



# Full wwPDB X-ray Structure Validation Report ⓘ

Apr 15, 2026 – 12:55 AM UTC

PDB ID : 9UEE / pdb\_00009uee  
Title : Crystal structure of (R)-carbonyl reductase mutant I51L/Y61F/D147E from Candida Parapsilosis in complex with NAD  
Authors : Gu, J.; Nie, Y.  
Deposited on : 2025-04-08  
Resolution : 2.50 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](mailto: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>.

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The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

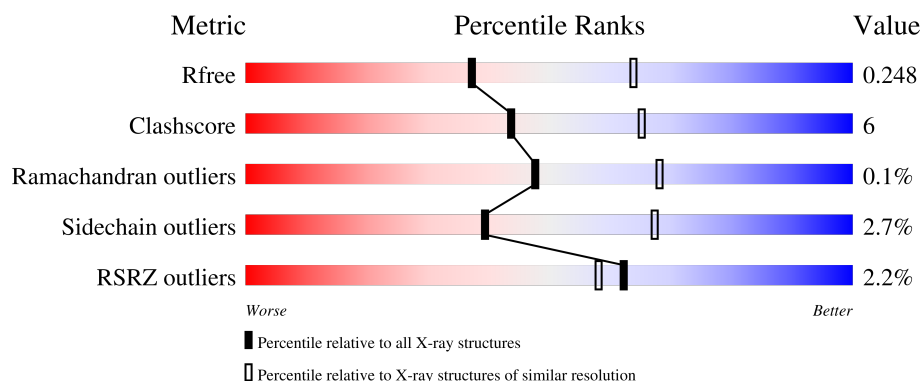
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.50 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
$R_{free}$	180053	5829 (2.50-2.50)
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)
RSRZ outliers	180081	5833 (2.50-2.50)

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  $\geq 3$ , 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	336	<div> <div>0%</div> <div>90%</div> <div>10%</div> <div>.</div> </div>
1	B	336	<div> <div>0%</div> <div>89%</div> <div>11%</div> <div>.</div> </div>
1	C	336	<div> <div>2%</div> <div>78%</div> <div>19%</div> <div>.</div> </div>
1	D	336	<div> <div>5%</div> <div>77%</div> <div>21%</div> <div>.</div> </div>

## 2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 10346 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 (R)-specific carbonyl reductase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	336	Total	C	N	O	S	0	0	0
			2530	1607	435	475	13			
1	B	336	Total	C	N	O	S	0	0	0
			2530	1607	435	475	13			
1	C	336	Total	C	N	O	S	0	0	0
			2530	1607	435	475	13			
1	D	336	Total	C	N	O	S	0	0	0
			2530	1607	435	475	13			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	51	LEU	ILE	engineered mutation	UNP A1X808
A	61	PHE	TYR	engineered mutation	UNP A1X808
A	147	GLU	ASP	engineered mutation	UNP A1X808
B	51	LEU	ILE	engineered mutation	UNP A1X808
B	61	PHE	TYR	engineered mutation	UNP A1X808
B	147	GLU	ASP	engineered mutation	UNP A1X808
C	51	LEU	ILE	engineered mutation	UNP A1X808
C	61	PHE	TYR	engineered mutation	UNP A1X808
C	147	GLU	ASP	engineered mutation	UNP A1X808
D	51	LEU	ILE	engineered mutation	UNP A1X808
D	61	PHE	TYR	engineered mutation	UNP A1X808
D	147	GLU	ASP	engineered mutation	UNP A1X808

- Molecule 2 is NICOTINAMIDE-ADENINE-DINUCLEOTIDE (CCD ID: NAD) (formula:  $C_{21}H_{27}N_7O_{14}P_2$ ) (labeled as "Ligand of Interest" by depositor).



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

- Molecule 3 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Zn	0	0
			1	1		
3	B	1	Total	Zn	0	0
			1	1		
3	C	1	Total	Zn	0	0
			1	1		
3	D	1	Total	Zn	0	0
			1	1		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	23	Total	O	0	0
			23	23		

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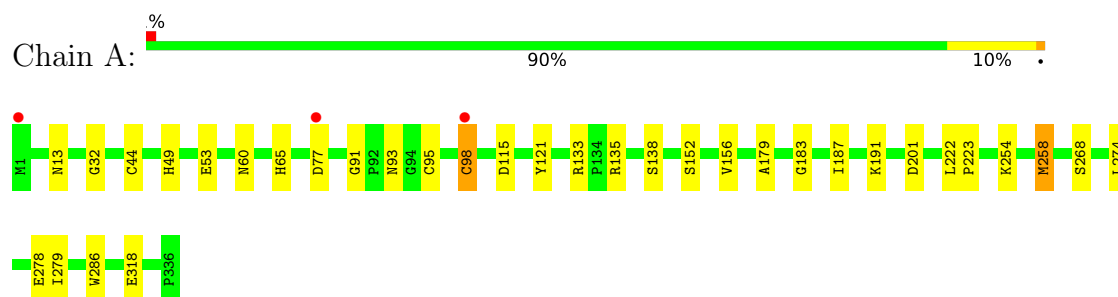
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	8	Total	O	0	0
			8	8		
4	C	9	Total	O	0	0
			9	9		
4	D	6	Total	O	0	0
			6	6		

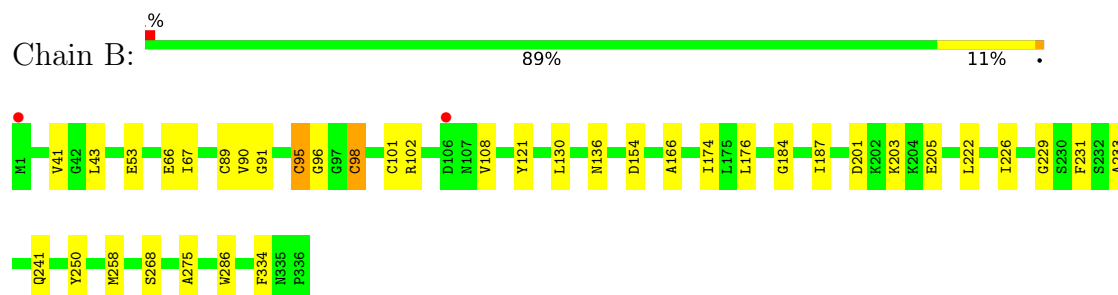
### 3 Residue-property plots [i](#)

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.

