1:B:71:ARG:HG3	1:B:71:ARG:NH1	2.15	0.59
1:A:279:THR:O	1:A:283:GLY:HA2	2.01	0.58
1:B:15:GLY:HA3	1:B:36:VAL:HG13	1.85	0.58
1:A:197:GLU:HA	1:A:257:GLU:O	2.04	0.58
2:B:400:UD1:H4'	3:B:401:NAD:C4N	2.34	0.57
1:B:9:ARG:NH2	4:B:530:HOH:O	2.24	0.56
1:A:197:GLU:HG2	1:A:257:GLU:OE2	2.06	0.56
1:B:306:ASP:C	1:B:306:ASP:OD1	2.45	0.56
1:A:300:GLU:HA	1:A:303:GLN:HE21	1.72	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:52:GLY:HA2	4:B:530:HOH:O	2.06	0.54
1:B:279:THR:O	1:B:283:GLY:CA	2.55	0.54
1:A:260:ARG:NH2	4:A:620:HOH:O	2.24	0.54
1:A:8:HIS:NE2	4:A:618:HOH:O	2.23	0.53
1:A:309:ASP:OD1	1:A:309:ASP:C	2.37	0.52
1:B:257:GLU:N	1:B:257:GLU:OE1	2.43	0.52
1:A:13:THR:OG1	1:A:75:HIS:HD2	1.93	0.50
1:A:242:VAL:O	1:A:246:ILE:HG12	2.11	0.50
2:A:400:UD1:H4'	3:A:401:NAD:C4N	2.41	0.50
1:B:308:ASP:OD1	1:B:308:ASP:O	2.30	0.49
1:B:37:LEU:HD13	1:B:62:LEU:HD13	1.95	0.48
1:A:133:GLU:HB3	1:A:231:VAL:HG13	1.96	0.48
1:A:25:ARG:NH2	4:A:533:HOH:O	2.47	0.48
1:B:300:GLU:HA	1:B:303:GLN:HE21	1.80	0.47
1:B:37:LEU:CD1	1:B:62:LEU:HD13	2.45	0.47
1:A:80:LYS:HA	1:A:143:TYR:CE1	2.50	0.46
1:B:283:GLY:C	1:B:284:GLU:HG2	2.25	0.46
1:B:71:ARG:HA	1:B:112:PRO:HD2	1.97	0.46
1:A:170:ARG:CZ	1:A:231:VAL:HG21	2.46	0.46
1:B:219:LYS:NZ	4:B:551:HOH:O	2.34	0.46
1:A:291:ILE:O	1:A:295:ILE:HD12	2.16	0.45
1:B:84:ARG:O	1:B:87:LYS:CG	2.63	0.45
1:A:306:ASP:O	1:A:309:ASP:HB3	2.17	0.45
1:B:71:ARG:NH1	1:B:71:ARG:CG	2.68	0.45
1:B:80:LYS:HA	1:B:143:TYR:CE1	2.52	0.45
1:A:75:HIS:HE1	4:A:513:HOH:O	1.98	0.45
1:A:291:ILE:HG13	1:A:295:ILE:HD12	1.99	0.44
1:A:291:ILE:HG13	1:A:295:ILE:CD1	2.47	0.44
1:A:71:ARG:HH11	1:A:71:ARG:HD3	1.60	0.44
1:B:242:VAL:O	1:B:246:ILE:HG12	2.16	0.44
1:A:279:THR:O	1:A:283:GLY:CA	2.64	0.43
1:B:54:PHE:CE2	1:B:56:GLU:HB2	2.53	0.43
1:B:173:ASN:ND2	4:B:506:HOH:O	2.48	0.43
1:A:54:PHE:CE2	1:A:56:GLU:HB2	2.53	0.43
1:A:237:GLY:HA2	1:A:273:ALA:O	2.18	0.43
1:B:8:HIS:CD2	1:B:33:GLU:OE1	2.72	0.43
1:B:246:ILE:O	1:B:250:GLN:HG3	2.19	0.43
1:A:71:ARG:HA	1:A:112:PRO:HD2	2.01	0.43
1:A:25:ARG:HD3	4:A:514:HOH:O	2.19	0.42
1:A:69:ASP:OD2	4:A:618:HOH:O	2.22	0.42
1:A:307:LEU:HD23	1:A:307:LEU:HA	1.89	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:296:ARG:NH1	1:A:300:GLU:OE1	2.53	0.42
1:A:308:ASP:CG	1:A:309:ASP:N	2.65	0.41
1:A:292:GLU:CD	1:A:292:GLU:H	2.23	0.41
1:A:234:PHE:HA	1:A:278:GLN:NE2	2.36	0.40
1:B:205:GLU:OE2	1:B:243:ASN:ND2	2.51	0.40
1:B:53:LYS:HE2	1:B:55:LEU:HD21	2.04	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)
4:A:597:HOH:O	4:A:617:HOH:O[1_655]	1.79	0.41
4:A:595:HOH:O	4:B:581:HOH:O[1_656]	1.98	0.22

