



# Full wwPDB X-ray Structure Validation Report ⓘ

Aug 8, 2022 – 11:43 AM EDT

PDB ID : 8DT1  
Title : Crystal Structure of a Putative D-beta-hydroxybutyrate dehydrogenase from Burkholderia cenocepacia J2315 in complex with NAD  
Authors : Seattle Structural Genomics Center for Infectious Disease (SSGCID)  
Deposited on : 2022-07-25  
Resolution : 1.80 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at [validation@mail.wwpdb.org](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>.

---

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)	:	1.13
EDS	:	2.29
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)	:	2.29

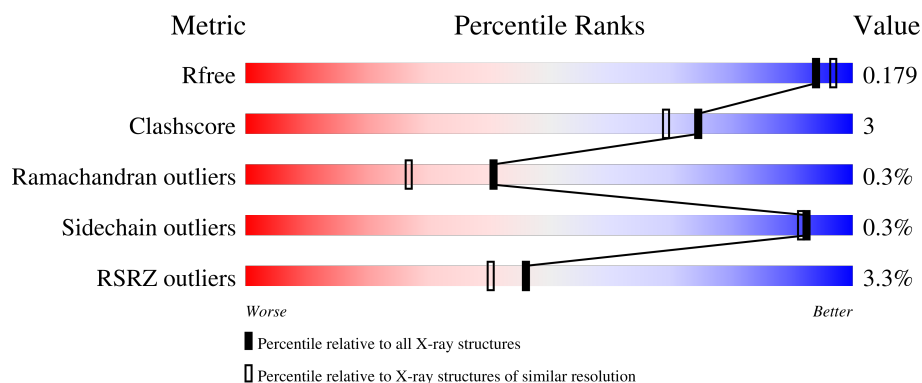
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 1.80 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
$R_{free}$	130704	5950 (1.80-1.80)
Clashscore	141614	6793 (1.80-1.80)
Ramachandran outliers	138981	6697 (1.80-1.80)
Sidechain outliers	138945	6696 (1.80-1.80)
RSRZ outliers	127900	5850 (1.80-1.80)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 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	269	<div> <div>5%</div> <div>87%</div> <div>6%</div> <div>7%</div> </div>
1	B	269	<div> <div>2%</div> <div>87%</div> <div>7%</div> <div>7%</div> </div>
1	C	269	<div> <div>%</div> <div>90%</div> <div>6%</div> <div>.</div> </div>
1	D	269	<div> <div>5%</div> <div>86%</div> <div>7%</div> <div>7%</div> </div>

## 2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 8605 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 3-hydroxybutyrate dehydrogenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	251	Total	C	N	O	S	0	7	0
			1847	1176	314	350	7			
1	B	251	Total	C	N	O	S	0	5	0
			1844	1172	316	349	7			
1	C	259	Total	C	N	O	S	0	4	0
			1889	1196	323	363	7			
1	D	251	Total	C	N	O	S	0	6	0
			1833	1165	313	348	7			

There are 32 discrepancies between the modelled and reference sequences:

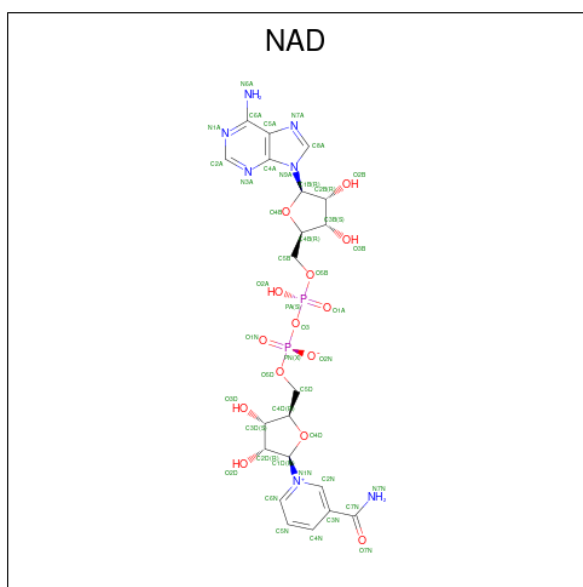
Chain	Residue	Modelled	Actual	Comment	Reference
A	-7	MET	-	initiating methionine	UNP A0A1V2Y0M0
A	-6	ALA	-	expression tag	UNP A0A1V2Y0M0
A	-5	HIS	-	expression tag	UNP A0A1V2Y0M0
A	-4	HIS	-	expression tag	UNP A0A1V2Y0M0
A	-3	HIS	-	expression tag	UNP A0A1V2Y0M0
A	-2	HIS	-	expression tag	UNP A0A1V2Y0M0
A	-1	HIS	-	expression tag	UNP A0A1V2Y0M0
A	0	HIS	-	expression tag	UNP A0A1V2Y0M0
B	-7	MET	-	initiating methionine	UNP A0A1V2Y0M0
B	-6	ALA	-	expression tag	UNP A0A1V2Y0M0
B	-5	HIS	-	expression tag	UNP A0A1V2Y0M0
B	-4	HIS	-	expression tag	UNP A0A1V2Y0M0
B	-3	HIS	-	expression tag	UNP A0A1V2Y0M0
B	-2	HIS	-	expression tag	UNP A0A1V2Y0M0
B	-1	HIS	-	expression tag	UNP A0A1V2Y0M0
B	0	HIS	-	expression tag	UNP A0A1V2Y0M0
C	-7	MET	-	initiating methionine	UNP A0A1V2Y0M0
C	-6	ALA	-	expression tag	UNP A0A1V2Y0M0
C	-5	HIS	-	expression tag	UNP A0A1V2Y0M0
C	-4	HIS	-	expression tag	UNP A0A1V2Y0M0
C	-3	HIS	-	expression tag	UNP A0A1V2Y0M0