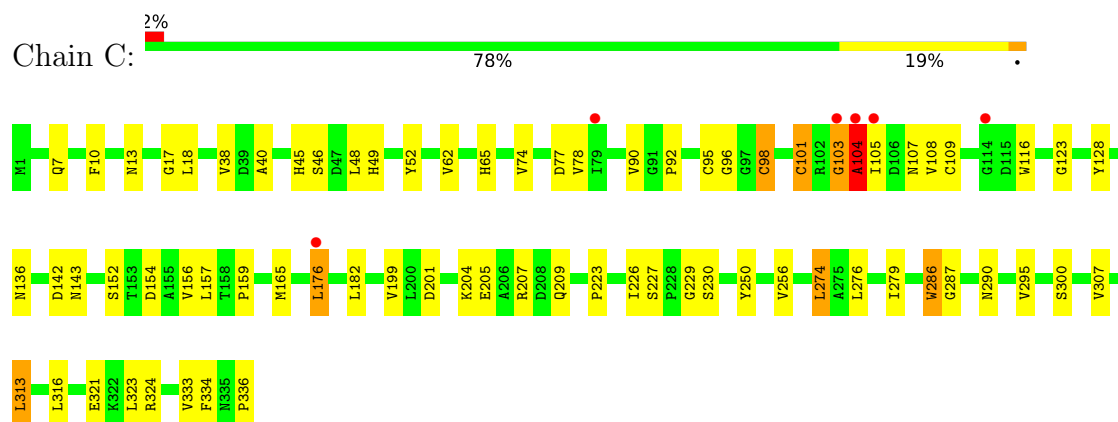
- Molecule 1: (R)-specific carbonyl reductase



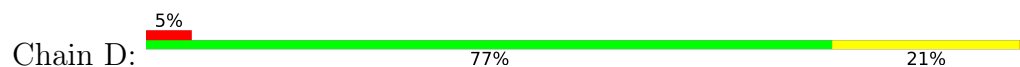
- Molecule 1: (R)-specific carbonyl reductase

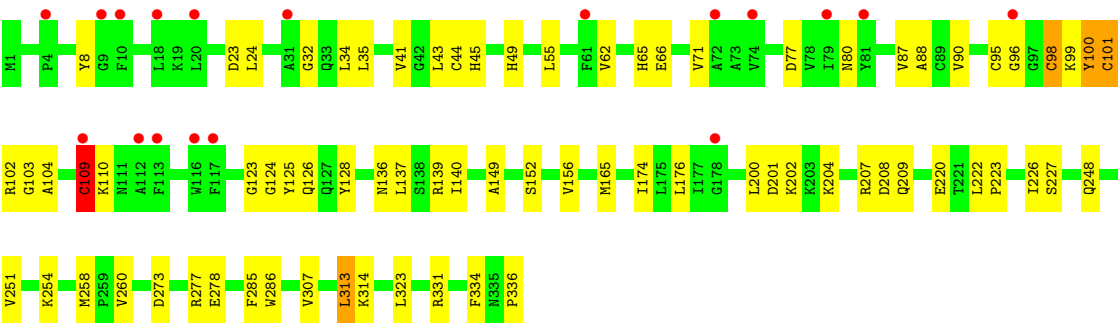


- Molecule 1: (R)-specific carbonyl reductase



- Molecule 1: (R)-specific carbonyl reductase





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	86.63Å 108.93Å 148.01Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	37.38 – 2.50 37.38 – 2.50	Depositor EDS
% Data completeness (in resolution range)	99.0 (37.38-2.50) 99.0 (37.38-2.50)	Depositor EDS
$R_{merge}$	0.12	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.85 (at 2.51Å)	Xtriage
Refinement program	PHENIX 1.15.1_3469	Depositor
R, $R_{free}$	0.210 , 0.249 0.210 , 0.248	Depositor DCC
$R_{free}$ test set	2438 reflections (4.96%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	42.5	Xtriage
Anisotropy	0.161	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.34 , 50.6	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.43$ , $\langle L^2 \rangle = 0.26$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.93	EDS
Total number of atoms	10346	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	59.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.31% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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: ZN, 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	0.11	0/2579	0.32	0/3492
1	B	0.23	0/2579	0.40	3/3492 (0.1%)
1	C	0.31	1/2579 (0.0%)	0.49	6/3492 (0.2%)
1	D	0.17	0/2579	0.42	3/3492 (0.1%)
All	All	0.22	1/10316 (0.0%)	0.41	12/13968 (0.1%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	109	CYS	N-CA	5.37	1.53	1.46

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	103	GLY	N-CA-C	8.14	122.50	112.73
1	D	110	LYS	N-CA-C	7.48	119.44	111.28
1	D	101	CYS	CA-CB-SG	7.44	131.51	114.40
1	B	95	CYS	N-CA-C	7.34	119.28	111.28
1	C	101	CYS	CA-CB-SG	6.81	130.06	114.40
1	C	101	CYS	N-CA-C	-6.15	104.58	111.28
1	B	95	CYS	N-CA-CB	-5.60	101.88	110.12
1	C	109	CYS	CA-CB-SG	5.52	127.10	114.40
1	C	103	GLY	CA-C-O	-5.27	115.32	120.75
1	B	108	VAL	N-CA-C	5.07	116.42	111.91
1	C	104	ALA	N-CA-C	-5.07	100.00	110.80
1	D	109	CYS	CB-CA-C	5.06	118.13	109.72

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	2530	0	2541	18	0
1	B	2530	0	2541	20	0
1	C	2530	0	2541	45	0
1	D	2530	0	2541	48	0
2	A	44	0	26	2	0
2	B	44	0	26	1	0
2	C	44	0	26	2	0
2	D	44	0	26	2	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
3	C	1	0	0	0	0
3	D	1	0	0	0	0
4	A	23	0	0	0	0
4	B	8	0	0	0	0
4	C	9	0	0	0	0
4	D	6	0	0	4	0
All	All	10346	0	10268	126	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 6.