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	301/321 (94%)	291 (97%)	9 (3%)	1 (0%)	41	31
1	B	297/321 (92%)	292 (98%)	5 (2%)	0	100	100
All	All	598/642 (93%)	583 (98%)	14 (2%)	1 (0%)	47	38

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	287	GLY

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	253/270 (94%)	240 (95%)	13 (5%)	24	14
1	B	252/270 (93%)	238 (94%)	14 (6%)	21	11
All	All	505/540 (94%)	478 (95%)	27 (5%)	22	13

All (27) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	46	ILE
1	A	161	SER
1	A	247	ARG
1	A	250	GLN
1	A	252	THR
1	A	253	SER
1	A	262	GLN
1	A	286	SER
1	A	292	GLU
1	A	295	ILE
1	A	306	ASP
1	A	307	LEU
1	A	309	ASP
1	B	51	THR
1	B	65	ARG
1	B	187	ARG
1	B	193	LEU
1	B	241	SER
1	B	247	ARG
1	B	252	THR
1	B	253	SER
1	B	281	GLN
1	B	284	GLU
1	B	292	GLU
1	B	306	ASP
1	B	308	ASP
1	B	310	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	75	HIS
1	A	88	GLN
1	A	125	GLN
1	A	158	GLN
1	A	173	ASN
1	A	278	GLN
1	A	303	GLN
1	B	22	HIS
1	B	125	GLN
1	B	173	ASN
1	B	278	GLN
1	B	281	GLN
1	B	303	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 monosaccharides in this entry.

5.6 Ligand geometry (i)

4 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	$\#\|Z\| > 2$
3	NAD	A	401	-	42,48,48	2.03	7 (16%)	50,73,73	2.28	7 (14%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\| > 2$	Counts	RMSZ	$\#\|Z\| > 2$
2	UD1	B	400	-	40,41,41	1.24	3 (7%)	59,62,62	1.62	9 (15%)
2	UD1	A	400	-	40,41,41	1.16	3 (7%)	59,62,62	1.74	12 (20%)
3	NAD	B	401	-	42,48,48	1.63	8 (19%)	50,73,73	2.73	15 (30%)

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
3	NAD	A	401	-	-	4/26/62/62	0/5/5/5
2	UD1	B	400	-	-	4/26/63/63	0/3/3/3
2	UD1	A	400	-	-	3/26/63/63	0/3/3/3
3	NAD	B	401	-	-	2/26/62/62	0/5/5/5

All (21) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	401	NAD	O7N-C7N	8.13	1.39	1.24
3	B	401	NAD	O7N-C7N	4.76	1.33	1.24
3	A	401	NAD	C7N-N7N	-4.29	1.25	1.33
3	A	401	NAD	C2A-N3A	3.97	1.38	1.32
2	B	400	UD1	PB-O3A	3.88	1.63	1.59
3	A	401	NAD	O4B-C1B	3.81	1.45	1.40
2	A	400	UD1	PB-O3A	3.59	1.63	1.59
2	A	400	UD1	C2-N1	3.55	1.44	1.38
3	B	401	NAD	O4D-C1D	3.31	1.45	1.40
3	B	401	NAD	C2A-N1A	3.16	1.39	1.33
3	B	401	NAD	C2N-N1N	3.07	1.38	1.35
3	A	401	NAD	C2A-N1A	3.00	1.39	1.33
3	B	401	NAD	C2A-N3A	2.83	1.36	1.32
2	B	400	UD1	PA-O3A	2.81	1.62	1.59
3	A	401	NAD	C2N-N1N	2.79	1.38	1.35
2	B	400	UD1	C2-N1	2.77	1.42	1.38
3	B	401	NAD	O4B-C1B	2.64	1.44	1.40
3	B	401	NAD	PA-O3	2.60	1.62	1.59
3	B	401	NAD	C4N-C3N	2.41	1.43	1.39
2	A	400	UD1	O2-C2	2.38	1.27	1.23
3	A	401	NAD	PN-O2N	-2.25	1.44	1.55