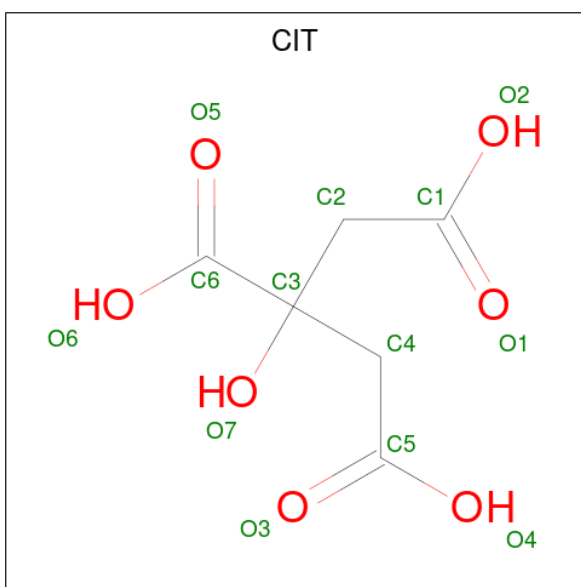
*Continued on next page...*

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
C	-2	HIS	-	expression tag	UNP A0A1V2Y0M0
C	-1	HIS	-	expression tag	UNP A0A1V2Y0M0
C	0	HIS	-	expression tag	UNP A0A1V2Y0M0
D	-7	MET	-	initiating methionine	UNP A0A1V2Y0M0
D	-6	ALA	-	expression tag	UNP A0A1V2Y0M0
D	-5	HIS	-	expression tag	UNP A0A1V2Y0M0
D	-4	HIS	-	expression tag	UNP A0A1V2Y0M0
D	-3	HIS	-	expression tag	UNP A0A1V2Y0M0
D	-2	HIS	-	expression tag	UNP A0A1V2Y0M0
D	-1	HIS	-	expression tag	UNP A0A1V2Y0M0
D	0	HIS	-	expression tag	UNP A0A1V2Y0M0

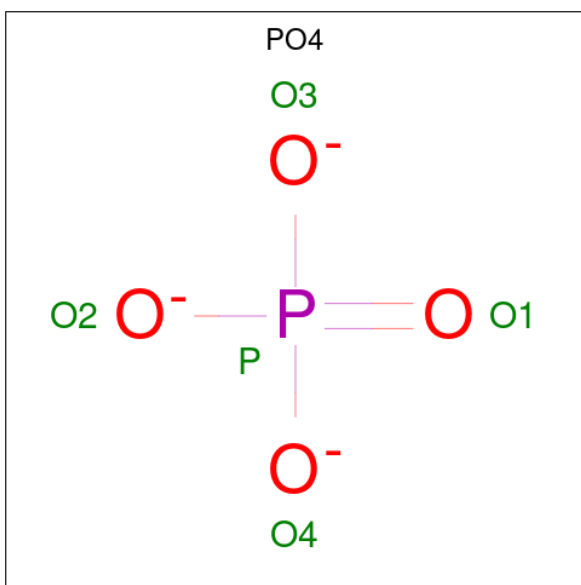
- Molecule 2 is NICOTINAMIDE-ADENINE-DINUCLEOTIDE (three-letter code: NAD) (formula:  $C_{21}H_{27}N_7O_{14}P_2$ ) (labeled as "Ligand of Interest" by depositor).





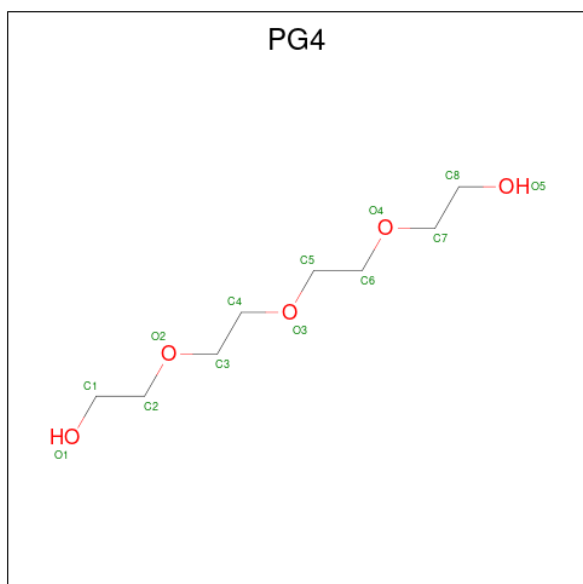
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			13	6	7		
3	B	1	Total	C	O	0	0
			13	6	7		
3	C	1	Total	C	O	0	0
			13	6	7		
3	D	1	Total	C	O	0	0
			13	6	7		

- Molecule 4 is PHOSPHATE ION (three-letter code: PO4) (formula: O<sub>4</sub>P).



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

- Molecule 5 is TETRAETHYLENE GLYCOL (three-letter code: PG4) (formula: C<sub>8</sub>H<sub>18</sub>O<sub>5</sub>).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			10	6	4		
5	A	1	Total	C	O	0	0
			7	4	3		
5	B	1	Total	C	O	0	0
			10	6	4		
5	B	1	Total	C	O	0	0
			7	4	3		
5	C	1	Total	C	O	0	0
			10	6	4		
5	C	1	Total	C	O	0	0
			7	4	3		
5	D	1	Total	C	O	0	0
			10	6	4		
5	D	1	Total	C	O	0	0
			7	4	3		

- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	202	Total	O	0	4
			206	206		

*Continued on next page...*

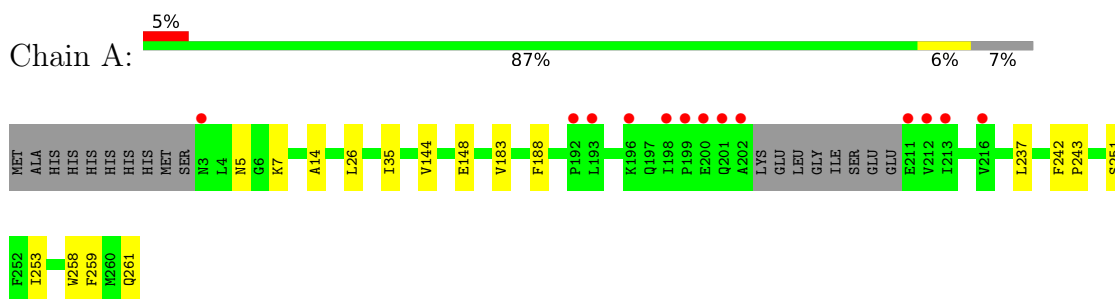
*Continued from previous page...*

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	B	207	Total 216	O 216	0	9
6	C	153	Total 158	O 158	0	5
6	D	176	Total 179	O 179	0	4

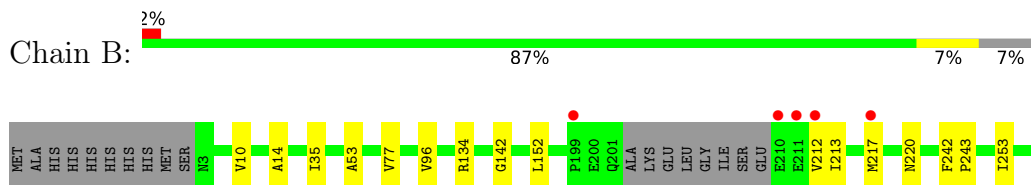
### 3 Residue-property plots [i](#)

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.