All (126) 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:D:96:GLY:N	1:D:101:CYS:SG	2.46	0.89
1:C:103:GLY:O	1:D:104:ALA:HB2	1.80	0.80
1:A:95:CYS:SG	1:A:98:CYS:N	2.59	0.75
1:C:98:CYS:HB2	1:C:101:CYS:H	1.51	0.74
1:C:46:SER:HA	1:C:49:HIS:HD2	1.56	0.71
1:D:49:HIS:HB3	1:D:55:LEU:HD23	1.74	0.69
1:D:200:LEU:HD11	1:D:222:LEU:HG	1.74	0.69
1:D:251:VAL:O	1:D:277:ARG:NH2	2.24	0.69
1:C:204:LYS:HD2	1:C:207:ARG:HH21	1.57	0.69
1:D:95:CYS:SG	1:D:98:CYS:N	2.67	0.68
1:D:223:PRO:HD2	1:D:226:ILE:HD13	1.75	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:49:HIS:HE1	4:D:505:HOH:O	1.79	0.65
1:C:90:VAL:HB	1:C:136:ASN:HB3	1.81	0.62
1:C:95:CYS:SG	1:C:98:CYS:N	2.73	0.62
1:D:165:MET:HE2	1:D:258:MET:HE1	1.81	0.62
1:A:32:GLY:N	1:A:77:ASP:OD1	2.33	0.61
1:B:166:ALA:HB2	1:B:258:MET:HE3	1.80	0.61
1:C:46:SER:HA	1:C:49:HIS:CD2	2.36	0.61
1:C:96:GLY:N	1:C:101:CYS:SG	2.63	0.59
1:A:49:HIS:HA	1:A:53:GLU:HG3	1.85	0.59
1:D:35:LEU:HD11	1:D:128:TYR:HB3	1.84	0.58
1:B:201:ASP:OD1	2:B:401:NAD:O2B	2.21	0.58
1:D:80:ASN:O	1:D:139:ARG:NH2	2.37	0.58
1:A:254:LYS:HE3	1:A:278:GLU:HB3	1.85	0.57
1:C:227:SER:O	1:C:230:SER:OG	2.21	0.56
1:B:98:CYS:HB2	1:B:101:CYS:H	1.71	0.55
1:C:7:GLN:OE1	1:C:128:TYR:N	2.37	0.55
1:B:121:TYR:OH	1:B:130:LEU:O	2.22	0.54
1:B:241:GLN:OE1	1:B:268:SER:N	2.40	0.54
1:C:274:LEU:HA	1:C:279:ILE:HD12	1.89	0.54
1:D:209:GLN:HE22	1:D:307:VAL:HG11	1.71	0.53
1:D:90:VAL:HB	1:D:136:ASN:HB3	1.90	0.53
1:C:104:ALA:HB2	1:D:103:GLY:HA2	1.89	0.53
1:C:209:GLN:NE2	1:C:307:VAL:HG21	2.23	0.53
1:C:201:ASP:OD1	2:C:401:NAD:O2B	2.22	0.52
1:C:95:CYS:N	1:C:101:CYS:SG	2.82	0.52
1:D:99:LYS:HE3	1:D:100:TYR:CZ	2.44	0.52
1:D:95:CYS:N	1:D:101:CYS:SG	2.83	0.52
1:D:88:ALA:HB2	1:D:140:ILE:HD11	1.92	0.52
1:D:45:HIS:HE1	4:D:505:HOH:O	1.92	0.51
1:C:13:ASN:O	1:C:17:GLY:N	2.43	0.51
1:A:13:ASN:ND2	1:A:60:ASN:OD1	2.40	0.51
1:D:204:LYS:NZ	1:D:208:ASP:OD2	2.43	0.51
1:B:95:CYS:SG	1:B:96:GLY:N	2.84	0.50
1:D:32:GLY:N	1:D:77:ASP:OD1	2.45	0.50
1:D:124:GLY:O	1:D:126:GLN:N	2.40	0.49
1:B:174:ILE:HG22	1:B:233:ALA:HB3	1.94	0.49
1:C:38:VAL:HB	1:C:336:PRO:HG2	1.94	0.49
1:D:44:CYS:HB3	1:D:65:HIS:CE1	2.47	0.49
1:D:254:LYS:NZ	1:D:278:GLU:OE1	2.39	0.49
1:C:209:GLN:HE22	1:C:307:VAL:HG21	1.77	0.49
1:B:275:ALA:O	1:D:285:PHE:HB2	2.13	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:274:LEU:HD22	1:A:279:ILE:HB	1.94	0.49
1:B:203:LYS:HD3	1:B:205:GLU:HB2	1.93	0.49
1:C:152:SER:HA	1:C:156:VAL:HB	1.94	0.49
1:C:290:ASN:OD1	1:D:102:ARG:NH2	2.46	0.48
1:D:43:LEU:HB2	1:D:334:PHE:HE2	1.78	0.48
1:D:62:VAL:O	1:D:123:GLY:N	2.44	0.48
1:C:45:HIS:HB3	2:C:401:NAD:H3D	1.95	0.48
1:D:201:ASP:OD1	2:D:401:NAD:O2B	2.29	0.48
1:C:143:ASN:HD21	1:C:300:SER:HB2	1.79	0.47
1:C:159:PRO:HD3	1:C:182:LEU:HD13	1.95	0.47
1:D:23:ASP:OD1	1:D:23:ASP:N	2.47	0.47
1:A:91:GLY:HA3	1:A:286:TRP:CZ2	2.49	0.47
1:C:105:ILE:O	1:C:108:VAL:HG22	2.15	0.47
1:D:34:LEU:HD13	1:D:71:VAL:HG13	1.96	0.47
1:D:260:VAL:O	2:D:401:NAD:H2N	2.15	0.47
1:C:52:TYR:CE2	1:C:324:ARG:HD2	2.49	0.47
1:D:66:GLU:OE2	1:D:331:ARG:NH2	2.45	0.47
1:B:91:GLY:HA3	1:B:286:TRP:CZ2	2.51	0.46
1:D:44:CYS:HB3	1:D:65:HIS:NE2	2.31	0.46
1:B:41:VAL:HA	1:B:66:GLU:O	2.15	0.46
1:D:49:HIS:CE1	4:D:505:HOH:O	2.61	0.46
1:C:103:GLY:O	1:C:104:ALA:CB	2.64	0.45
1:C:40:ALA:HB1	1:C:333:VAL:HG12	1.98	0.45
1:C:65:HIS:CD2	1:C:154:ASP:OD2	2.68	0.45
1:D:207:ARG:NH1	1:D:220:GLU:HB2	2.31	0.45
1:C:157:LEU:HA	1:C:295:VAL:HG21	1.99	0.45
1:D:140:ILE:HD12	1:D:149:ALA:HB1	1.98	0.45
1:B:231:PHE:HD2	1:B:250:TYR:HB3	1.80	0.45
1:C:286:TRP:HB3	1:C:287:GLY:H	1.60	0.44
1:D:87:VAL:HG21	1:D:137:LEU:HD13	2.00	0.44
1:A:201:ASP:OD1	2:A:401:NAD:O2B	2.31	0.44
1:C:48:LEU:HD13	1:C:323:LEU:HD23	1.99	0.44
1:D:8:TYR:OH	1:D:336:PRO:OXT	2.25	0.44
1:D:152:SER:HA	1:D:156:VAL:HB	2.00	0.44
1:A:115:ASP:OD1	1:A:121:TYR:HB2	2.18	0.43
1:C:176:LEU:O	1:C:199:VAL:HA	2.18	0.43
1:A:152:SER:HA	1:A:156:VAL:HB	1.99	0.43
1:C:74:VAL:HG22	1:C:78:VAL:HB	2.00	0.43
1:B:184:GLY:HA2	1:B:187:ILE:HD12	2.01	0.43
1:A:44:CYS:HB3	1:A:65:HIS:CE1	2.53	0.43
1:B:90:VAL:HB	1:B:136:ASN:HB3	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:222:LEU:HD22	1:B:226:ILE:HD12	2.01	0.43
1:D:98:CYS:HB2	1:D:101:CYS:H	1.82	0.43
1:A:187:ILE:O	1:A:191:LYS:HG2	2.19	0.42
1:A:135:ARG:NH2	1:B:102:ARG:HD2	2.34	0.42
1:A:201:ASP:CG	2:A:401:NAD:HO2A	2.26	0.42
1:D:313:LEU:HD12	1:D:336:PRO:HA	2.02	0.42
1:A:258:MET:HE2	1:A:258:MET:HB2	1.71	0.42
1:B:66:GLU:HG2	1:B:154:ASP:HB3	2.02	0.42
1:C:165:MET:HE2	1:C:165:MET:HB3	1.88	0.42
1:D:45:HIS:CE1	4:D:505:HOH:O	2.70	0.42
1:C:223:PRO:HD2	1:C:226:ILE:HD12	2.02	0.42
1:C:142:ASP:OD1	1:C:142:ASP:N	2.53	0.42
1:D:248:GLN:HE21	1:D:273:ASP:HB3	1.85	0.41
1:D:313:LEU:HB2	1:D:336:PRO:HA	2.01	0.41
1:D:98:CYS:HB2	1:D:109:CYS:HB2	1.80	0.41
1:A:93:ASN:O	1:A:133:ARG:NH1	2.54	0.41
1:C:62:VAL:O	1:C:123:GLY:N	2.54	0.41
1:C:92:PRO:O	1:C:116:TRP:HD1	2.04	0.41
1:C:229:GLY:H	1:C:250:TYR:HA	1.85	0.41
1:A:222:LEU:HA	1:A:223:PRO:HD3	1.95	0.41
1:C:107:ASN:OD1	1:C:107:ASN:N	2.52	0.41
1:C:10:PHE:HB3	1:C:18:LEU:HD11	2.01	0.41
1:C:205:GLU:N	1:C:205:GLU:OE1	2.54	0.41
1:B:229:GLY:H	1:B:250:TYR:HA	1.86	0.41
1:D:201:ASP:OD1	1:D:202:LYS:N	2.54	0.41
1:D:222:LEU:HA	1:D:223:PRO:HD3	1.97	0.41
1:A:179:ALA:HA	1:A:183:GLY:HA3	2.03	0.40
1:B:67:ILE:HB	1:B:89:CYS:HB2	2.02	0.40
1:C:103:GLY:O	1:C:104:ALA:HB3	2.21	0.40
1:D:41:VAL:HA	1:D:66:GLU:O	2.21	0.40
1:B:43:LEU:HB2	1:B:334:PHE:HE2	1.86	0.40
1:C:95:CYS:SG	1:C:96:GLY:N	2.93	0.40
1:C:313:LEU:HA	1:C:334:PHE:HB3	2.04	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	334/336 (99%)	323 (97%)	11 (3%)	0	100	100
1	B	334/336 (99%)	315 (94%)	19 (6%)	0	100	100
1	C	334/336 (99%)	315 (94%)	18 (5%)	1 (0%)	36	55
1	D	334/336 (99%)	322 (96%)	11 (3%)	1 (0%)	36	55
All	All	1336/1344 (99%)	1275 (95%)	59 (4%)	2 (0%)	48	68