All (43) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
3	B	401	NAD	C3N-C7N-N7N	9.86	129.89	117.74
3	A	401	NAD	O4B-C1B-N9A	-8.33	97.70	108.75
3	B	401	NAD	O7N-C7N-N7N	-7.41	111.90	122.62
3	B	401	NAD	N3A-C2A-N1A	-7.35	118.69	128.67
3	B	401	NAD	O4B-C1B-N9A	-7.16	99.26	108.75
3	A	401	NAD	O7N-C7N-C3N	6.73	127.83	119.60
3	A	401	NAD	O7N-C7N-N7N	-6.55	113.15	122.62
3	A	401	NAD	N3A-C2A-N1A	-6.35	120.05	128.67
2	B	400	UD1	C4-N3-C2	-5.97	119.20	126.61
2	A	400	UD1	C5-C4-N3	5.74	122.84	114.80
2	A	400	UD1	C4-N3-C2	-5.63	119.62	126.61
2	B	400	UD1	C5-C4-N3	4.81	121.53	114.80
2	A	400	UD1	O4-C4-C5	-4.35	117.66	125.16
2	B	400	UD1	N3-C2-N1	4.22	120.39	114.89
3	B	401	NAD	C4A-C5A-N7A	-3.39	105.75	109.34
3	B	401	NAD	C4D-O4D-C1D	3.17	112.83	109.92
2	A	400	UD1	O5'-C1'-O1'	-3.03	107.40	111.36
3	B	401	NAD	C4B-O4B-C1B	-3.02	107.16	109.92
2	A	400	UD1	N3-C2-N1	3.00	118.80	114.89
2	B	400	UD1	O3A-PA-O1A	-2.98	101.75	110.70
3	B	401	NAD	C5A-C6A-N6A	2.94	124.80	120.31
3	B	401	NAD	C6N-N1N-C2N	-2.91	119.41	121.88
2	B	400	UD1	O4-C4-C5	-2.80	120.33	125.16
2	A	400	UD1	O6'-C6'-C5'	-2.76	101.95	111.33
3	B	401	NAD	C5N-C4N-C3N	-2.75	117.67	120.36
2	B	400	UD1	O6'-C6'-C5'	-2.67	102.24	111.33
3	B	401	NAD	O3-PA-O1A	-2.66	102.70	110.70
3	A	401	NAD	C1B-N9A-C4A	-2.62	122.05	126.64
3	B	401	NAD	C1B-N9A-C4A	-2.61	122.06	126.64
3	A	401	NAD	C4B-O4B-C1B	-2.60	107.55	109.92
2	A	400	UD1	O2B-PB-O3A	2.56	114.20	107.27
3	B	401	NAD	C2N-C3N-C4N	2.53	121.20	118.26
3	B	401	NAD	O3D-C3D-C4D	-2.43	104.11	111.08
3	A	401	NAD	O3-PA-O1A	-2.42	103.43	110.70
2	A	400	UD1	O5'-C5'-C4'	2.41	114.04	109.70
2	B	400	UD1	C1'-O5'-C5'	-2.35	109.12	113.72
2	A	400	UD1	C1'-C2'-N2'	-2.34	106.98	110.92
2	A	400	UD1	O3A-PB-O1B	-2.18	104.15	110.70
2	A	400	UD1	C2B-C1B-N1	2.17	119.29	113.25
2	A	400	UD1	C3'-C2'-N2'	-2.11	106.73	110.62
2	B	400	UD1	C5-C6-N1	-2.10	118.42	121.84
2	B	400	UD1	O2-C2-N3	-2.02	117.77	121.49
3	B	401	NAD	O3B-C3B-C4B	-2.00	105.33	111.08

There are no chirality outliers.

All (13) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	400	UD1	C1'-O1'-PB-O3A
2	A	400	UD1	PB-O3A-PA-O5B
2	B	400	UD1	C1'-O1'-PB-O3A
3	A	401	NAD	C5D-O5D-PN-O2N
3	B	401	NAD	C5D-O5D-PN-O2N
2	B	400	UD1	PB-O3A-PA-O5B
3	A	401	NAD	C5D-O5D-PN-O3
2	A	400	UD1	C1'-O1'-PB-O1B
3	B	401	NAD	O4B-C4B-C5B-O5B
2	B	400	UD1	PA-O3A-PB-O1B
2	B	400	UD1	PA-O3A-PB-O2B
3	A	401	NAD	O4B-C4B-C5B-O5B
3	A	401	NAD	PA-O3-PN-O1N