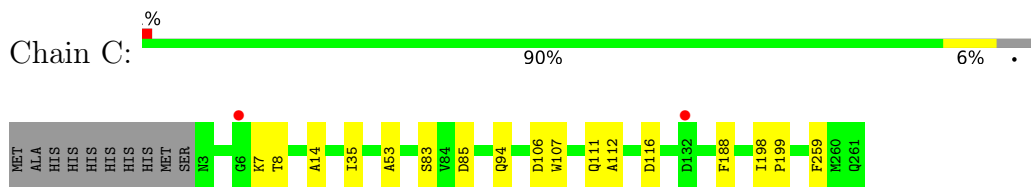
- Molecule 1: 3-hydroxybutyrate dehydrogenase



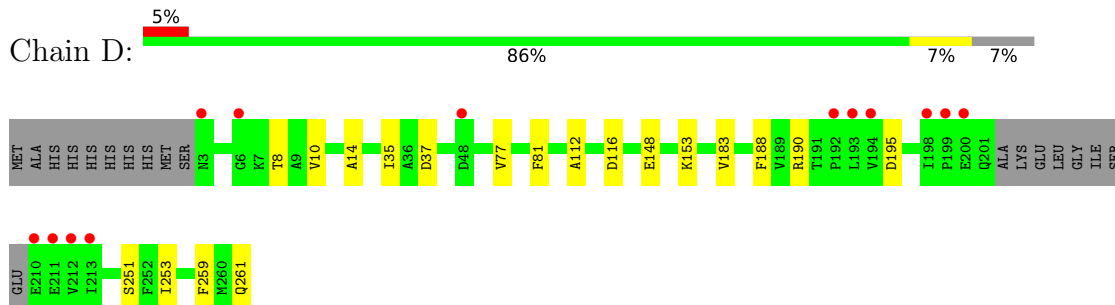
- Molecule 1: 3-hydroxybutyrate dehydrogenase



- Molecule 1: 3-hydroxybutyrate dehydrogenase



- Molecule 1: 3-hydroxybutyrate dehydrogenase





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	59.91Å 65.60Å 68.32Å 97.15° 98.38° 107.47°	Depositor
Resolution (Å)	22.17 – 1.80 49.83 – 1.80	Depositor EDS
% Data completeness (in resolution range)	99.4 (22.17-1.80) 99.4 (49.83-1.80)	Depositor EDS
$R_{merge}$	0.04	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.34 (at 1.79Å)	Xtriage
Refinement program	PHENIX 1.20.1	Depositor
R, $R_{free}$	0.146 , 0.178 0.148 , 0.179	Depositor DCC
$R_{free}$ test set	2013 reflections (2.26%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	19.5	Xtriage
Anisotropy	0.523	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.35 , 55.0	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.97	EDS
Total number of atoms	8605	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	26.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.67% 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: NAD, CIT, PG4, PO4

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with  $|Z| > 5$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z  > 5$	RMSZ	$\# Z  > 5$
1	A	0.43	0/1897	0.59	0/2580
1	B	0.46	0/1888	0.60	0/2564
1	C	0.39	0/1931	0.57	0/2624
1	D	0.39	0/1880	0.57	0/2557
All	All	0.42	0/7596	0.58	0/10325

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

### 5.2 Too-close contacts [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	1847	0	1851	11	0
1	B	1844	0	1851	11	0
1	C	1889	0	1877	14	0
1	D	1833	0	1817	13	0
2	A	88	0	52	0	0
2	B	44	0	26	1	0
2	C	88	0	52	0	0
2	D	88	0	52	1	0
3	A	13	0	5	0	0

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	13	0	5	0	0
3	C	13	0	5	4	0
3	D	13	0	5	2	0
4	A	5	0	0	0	0
5	A	17	0	22	0	0
5	B	17	0	22	1	0
5	C	17	0	22	2	0
5	D	17	0	22	1	0
6	A	206	0	0	3	0
6	B	216	0	0	1	0
6	C	158	0	0	2	0
6	D	179	0	0	1	0
All	All	8605	0	7686	50	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 (50) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:134:ARG:NH2	6:B:401:HOH:O	2.06	0.86
1:C:106:ASP:OD2	6:C:401:HOH:O	2.03	0.76
1:B:14:ALA:HB3	1:B:35:ILE:HG23	1.75	0.69
1:C:8[B]:THR:HG23	1:C:85:ASP:H	1.60	0.67
1:D:14:ALA:HB3	1:D:35[A]:ILE:HG23	1.80	0.63
1:B:96[B]:VAL:HG13	1:B:152:LEU:HB3	1.83	0.61
1:A:14:ALA:HB3	1:A:35[A]:ILE:HG23	1.82	0.61
1:A:5:ASN:ND2	6:A:405:HOH:O	2.35	0.59
1:A:7:LYS:NZ	6:A:404:HOH:O	2.33	0.59
1:A:148:GLU:OE1	6:A:401:HOH:O	2.16	0.59
1:B:213:ILE:HA	1:B:217:MET:HE2	1.84	0.58
1:D:153:LYS:NZ	3:D:302:CIT:H42	2.18	0.58
1:C:94:GLN:HE21	3:C:302:CIT:C6	2.17	0.57
1:B:220:ASN:HB3	1:B:258:TRP:CD1	2.40	0.57
1:C:14:ALA:HB3	1:C:35:ILE:HG23	1.85	0.57
1:C:8[B]:THR:HG21	1:C:83:SER:O	2.05	0.56
1:D:14:ALA:HB3	1:D:35[B]:ILE:HG23	1.87	0.55
1:B:10[A]:VAL:HG22	1:B:77:VAL:HG11	1.90	0.54
1:D:10[B]:VAL:HG22	1:D:77:VAL:HG11	1.89	0.54
1:C:53:ALA:HB3	5:C:304:PG4:H21	1.90	0.53
1:C:8[B]:THR:HG21	1:C:83:SER:C	2.30	0.52