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	104	ALA
1	D	125	TYR

### 5.3.2 Protein sidechains

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	270/270 (100%)	265 (98%)	5 (2%)	50	76
1	B	270/270 (100%)	267 (99%)	3 (1%)	65	84
1	C	270/270 (100%)	260 (96%)	10 (4%)	30	57
1	D	270/270 (100%)	259 (96%)	11 (4%)	27	53
All	All	1080/1080 (100%)	1051 (97%)	29 (3%)	39	67

All (29) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	98	CYS
1	A	138	SER
1	A	258	MET
1	A	268	SER
1	A	318	GLU
1	B	53	GLU
1	B	98	CYS
1	B	176	LEU
1	C	77	ASP
1	C	98	CYS
1	C	176	LEU
1	C	256	VAL
1	C	274	LEU
1	C	276	LEU
1	C	286	TRP
1	C	313	LEU
1	C	316	LEU
1	C	321	GLU
1	D	24	LEU
1	D	98	CYS
1	D	100	TYR
1	D	109	CYS
1	D	174	ILE
1	D	176	LEU
1	D	227	SER
1	D	286	TRP
1	D	313	LEU
1	D	314	LYS
1	D	323	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	325	ASN
1	C	335	ASN
1	D	60	ASN
1	D	209	GLN
1	D	325	ASN

### 5.3.3 RNA ⓘ

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 [i](#)

Of 8 ligands modelled in this entry, 4 are monoatomic - leaving 4 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	A	401	-	46,48,48	1.22	5 (10%)	64,73,73	1.56	10 (15%)
2	NAD	D	401	-	46,48,48	1.18	4 (8%)	64,73,73	1.58	10 (15%)
2	NAD	C	401	-	46,48,48	1.20	4 (8%)	64,73,73	1.56	9 (14%)
2	NAD	B	401	-	46,48,48	1.18	4 (8%)	64,73,73	1.57	10 (15%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	NAD	A	401	-	-	11/30/62/62	0/5/5/5
2	NAD	D	401	-	-	12/30/62/62	0/5/5/5
2	NAD	C	401	-	-	7/30/62/62	0/5/5/5
2	NAD	B	401	-	-	7/30/62/62	0/5/5/5

All (17) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	401	NAD	C5A-C4A	4.79	1.47	1.39

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	401	NAD	C5A-C4A	4.71	1.47	1.39
2	D	401	NAD	C5A-C4A	4.71	1.47	1.39
2	B	401	NAD	C5A-C4A	4.68	1.47	1.39
2	A	401	NAD	C5A-C6A	2.81	1.48	1.41
2	B	401	NAD	C5A-C6A	2.80	1.48	1.41
2	D	401	NAD	C5A-C6A	2.79	1.48	1.41
2	C	401	NAD	C5A-C6A	2.75	1.48	1.41
2	D	401	NAD	C8A-N7A	2.48	1.36	1.31
2	A	401	NAD	C8A-N7A	2.42	1.36	1.31
2	B	401	NAD	C8A-N7A	2.41	1.36	1.31
2	C	401	NAD	C8A-N7A	2.34	1.36	1.31
2	C	401	NAD	C5A-N7A	-2.31	1.34	1.39
2	A	401	NAD	C5A-N7A	-2.23	1.35	1.39
2	D	401	NAD	C5A-N7A	-2.17	1.35	1.39
2	B	401	NAD	C5A-N7A	-2.16	1.35	1.39
2	A	401	NAD	PA-O3	2.07	1.61	1.59

All (39) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	401	NAD	C5A-C4A-N3A	-5.98	118.48	126.72
2	D	401	NAD	C5A-C4A-N3A	-5.98	118.49	126.72
2	A	401	NAD	C5A-C4A-N3A	-5.79	118.75	126.72
2	B	401	NAD	C5A-C4A-N3A	-5.58	119.04	126.72
2	C	401	NAD	N3A-C4A-N9A	4.77	135.27	127.17
2	D	401	NAD	N3A-C4A-N9A	4.57	134.94	127.17
2	A	401	NAD	N3A-C4A-N9A	4.48	134.79	127.17
2	B	401	NAD	N3A-C4A-N9A	4.37	134.59	127.17
2	D	401	NAD	C2A-N3A-C4A	3.78	121.07	111.83
2	C	401	NAD	C2A-N3A-C4A	3.72	120.91	111.83
2	B	401	NAD	C2A-N3A-C4A	3.71	120.90	111.83
2	D	401	NAD	C4A-C5A-N7A	-3.71	106.34	110.58
2	A	401	NAD	C4A-C5A-N7A	-3.70	106.35	110.58
2	A	401	NAD	C2A-N3A-C4A	3.69	120.83	111.83
2	C	401	NAD	C4A-C5A-N7A	-3.62	106.45	110.58
2	B	401	NAD	C4A-C5A-N7A	-3.61	106.45	110.58
2	B	401	NAD	N3A-C2A-N1A	-3.41	123.41	128.58
2	D	401	NAD	N3A-C2A-N1A	-3.28	123.62	128.58
2	C	401	NAD	N3A-C2A-N1A	-3.22	123.70	128.58
2	A	401	NAD	N3A-C2A-N1A	-3.16	123.81	128.58
2	B	401	NAD	C4A-N9A-C8A	2.82	108.70	105.74
2	C	401	NAD	C5A-N7A-C8A	2.69	107.68	103.45