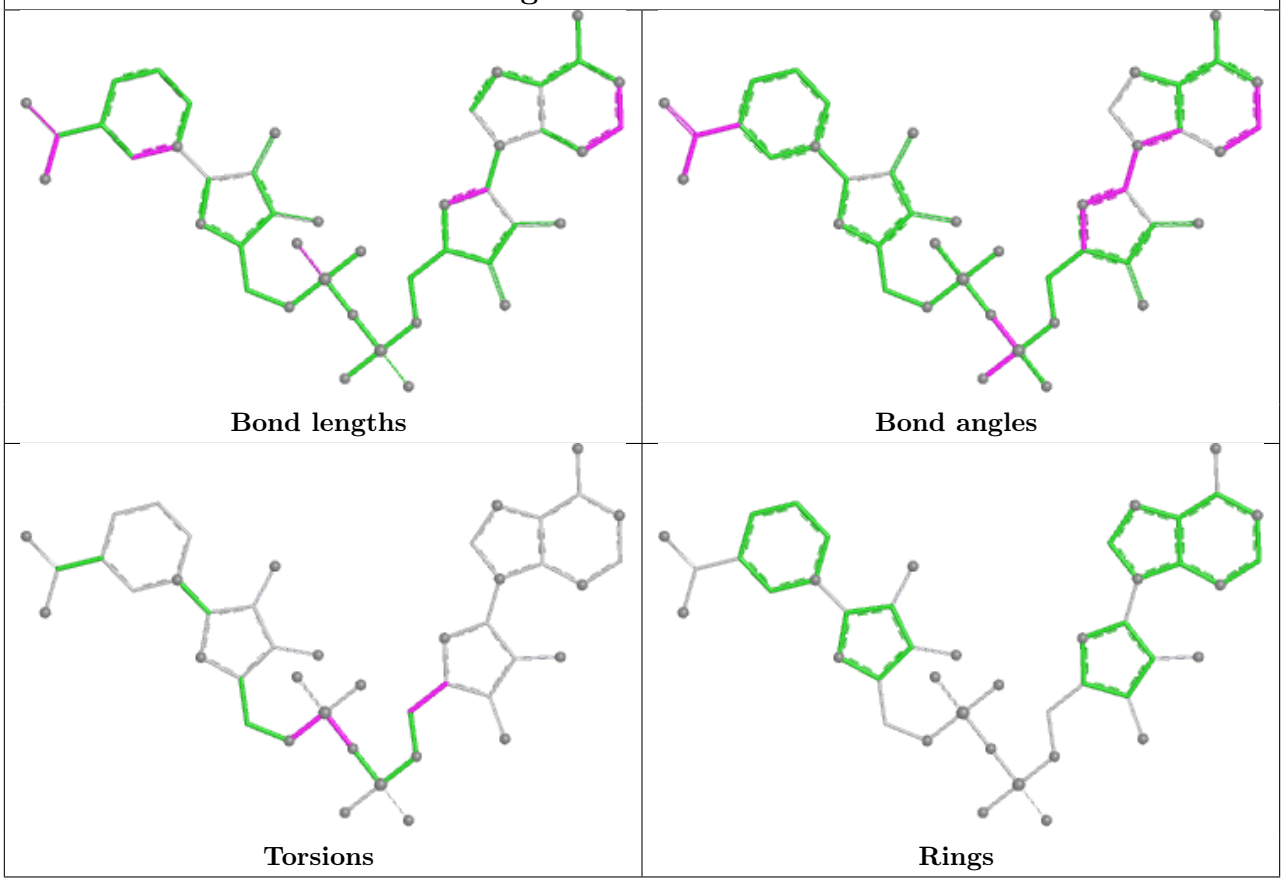
There are no ring outliers.

4 monomers are involved in 4 short contacts:

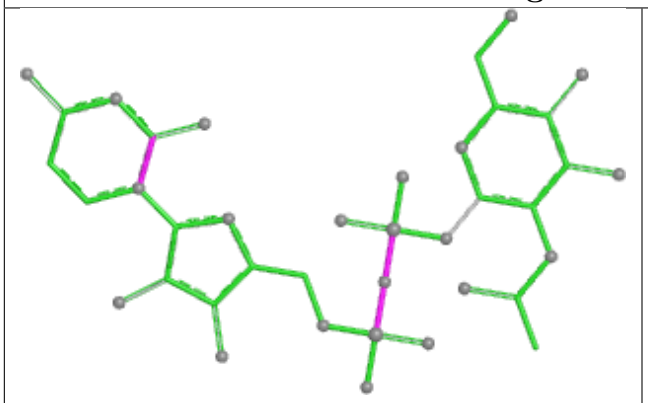
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	401	NAD	2	0
2	B	400	UD1	2	0
2	A	400	UD1	2	0
3	B	401	NAD	2	0

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 then 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.