*Continued on next page...*

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:14:ALA:HB3	1:A:35[B]:ILE:HG23	1.91	0.52
1:C:188:PHE:CE1	3:C:302:CIT:H42	2.44	0.51
6:C:489[A]:HOH:O	1:D:253[A]:ILE:HG12	2.10	0.51
1:D:190:ARG:NE	1:D:195:ASP:OD1	2.41	0.49
1:D:253[B]:ILE:HD12	1:D:261:GLN:HG2	1.94	0.48
1:C:198:ILE:HB	1:C:199:PRO:HD3	1.96	0.48
1:B:53:ALA:HB3	5:B:304:PG4:H31	1.96	0.48
1:C:7:LYS:HE2	5:C:303:PG4:H12	1.96	0.47
1:D:8:THR:HG21	1:D:81:PHE:HB3	1.97	0.47
5:D:303:PG4:H21	5:D:303:PG4:H41	1.60	0.45
1:A:26:LEU:HD23	1:A:237[A]:LEU:HD23	2.00	0.44
1:D:148:GLU:OE1	6:D:401:HOH:O	2.20	0.44
1:B:142:GLY:O	2:B:301:NAD:H6N	2.18	0.44
1:D:37:ASP:OD1	2:D:301[A]:NAD:O2B	2.29	0.44
1:B:253:ILE:CD1	1:B:259:PHE:HD2	2.30	0.44
1:A:242:PHE:HA	1:A:243:PRO:HD3	1.91	0.43
1:D:153:LYS:HZ3	3:D:302:CIT:H42	1.83	0.43
1:A:144:VAL:HG11	1:A:258:TRP:CD2	2.55	0.42
1:C:188:PHE:CD1	3:C:302:CIT:H42	2.55	0.42
1:C:188:PHE:HE1	3:C:302:CIT:C5	2.33	0.42
1:D:112:ALA:O	1:D:116:ASP:HB2	2.20	0.42
1:A:144:VAL:HG11	1:A:258:TRP:CE2	2.55	0.42
1:D:183:VAL:O	1:D:251:SER:HA	2.20	0.42
1:B:212:VAL:O	1:B:217:MET:HG3	2.20	0.41
1:C:107:TRP:O	1:C:111[B]:GLN:HG2	2.20	0.41
1:A:253[A]:ILE:HD12	1:A:261:GLN:HG2	2.03	0.41
1:C:112:ALA:O	1:C:116:ASP:HB2	2.21	0.40
1:A:183:VAL:O	1:A:251:SER:HA	2.21	0.40
1:B:242:PHE:HA	1:B:243:PRO:HD3	1.89	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	254/269 (94%)	248 (98%)	5 (2%)	1 (0%)	34	21
1	B	252/269 (94%)	244 (97%)	8 (3%)	0	100	100
1	C	261/269 (97%)	257 (98%)	3 (1%)	1 (0%)	34	21
1	D	253/269 (94%)	247 (98%)	5 (2%)	1 (0%)	34	21
All	All	1020/1076 (95%)	996 (98%)	21 (2%)	3 (0%)	41	27

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	259	PHE
1	D	259	PHE
1	C	259	PHE

### 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	189/207 (91%)	188 (100%)	1 (0%)	88	87
1	B	189/207 (91%)	189 (100%)	0	100	100
1	C	192/207 (93%)	192 (100%)	0	100	100
1	D	185/207 (89%)	184 (100%)	1 (0%)	88	87
All	All	755/828 (91%)	753 (100%)	2 (0%)	92	91

All (2) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	188	PHE
1	D	188	PHE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	145	HIS

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type
1	C	145	HIS
1	D	145	HIS

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

## 5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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

## 5.5 Carbohydrates ⓘ

There are no monosaccharides in this entry.

## 5.6 Ligand geometry ⓘ

20 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	CIT	D	302	-	12,12,12	1.08	0	17,17,17	1.61	3 (17%)
2	NAD	D	301[B]	-	42,48,48	0.56	0	50,73,73	0.62	1 (2%)
5	PG4	D	303	-	9,9,12	0.13	0	8,8,11	0.55	0
5	PG4	B	304	-	6,6,12	0.11	0	5,5,11	0.16	0
5	PG4	C	304	-	6,6,12	0.14	0	5,5,11	0.16	0
2	NAD	D	301[A]	-	42,48,48	0.56	0	50,73,73	0.66	1 (2%)
2	NAD	A	301[B]	-	42,48,48	0.55	0	50,73,73	0.63	1 (2%)
3	CIT	C	302	-	12,12,12	1.48	1 (8%)	17,17,17	1.78	4 (23%)
5	PG4	D	304	-	6,6,12	0.18	0	5,5,11	0.13	0
3	CIT	B	302	-	12,12,12	1.12	0	17,17,17	1.94	6 (35%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
2	NAD	A	301[A]	-	42,48,48	0.58	0	50,73,73	0.71	1 (2%)
4	PO4	A	303	-	4,4,4	0.91	0	6,6,6	1.06	1 (16%)
5	PG4	A	305	-	6,6,12	0.15	0	5,5,11	0.16	0
2	NAD	C	301[B]	-	42,48,48	0.56	0	50,73,73	0.64	1 (2%)
2	NAD	B	301	-	42,48,48	0.61	0	50,73,73	0.73	1 (2%)
2	NAD	C	301[A]	-	42,48,48	0.57	0	50,73,73	0.66	1 (2%)
5	PG4	A	304	-	9,9,12	0.16	0	8,8,11	0.50	0
5	PG4	B	303	-	9,9,12	0.14	0	8,8,11	0.56	0
3	CIT	A	302	-	12,12,12	1.10	0	17,17,17	1.48	1 (5%)
5	PG4	C	303	-	9,9,12	0.16	0	8,8,11	0.55	0

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	CIT	D	302	-	-	10/16/16/16	-
2	NAD	D	301[B]	-	-	10/26/62/62	0/5/5/5
5	PG4	D	303	-	-	5/7/7/10	-
5	PG4	B	304	-	-	3/4/4/10	-
5	PG4	C	304	-	-	1/4/4/10	-
2	NAD	D	301[A]	-	-	6/26/62/62	0/5/5/5
2	NAD	A	301[B]	-	-	10/26/62/62	0/5/5/5
3	CIT	C	302	-	-	8/16/16/16	-
5	PG4	D	304	-	-	1/4/4/10	-
3	CIT	B	302	-	-	7/16/16/16	-
2	NAD	A	301[A]	-	-	6/26/62/62	0/5/5/5
5	PG4	A	305	-	-	1/4/4/10	-
2	NAD	C	301[B]	-	-	4/26/62/62	0/5/5/5
2	NAD	B	301	-	-	7/26/62/62	0/5/5/5
2	NAD	C	301[A]	-	-	7/26/62/62	0/5/5/5
5	PG4	A	304	-	-	5/7/7/10	-
5	PG4	B	303	-	-	3/7/7/10	-
3	CIT	A	302	-	-	8/16/16/16	-
5	PG4	C	303	-	-	2/7/7/10	-

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	302	CIT	C3-C6	-3.19	1.50	1.53