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	401	NAD	C4A-N9A-C8A	2.69	108.56	105.74
2	D	401	NAD	C5A-N7A-C8A	2.68	107.66	103.45
2	A	401	NAD	C4A-N9A-C8A	2.68	108.55	105.74
2	A	401	NAD	C5A-N7A-C8A	2.67	107.65	103.45
2	B	401	NAD	C5A-N7A-C8A	2.63	107.59	103.45
2	B	401	NAD	O4B-C1B-N9A	2.57	113.03	108.09
2	D	401	NAD	C4A-N9A-C8A	2.54	108.41	105.74
2	B	401	NAD	C6A-C5A-N7A	2.36	136.65	132.09
2	A	401	NAD	C6A-C5A-N7A	2.29	136.51	132.09
2	D	401	NAD	C6A-C5A-N7A	2.19	136.31	132.09
2	B	401	NAD	N9A-C8A-N7A	-2.14	110.90	113.94
2	D	401	NAD	O4B-C1B-N9A	2.13	112.19	108.09
2	A	401	NAD	N9A-C8A-N7A	-2.08	110.99	113.94
2	D	401	NAD	N9A-C8A-N7A	-2.06	111.01	113.94
2	A	401	NAD	O4B-C1B-N9A	2.06	112.05	108.09
2	C	401	NAD	C6A-C5A-N7A	2.05	136.05	132.09
2	C	401	NAD	N9A-C8A-N7A	-2.02	111.06	113.94

There are no chirality outliers.

All (37) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	401	NAD	C5B-O5B-PA-O2A
2	A	401	NAD	C5B-O5B-PA-O3
2	A	401	NAD	O4B-C4B-C5B-O5B
2	A	401	NAD	C3B-C4B-C5B-O5B
2	A	401	NAD	O4D-C1D-N1N-C2N
2	A	401	NAD	O4D-C1D-N1N-C6N
2	A	401	NAD	C2D-C1D-N1N-C2N
2	A	401	NAD	C2D-C1D-N1N-C6N
2	B	401	NAD	C5B-O5B-PA-O3
2	B	401	NAD	O4D-C1D-N1N-C2N
2	B	401	NAD	O4D-C1D-N1N-C6N
2	B	401	NAD	C2D-C1D-N1N-C2N
2	B	401	NAD	C2D-C1D-N1N-C6N
2	C	401	NAD	O4D-C1D-N1N-C2N
2	C	401	NAD	O4D-C1D-N1N-C6N
2	C	401	NAD	C2D-C1D-N1N-C2N
2	D	401	NAD	C5B-O5B-PA-O3
2	D	401	NAD	O4D-C1D-N1N-C2N
2	D	401	NAD	O4D-C1D-N1N-C6N
2	D	401	NAD	C2D-C1D-N1N-C2N

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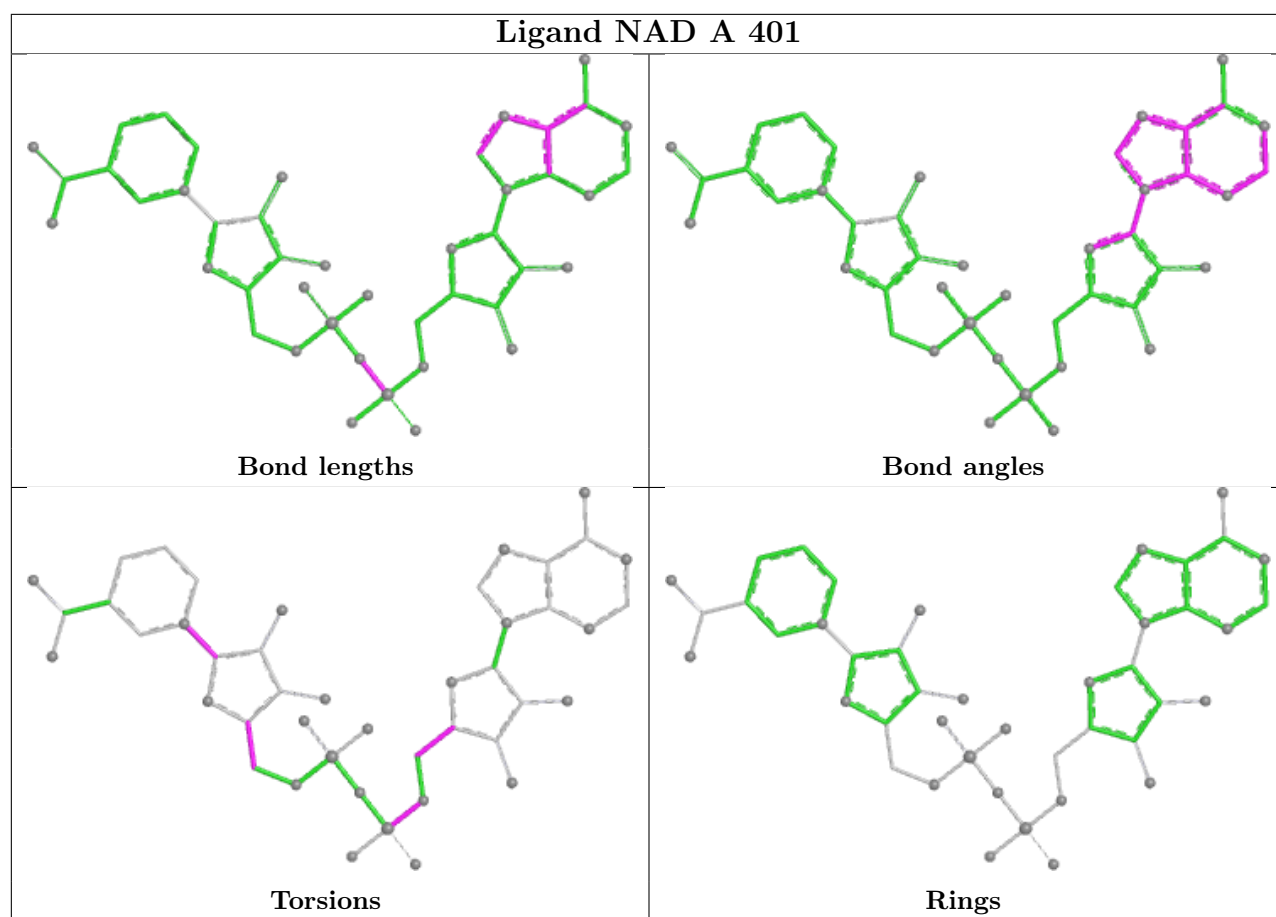
Mol	Chain	Res	Type	Atoms
2	D	401	NAD	C3B-C4B-C5B-O5B
2	A	401	NAD	O4D-C4D-C5D-O5D
2	C	401	NAD	O4B-C4B-C5B-O5B
2	D	401	NAD	O4B-C4B-C5B-O5B
2	A	401	NAD	C3D-C4D-C5D-O5D
2	C	401	NAD	C3B-C4B-C5B-O5B
2	D	401	NAD	C4N-C3N-C7N-O7N
2	D	401	NAD	C2N-C3N-C7N-O7N
2	D	401	NAD	C4N-C3N-C7N-N7N
2	A	401	NAD	C5B-O5B-PA-O1A
2	B	401	NAD	C5B-O5B-PA-O1A
2	D	401	NAD	C5B-O5B-PA-O1A
2	C	401	NAD	PN-O3-PA-O1A
2	D	401	NAD	C2N-C3N-C7N-N7N
2	C	401	NAD	C2D-C1D-N1N-C6N
2	D	401	NAD	C2D-C1D-N1N-C6N
2	B	401	NAD	PN-O3-PA-O2A