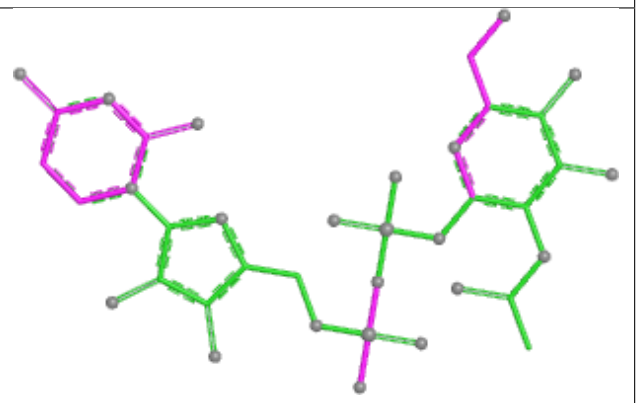
Ligand NAD A 401



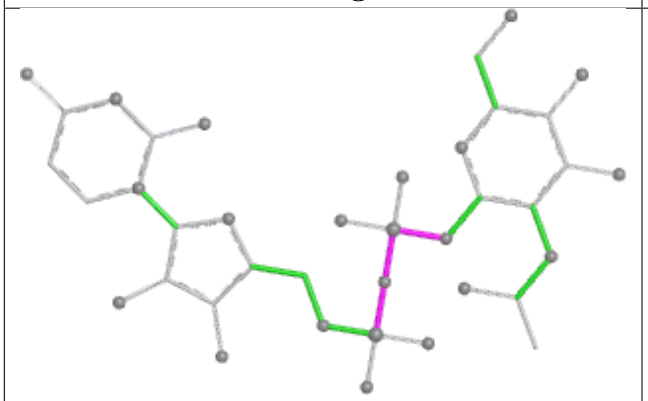
Ligand UD1 B 400



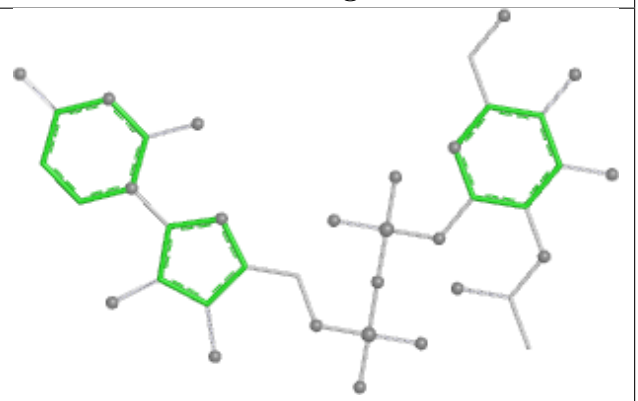
Bond lengths



Bond angles

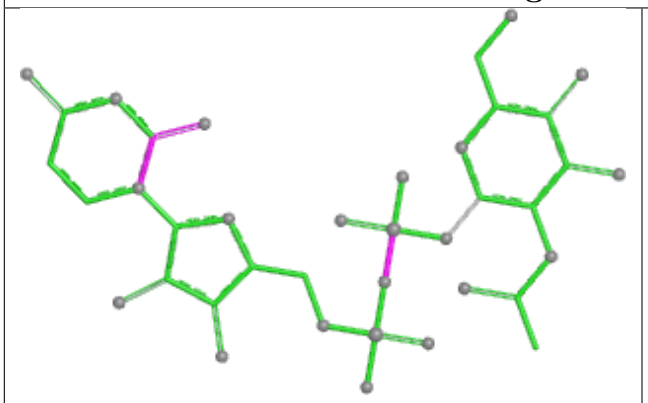


Torsions

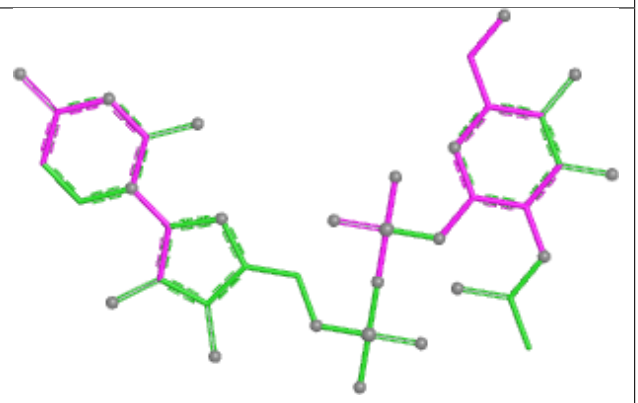


Rings

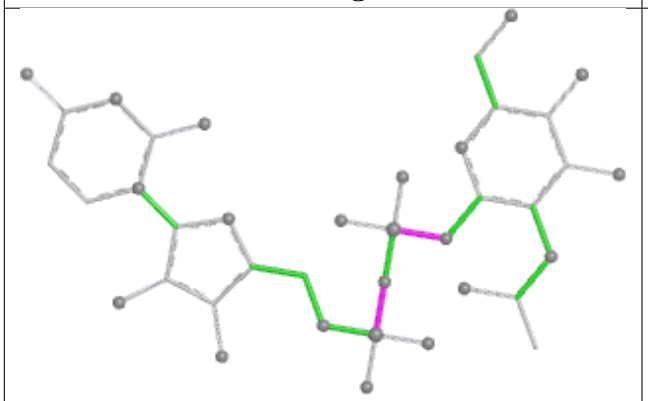
Ligand UD1 A 400