All (22) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	302	CIT	O6-C6-C3	4.97	121.67	113.05
3	A	302	CIT	O6-C6-C3	4.04	120.06	113.05
3	C	302	CIT	O7-C3-C4	3.79	118.27	109.40
3	C	302	CIT	O6-C6-C3	3.14	118.50	113.05
3	D	302	CIT	O6-C6-C3	3.01	118.28	113.05
3	B	302	CIT	O7-C3-C6	2.94	113.00	108.86
3	D	302	CIT	O7-C3-C6	-2.92	104.76	108.86
3	B	302	CIT	C2-C3-C6	-2.57	104.59	110.11
2	C	301[B]	NAD	C5A-C6A-N6A	2.41	124.01	120.35
2	C	301[A]	NAD	C5A-C6A-N6A	2.35	123.92	120.35
3	B	302	CIT	O4-C5-C4	2.30	121.73	114.35
2	D	301[B]	NAD	C5A-C6A-N6A	2.29	123.83	120.35
2	B	301	NAD	C5A-C6A-N6A	2.27	123.80	120.35
2	D	301[A]	NAD	C5A-C6A-N6A	2.27	123.80	120.35
2	A	301[B]	NAD	C5A-C6A-N6A	2.23	123.74	120.35
3	C	302	CIT	C3-C2-C1	-2.23	108.43	113.81
3	B	302	CIT	O5-C6-C3	-2.20	119.14	122.25
2	A	301[A]	NAD	C5A-C6A-N6A	2.20	123.69	120.35
4	A	303	PO4	O4-P-O3	2.06	114.58	107.97
3	C	302	CIT	O4-C5-C4	2.02	120.84	114.35
3	B	302	CIT	O4-C5-O3	-2.02	118.27	123.30
3	D	302	CIT	O4-C5-C4	2.01	120.80	114.35

There are no chirality outliers.

All (104) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	301[A]	NAD	O4D-C1D-N1N-C2N
2	A	301[B]	NAD	C5B-O5B-PA-O3
2	A	301[B]	NAD	C5D-O5D-PN-O2N
2	A	301[B]	NAD	O4D-C1D-N1N-C2N
2	B	301	NAD	C5D-O5D-PN-O1N
2	B	301	NAD	C5D-O5D-PN-O2N
2	B	301	NAD	O4D-C1D-N1N-C2N
2	C	301[A]	NAD	C5D-O5D-PN-O1N
2	C	301[A]	NAD	C5D-O5D-PN-O2N
2	C	301[A]	NAD	O4D-C1D-N1N-C2N

*Continued on next page...*



*Continued from previous page...*

Mol	Chain	Res	Type	Atoms
2	C	301[B]	NAD	O4D-C1D-N1N-C2N
2	C	301[B]	NAD	O4D-C1D-N1N-C6N
2	D	301[A]	NAD	C5D-O5D-PN-O1N
2	D	301[A]	NAD	C5D-O5D-PN-O2N
2	D	301[A]	NAD	O4D-C1D-N1N-C2N
2	D	301[B]	NAD	C5B-O5B-PA-O3
2	D	301[B]	NAD	C5D-O5D-PN-O2N
2	D	301[B]	NAD	O4D-C1D-N1N-C2N
3	A	302	CIT	C1-C2-C3-O7
3	A	302	CIT	C1-C2-C3-C4
3	A	302	CIT	C1-C2-C3-C6
3	A	302	CIT	C2-C3-C6-O5
3	A	302	CIT	C2-C3-C6-O6
3	A	302	CIT	O7-C3-C6-O5
3	A	302	CIT	O7-C3-C6-O6
3	C	302	CIT	C1-C2-C3-C4
3	C	302	CIT	C1-C2-C3-C6
3	C	302	CIT	C2-C3-C6-O5
3	C	302	CIT	C2-C3-C6-O6
3	C	302	CIT	O7-C3-C6-O5
3	C	302	CIT	O7-C3-C6-O6
3	D	302	CIT	C2-C3-C6-O6
3	D	302	CIT	O7-C3-C6-O5
3	D	302	CIT	O7-C3-C6-O6
5	D	303	PG4	C4-C3-O2-C2
5	D	303	PG4	O2-C3-C4-O3
5	B	303	PG4	O2-C3-C4-O3
5	B	304	PG4	O1-C1-C2-O2
5	A	304	PG4	O2-C3-C4-O3
3	C	302	CIT	C1-C2-C3-O7
3	D	302	CIT	C1-C2-C3-C6
5	A	305	PG4	O1-C1-C2-O2
5	B	303	PG4	O3-C5-C6-O4
5	D	303	PG4	O1-C1-C2-O2
5	D	304	PG4	O1-C1-C2-O2
5	A	304	PG4	O3-C5-C6-O4
3	D	302	CIT	C1-C2-C3-C4
5	C	303	PG4	O1-C1-C2-O2
3	D	302	CIT	C4-C3-C6-O5
3	D	302	CIT	C4-C3-C6-O6
3	B	302	CIT	C2-C3-C6-O5
3	B	302	CIT	C2-C3-C6-O6

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type	Atoms
3	B	302	CIT	C4-C3-C6-O5
3	B	302	CIT	C4-C3-C6-O6
3	D	302	CIT	C2-C3-C6-O5
2	A	301[B]	NAD	PA-O3-PN-O5D
2	D	301[B]	NAD	PA-O3-PN-O5D
2	A	301[A]	NAD	C5D-O5D-PN-O3
2	C	301[A]	NAD	C5D-O5D-PN-O3
2	D	301[A]	NAD	C5D-O5D-PN-O3
2	D	301[B]	NAD	C5D-O5D-PN-O3
5	B	304	PG4	C1-C2-O2-C3
2	A	301[B]	NAD	PN-O3-PA-O1A
2	A	301[A]	NAD	C5D-O5D-PN-O2N
2	A	301[B]	NAD	C5B-O5B-PA-O1A
2	A	301[B]	NAD	C5B-O5B-PA-O2A
2	D	301[B]	NAD	C5B-O5B-PA-O2A
2	D	301[B]	NAD	C5D-O5D-PN-O1N
5	C	304	PG4	O1-C1-C2-O2
5	A	304	PG4	C3-C4-O3-C5
5	D	303	PG4	C3-C4-O3-C5
5	D	303	PG4	O3-C5-C6-O4
5	A	304	PG4	C4-C3-O2-C2
2	D	301[A]	NAD	O4B-C4B-C5B-O5B
2	B	301	NAD	PA-O3-PN-O2N
2	D	301[B]	NAD	PN-O3-PA-O2A
3	D	302	CIT	O7-C3-C4-C5
3	B	302	CIT	C3-C4-C5-O4
3	B	302	CIT	O7-C3-C6-O6
3	C	302	CIT	C4-C3-C6-O5
3	B	302	CIT	C3-C4-C5-O3
5	B	303	PG4	C6-C5-O3-C4
2	C	301[A]	NAD	O4B-C4B-C5B-O5B
2	A	301[A]	NAD	PN-O3-PA-O1A
2	D	301[B]	NAD	PA-O3-PN-O2N
3	A	302	CIT	C4-C3-C6-O6
2	A	301[B]	NAD	C5D-O5D-PN-O3
2	B	301	NAD	C5D-O5D-PN-O3
2	C	301[B]	NAD	C2D-C1D-N1N-C2N
2	C	301[B]	NAD	C2D-C1D-N1N-C6N
5	B	304	PG4	O2-C3-C4-O3
2	A	301[B]	NAD	PA-O3-PN-O2N
2	B	301	NAD	PA-O3-PN-O1N
2	C	301[A]	NAD	PA-O3-PN-O1N