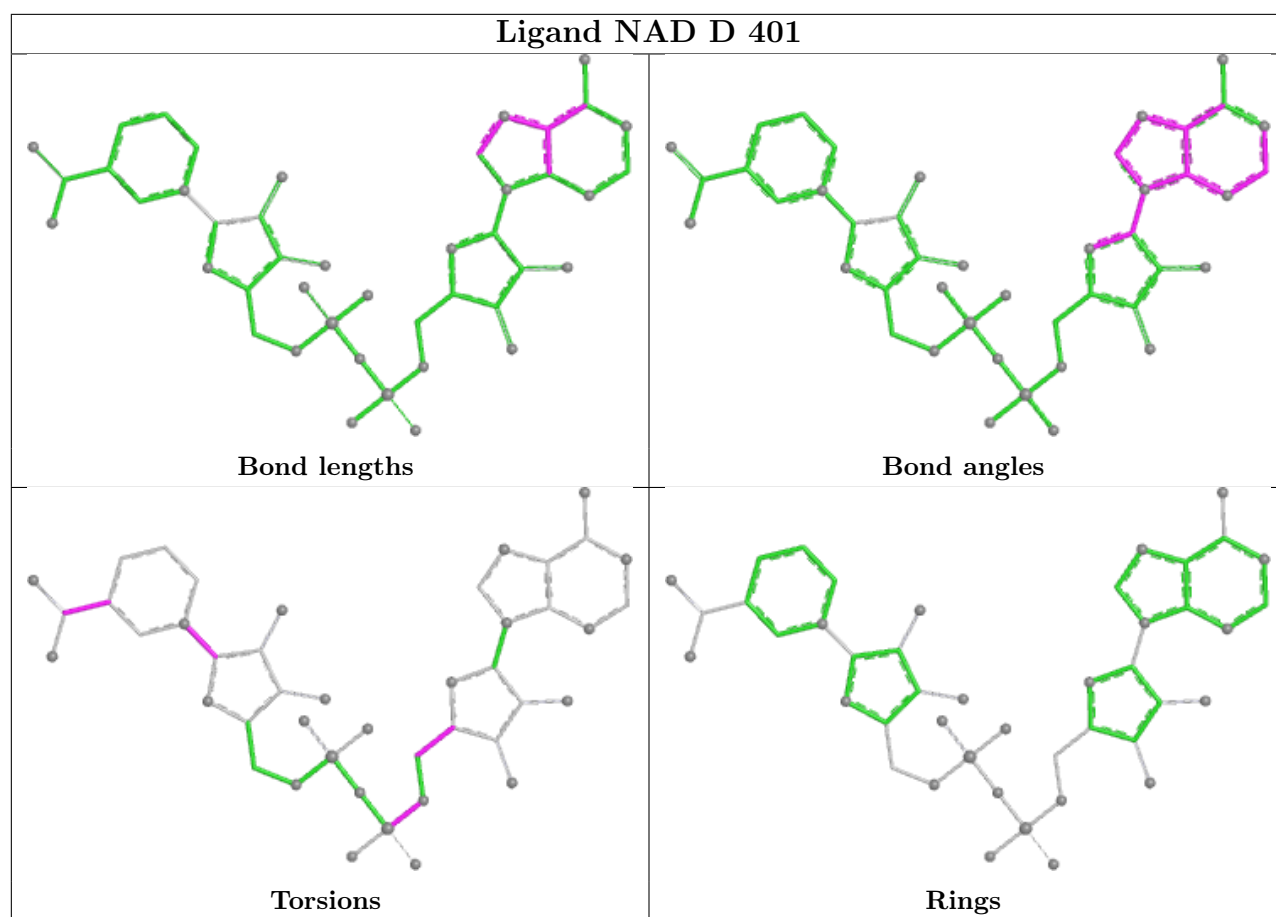
There are no ring outliers.

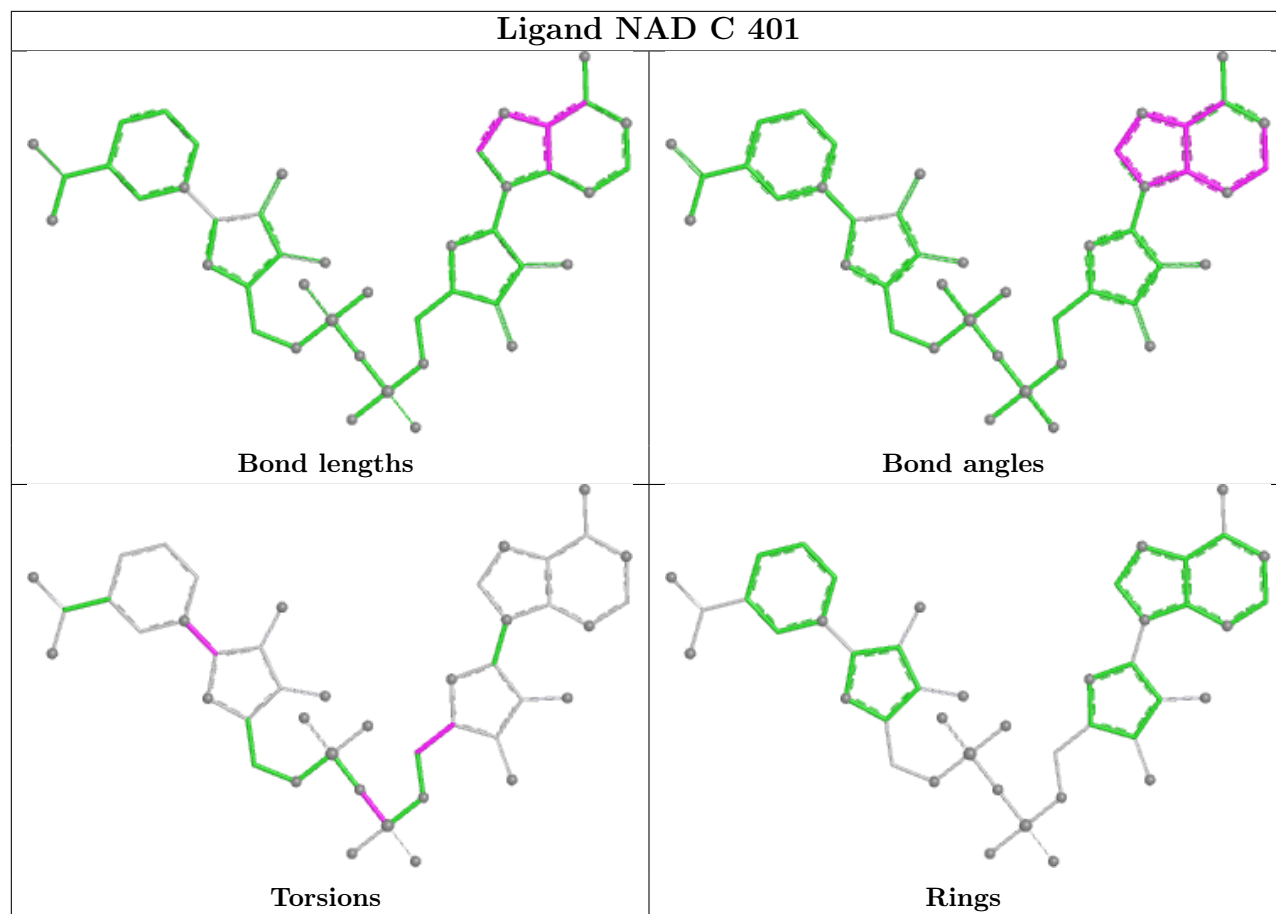
4 monomers are involved in 7 short contacts:

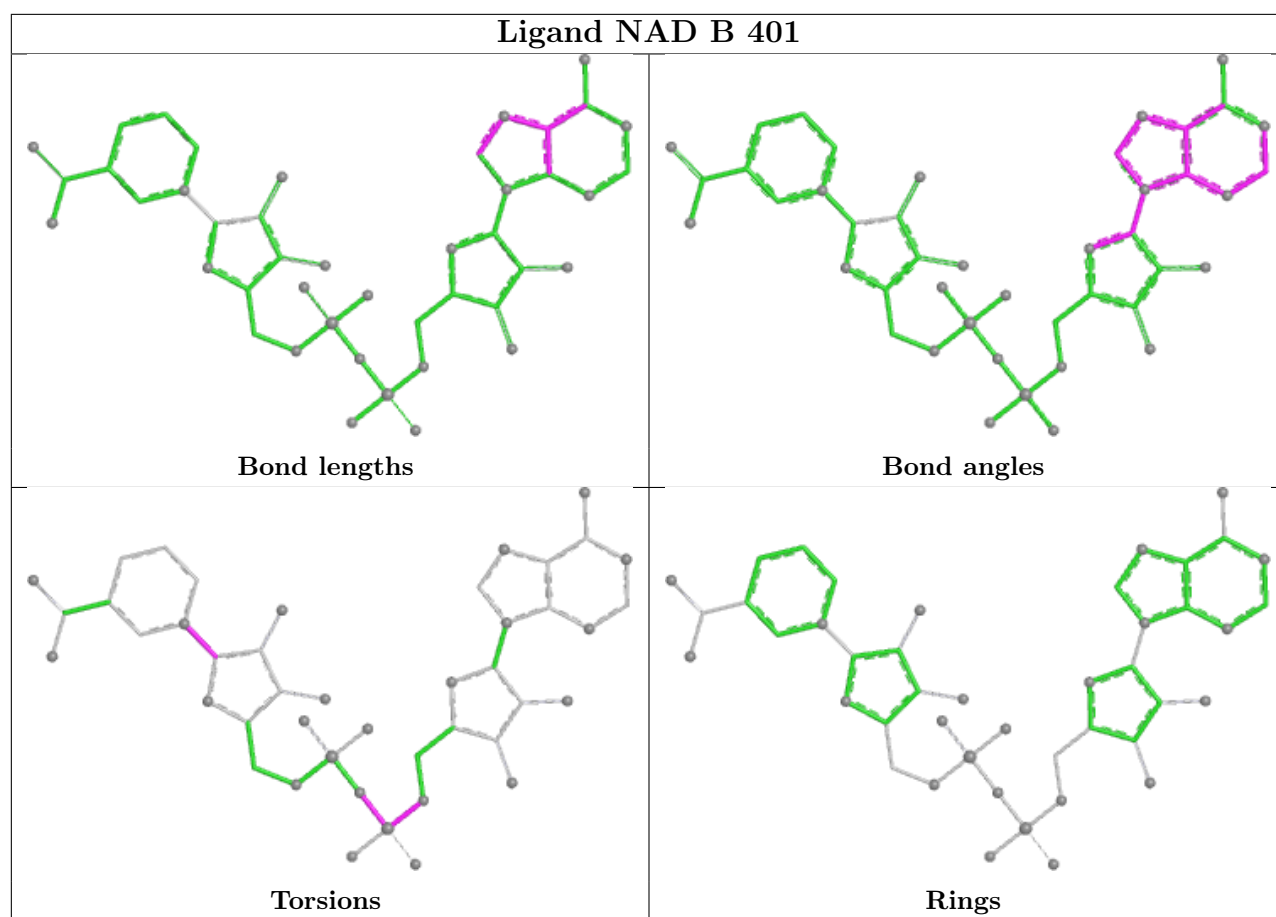
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	401	NAD	2	0
2	D	401	NAD	2	0
2	C	401	NAD	2	0
2	B	401	NAD	1	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.









## 5.7 Other polymers [i](#)

There are no such residues in this entry.

## 5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

## 6 Fit of model and data

### 6.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	336/336 (100%)	-0.26	3 (0%) 81 78	23, 35, 63, 132	0
1	B	336/336 (100%)	0.03	2 (0%) 85 83	29, 54, 85, 112	0
1	C	336/336 (100%)	0.25	6 (1%) 67 64	23, 62, 110, 135	0
1	D	336/336 (100%)	0.61	18 (5%) 31 27	35, 78, 120, 133	0
All	All	1344/1344 (100%)	0.16	29 (2%) 62 58	23, 55, 106, 135	0

All (29) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	98	CYS	3.7
1	D	109	CYS	3.5
1	D	113	PHE	3.1
1	D	96	GLY	2.9
1	D	81	TYR	2.8
1	C	79	ILE	2.8
1	D	79	ILE	2.8
1	D	112	ALA	2.7
1	C	105	ILE	2.7
1	D	116	TRP	2.7
1	D	18	LEU	2.7
1	D	74	VAL	2.5
1	A	1	MET	2.4
1	B	1	MET	2.4
1	C	103	GLY	2.4
1	C	114	GLY	2.3
1	D	10	PHE	2.3
1	D	61	PHE	2.3
1	B	106	ASP	2.3
1	C	176	LEU	2.2
1	D	31	ALA	2.2

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Mol	Chain	Res	Type	RSRZ
1	D	178	GLY	2.1
1	A	77	ASP	2.1
1	D	20	LEU	2.1
1	C	104	ALA	2.1
1	D	117	PHE	2.1
1	D	9	GLY	2.1
1	D	4	PRO	2.0
1	D	72	ALA	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 oligosaccharides in this entry.