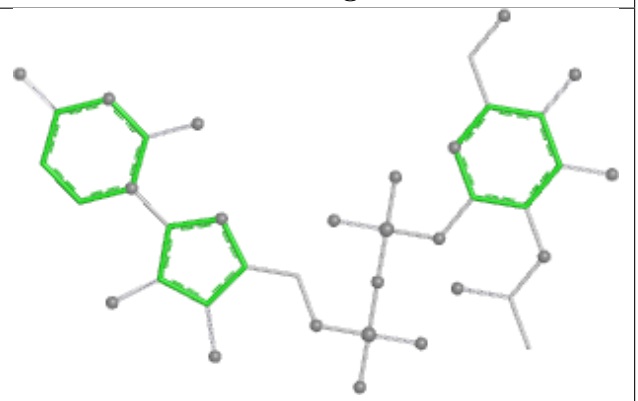
Bond lengths



Bond angles

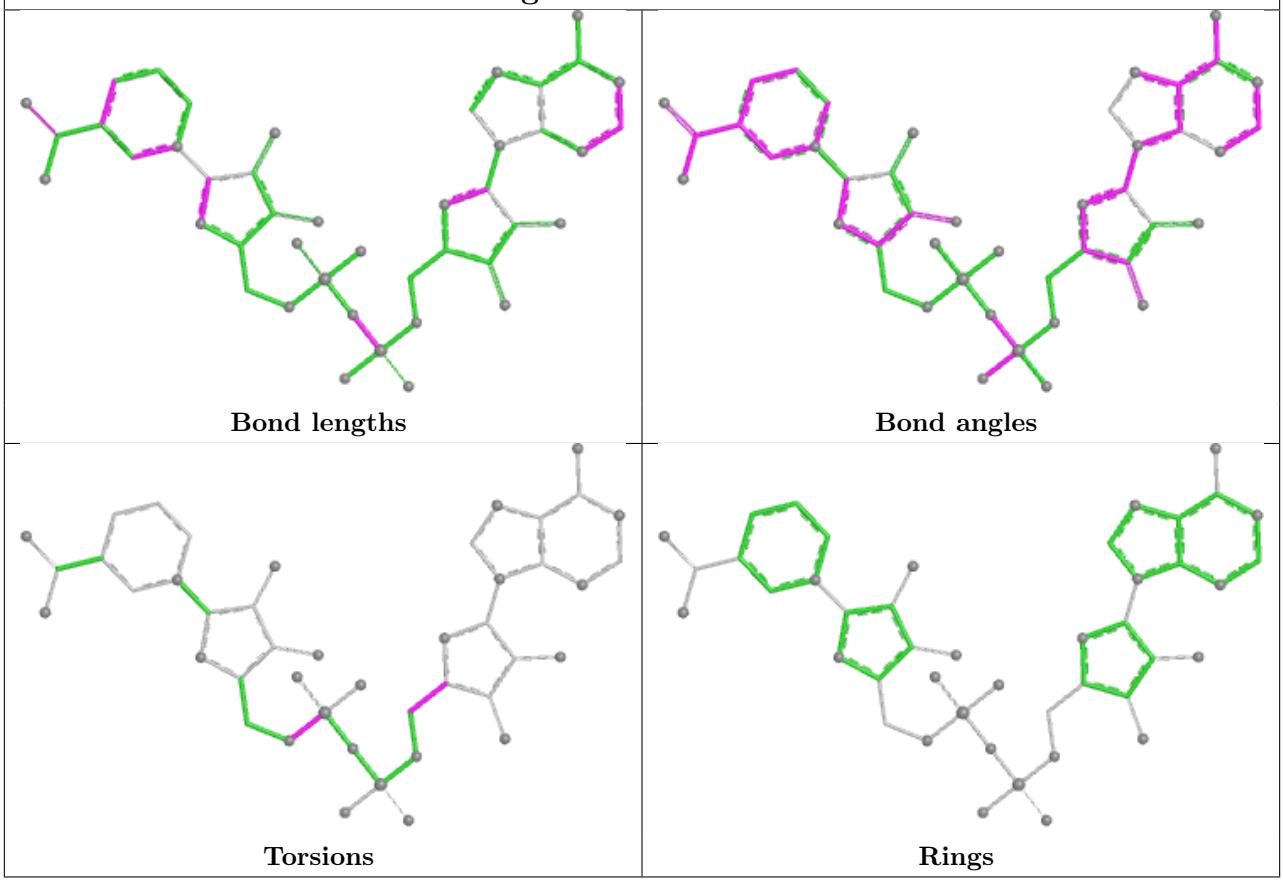


Torsions



Rings

Ligand NAD B 401



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	303/321 (94%)	0.47	33 (10%)	5	6	21, 36, 65, 84	0
1	B	301/321 (93%)	0.49	33 (10%)	5	6	23, 39, 69, 83	0
All	All	604/642 (94%)	0.48	66 (10%)	5	6	21, 38, 68, 84	0