*Continued on next page...*

*Continued from previous page...*

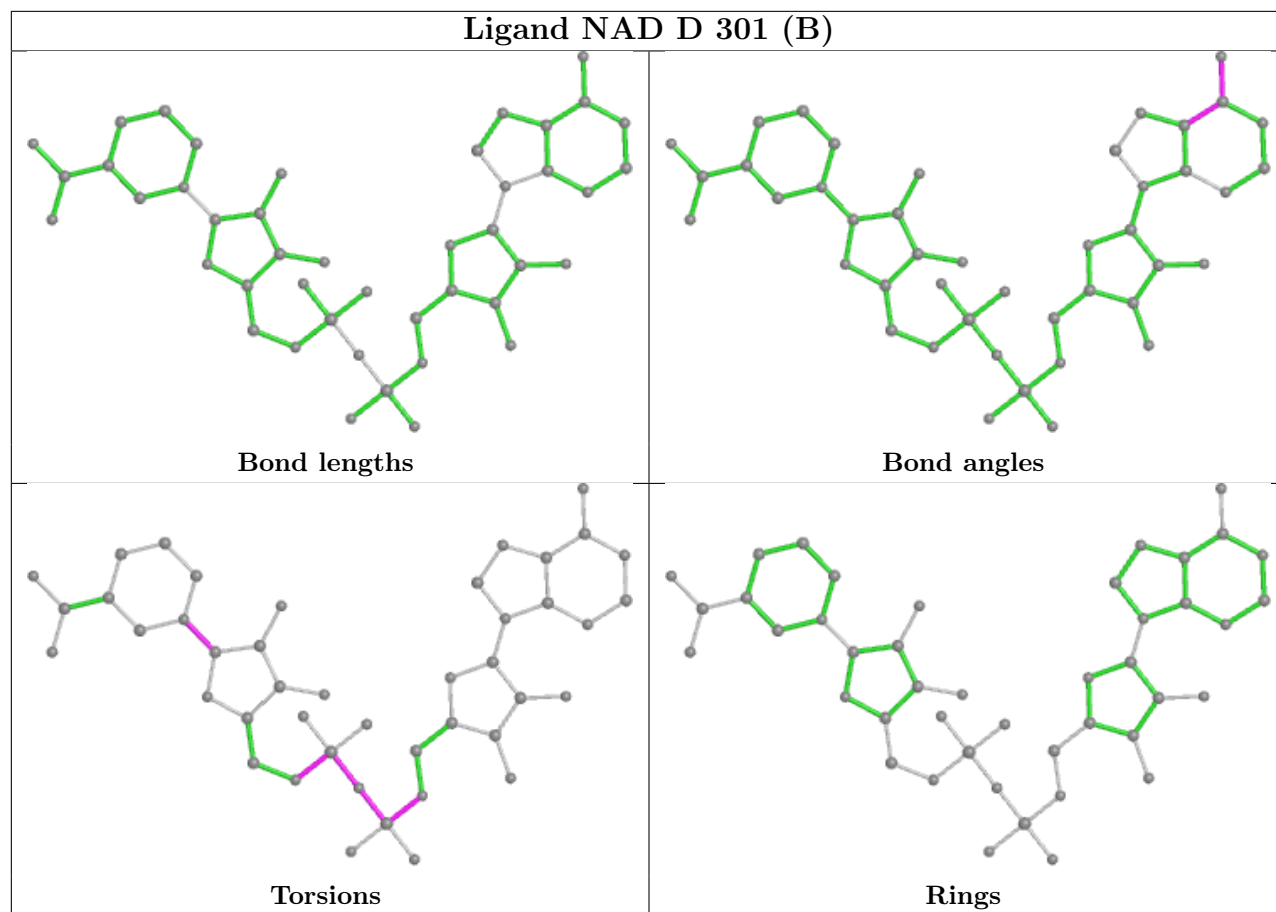
Mol	Chain	Res	Type	Atoms
2	C	301[A]	NAD	PA-O3-PN-O2N
2	D	301[A]	NAD	PA-O3-PN-O2N
5	C	303	PG4	O3-C5-C6-O4
3	D	302	CIT	C1-C2-C3-O7
2	A	301[A]	NAD	C5D-O5D-PN-O1N
2	A	301[B]	NAD	C5D-O5D-PN-O1N
2	D	301[B]	NAD	C5B-O5B-PA-O1A
2	A	301[A]	NAD	O4B-C4B-C5B-O5B
2	B	301	NAD	O4B-C4B-C5B-O5B
5	A	304	PG4	O1-C1-C2-O2

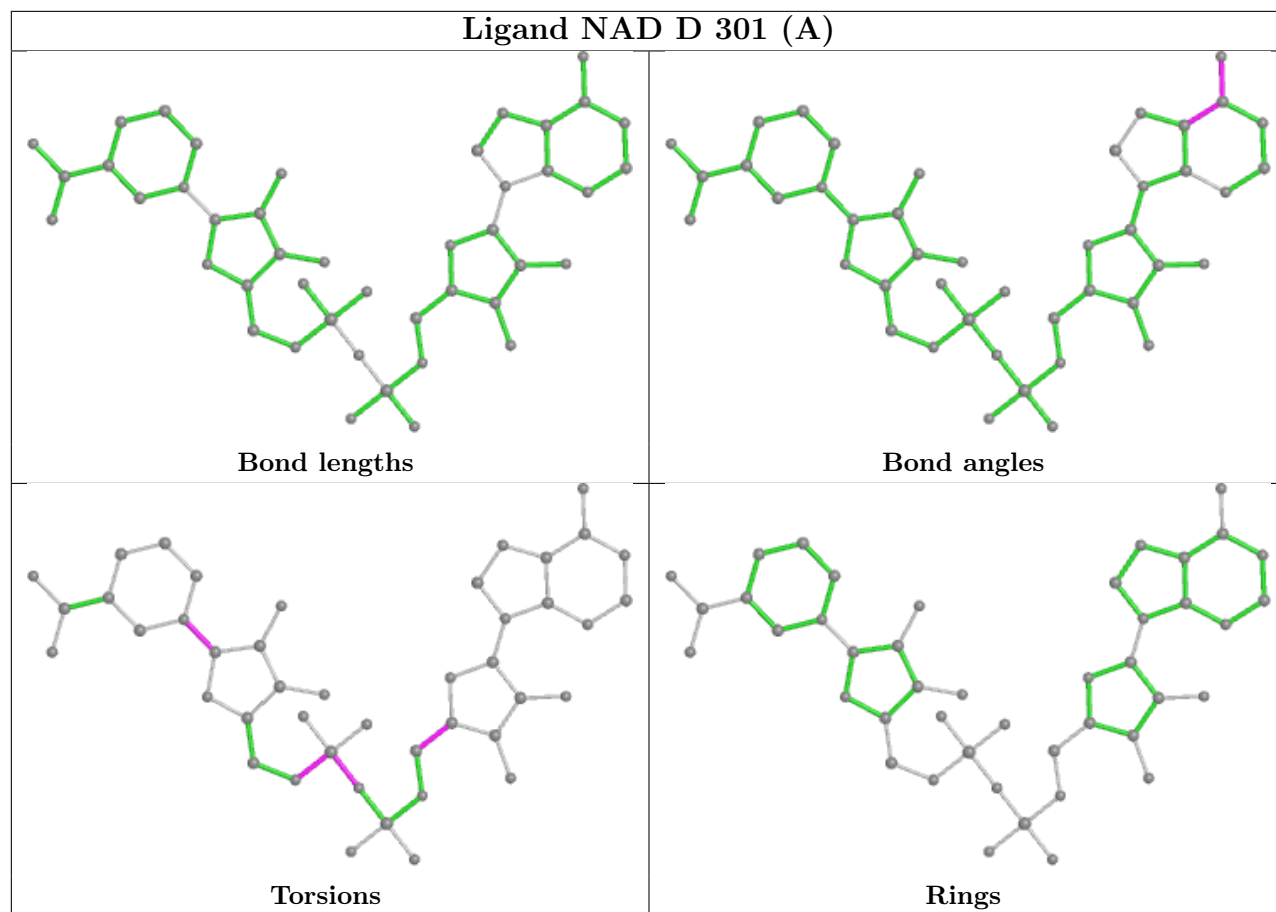
There are no ring outliers.