## 6.4 Ligands [i](#)

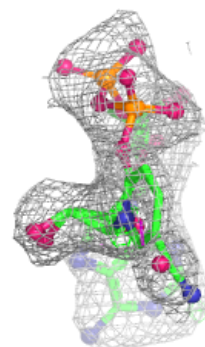
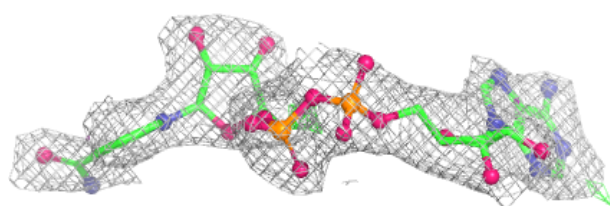
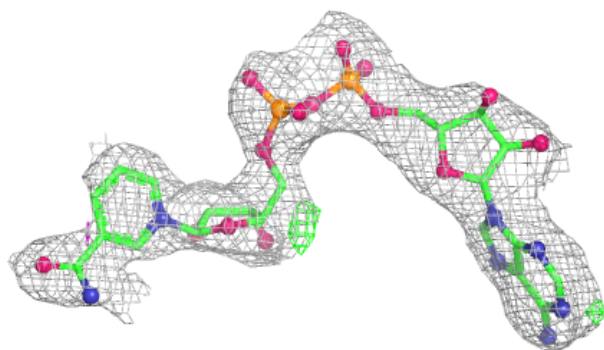
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
2	NAD	D	401	44/44	0.87	0.11	58,69,74,76	0
2	NAD	C	401	44/44	0.93	0.08	36,41,46,57	0
2	NAD	B	401	44/44	0.93	0.09	44,51,62,65	0
2	NAD	A	401	44/44	0.96	0.07	21,30,42,49	0
3	ZN	B	402	1/1	0.98	0.03	51,51,51,51	0
3	ZN	D	402	1/1	0.98	0.08	61,61,61,61	0
3	ZN	C	402	1/1	0.99	0.03	55,55,55,55	0
3	ZN	A	402	1/1	0.99	0.06	45,45,45,45	0

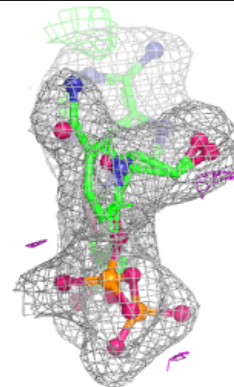
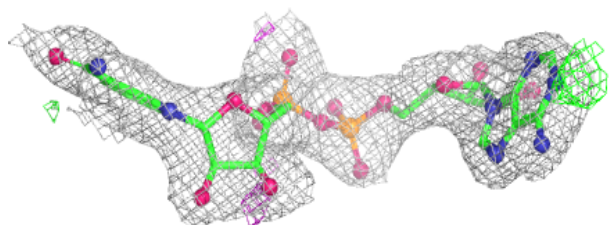
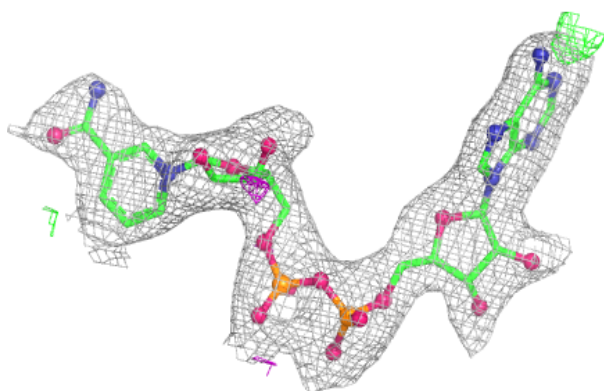
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

**Electron density around NAD D 401:**

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

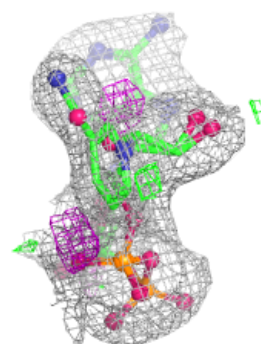
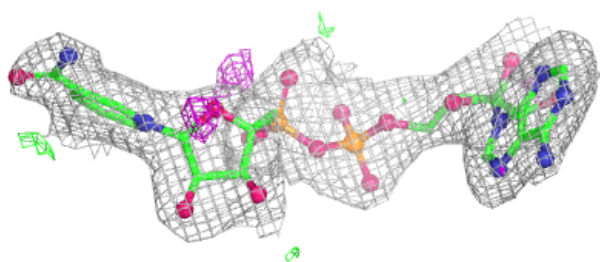
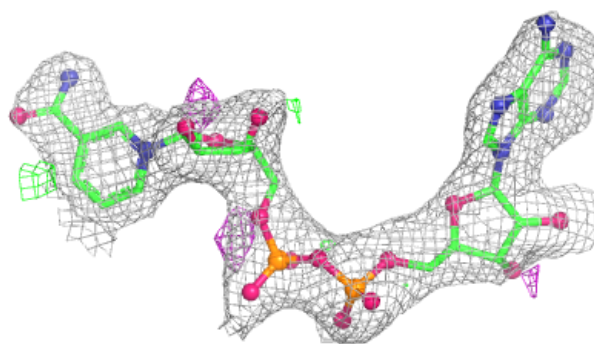
**Electron density around NAD C 401:**

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

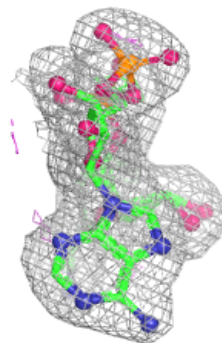
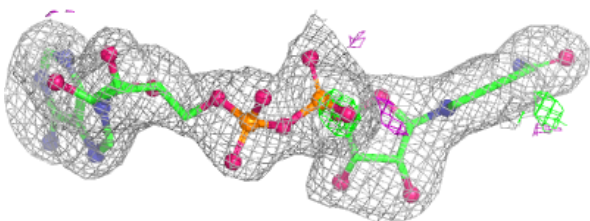
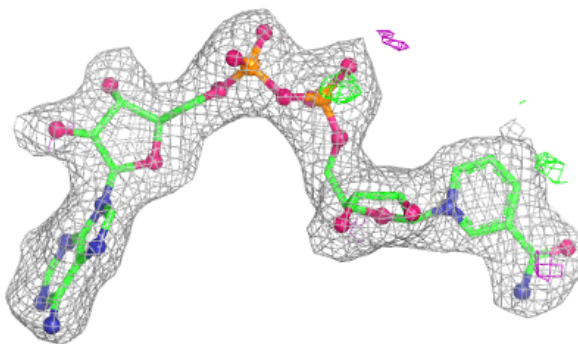


**Electron density around NAD B 401:**

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

**Electron density around NAD 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.