All (66) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	286	SER	7.8
1	B	50	GLY	6.7
1	A	257	GLU	6.6
1	B	51	THR	5.9
1	B	87	LYS	5.7
1	B	47	PRO	5.6
1	A	87	LYS	5.5
1	B	283	GLY	5.5
1	A	287	GLY	5.2
1	A	161	SER	5.1
1	A	285	ARG	4.9
1	A	308	ASP	4.8
1	A	256	ALA	4.7
1	B	310	ILE	4.6
1	B	287	GLY	4.6
1	B	308	ASP	4.5
1	B	262	GLN	4.4
1	B	309	ASP	4.3
1	A	307	LEU	4.2
1	B	49	GLU	4.0
1	A	284	GLU	4.0
1	B	254	PRO	4.0
1	A	306	ASP	3.9
1	A	254	PRO	3.8

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Mol	Chain	Res	Type	RSRZ
1	B	307	LEU	3.7
1	B	88	GLN	3.5
1	B	284	GLU	3.4
1	B	48	PRO	3.4
1	A	88	GLN	3.3
1	B	257	GLU	3.3
1	A	288	GLY	3.3
1	A	185	VAL	3.2
1	A	262	GLN	3.2
1	B	306	ASP	2.9
1	B	197	GLU	2.8
1	B	251	ALA	2.8
1	A	309	ASP	2.7
1	B	185	VAL	2.7
1	B	161	SER	2.7
1	B	266	ASN	2.7
1	B	255	ALA	2.6
1	B	52	GLY	2.6
1	B	260	ARG	2.6
1	A	251	ALA	2.5
1	A	188	LEU	2.5
1	B	162	VAL	2.5
1	A	255	ALA	2.5
1	B	304	SER	2.5
1	A	195	ARG	2.4
1	A	84	ARG	2.4
1	A	252	THR	2.4
1	A	12	ILE	2.4
1	B	188	LEU	2.3
1	A	83	PRO	2.3
1	B	303	GLN	2.3
1	B	265	PRO	2.3
1	A	50	GLY	2.3
1	B	43	PRO	2.2
1	A	76	LEU	2.2
1	A	20	GLY	2.1
1	A	238	GLN	2.1
1	B	194	THR	2.1
1	A	197	GLU	2.0
1	A	259	ALA	2.0
1	A	186	PRO	2.0
1	A	86	PHE	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)

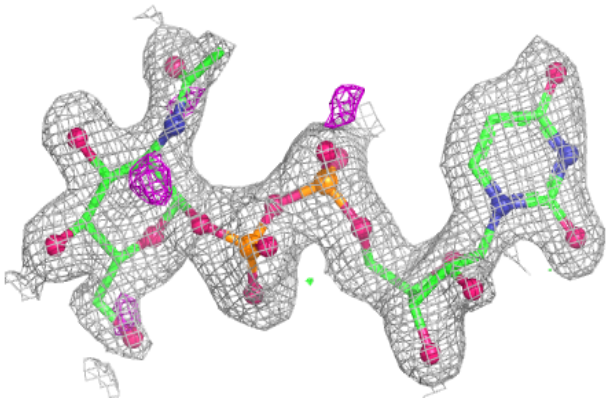
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

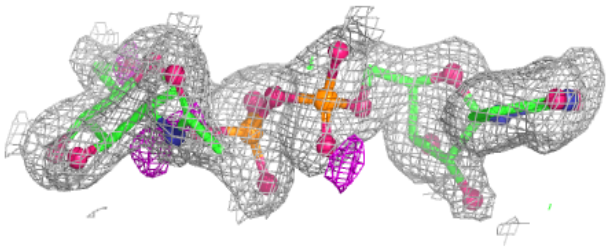
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
2	UD1	A	400	39/39	0.94	0.10	33,40,50,53	0
2	UD1	B	400	39/39	0.95	0.09	36,43,53,56	0
3	NAD	B	401	44/44	0.96	0.10	28,31,33,38	0
3	NAD	A	401	44/44	0.98	0.11	21,25,29,34	0

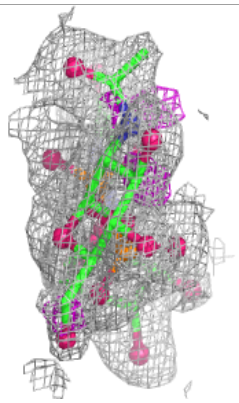
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around UD1 A 400:

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

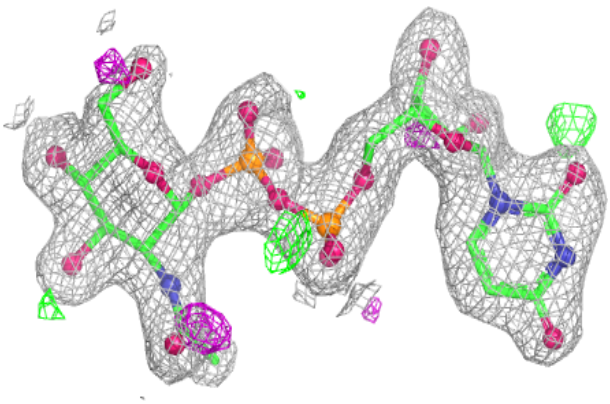


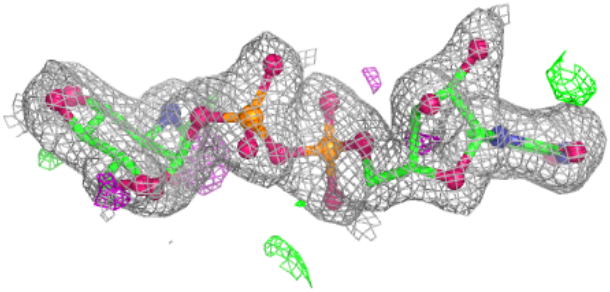


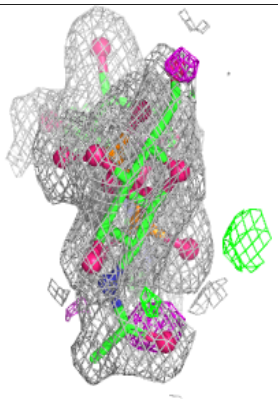


Electron density around UD1 B 400:

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

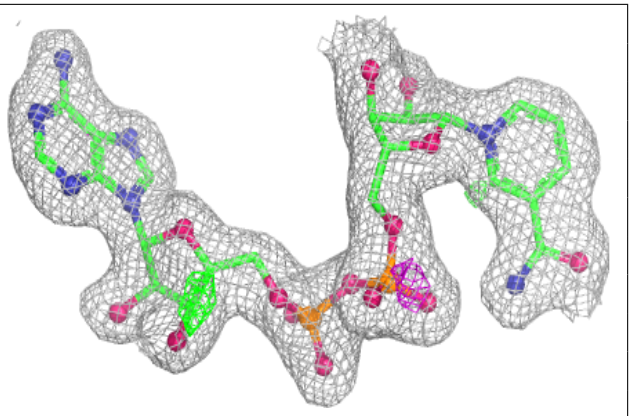


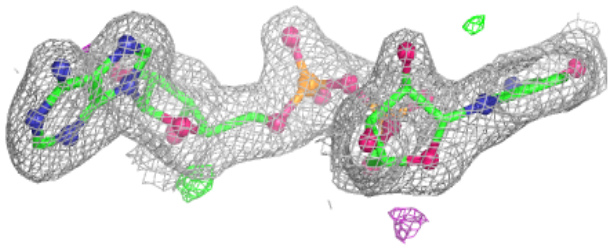


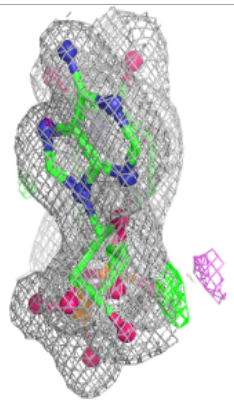


Electron density around NAD B 401:

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

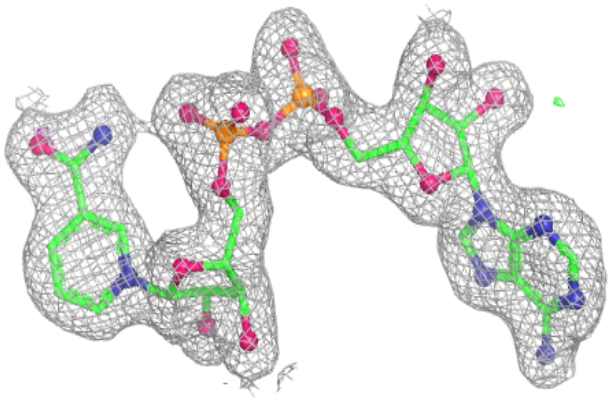


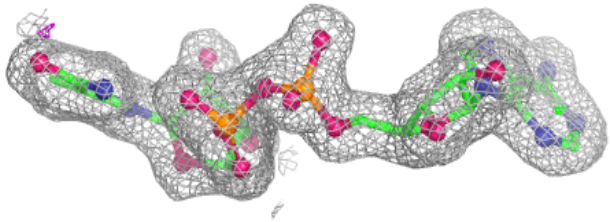


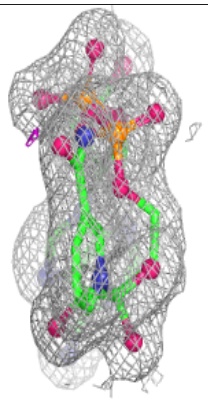


Electron density around NAD A 401:

$2mF_o$-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.