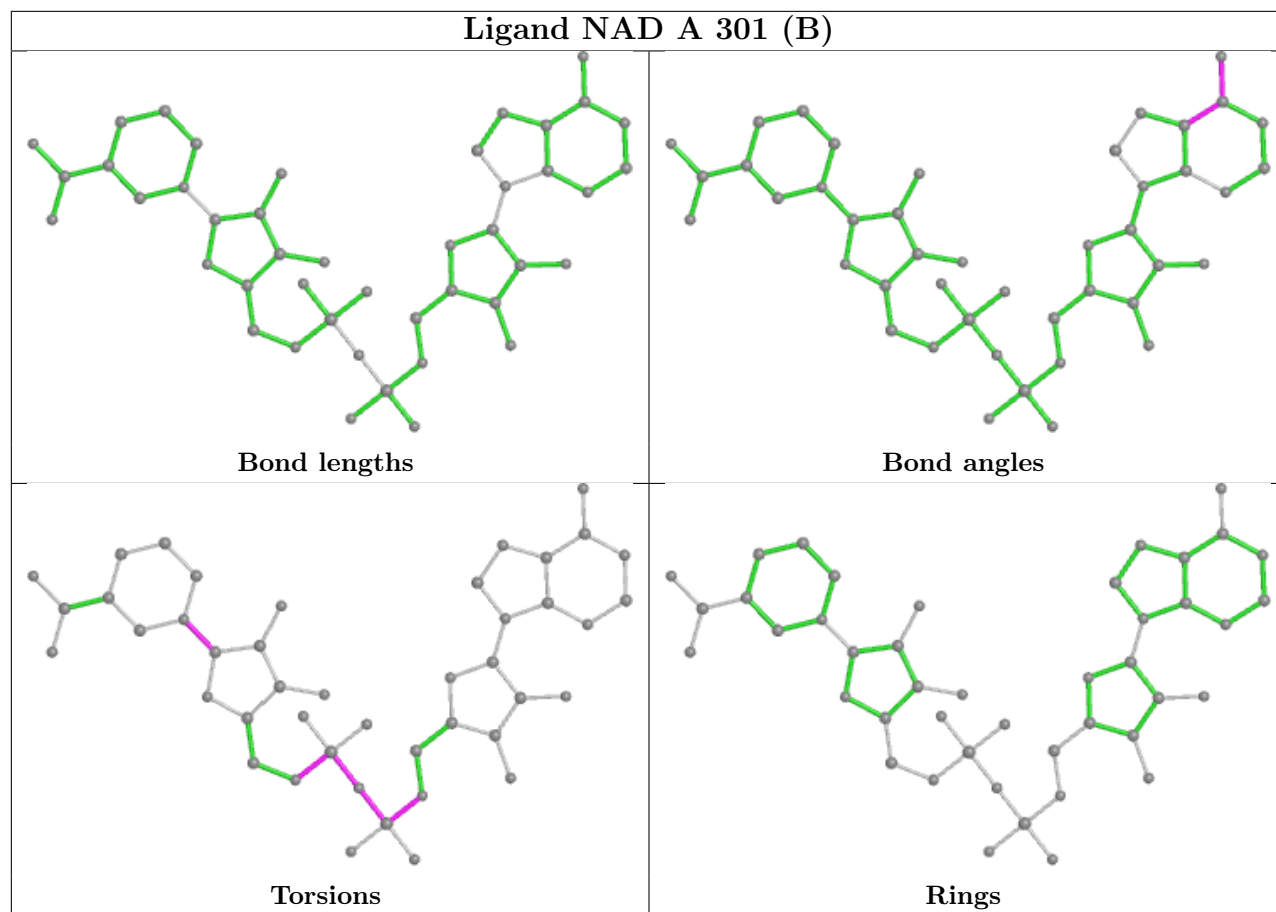
8 monomers are involved in 12 short contacts:

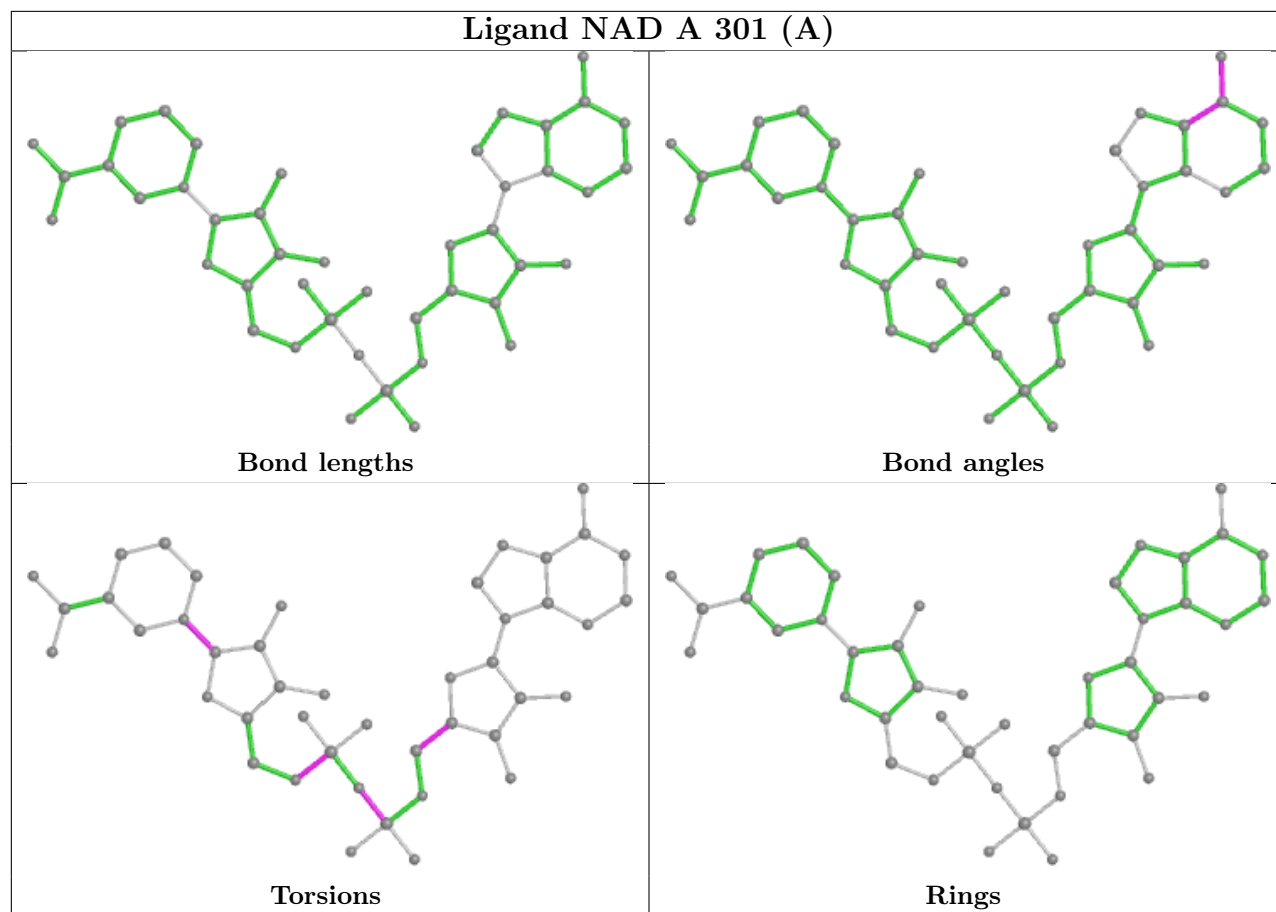
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	D	302	CIT	2	0
5	D	303	PG4	1	0
5	B	304	PG4	1	0
5	C	304	PG4	1	0
2	D	301[A]	NAD	1	0
3	C	302	CIT	4	0
2	B	301	NAD	1	0
5	C	303	PG4	1	0

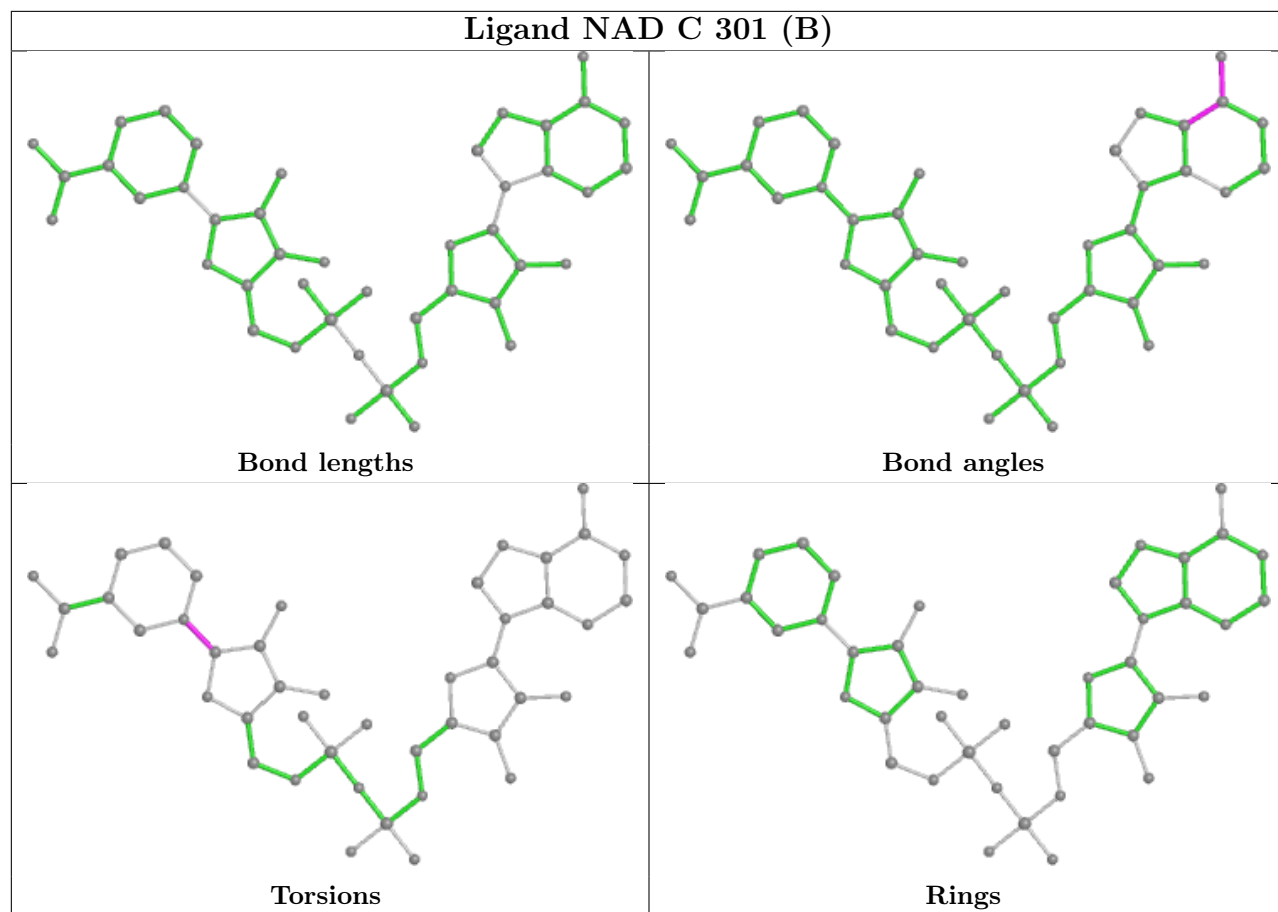
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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



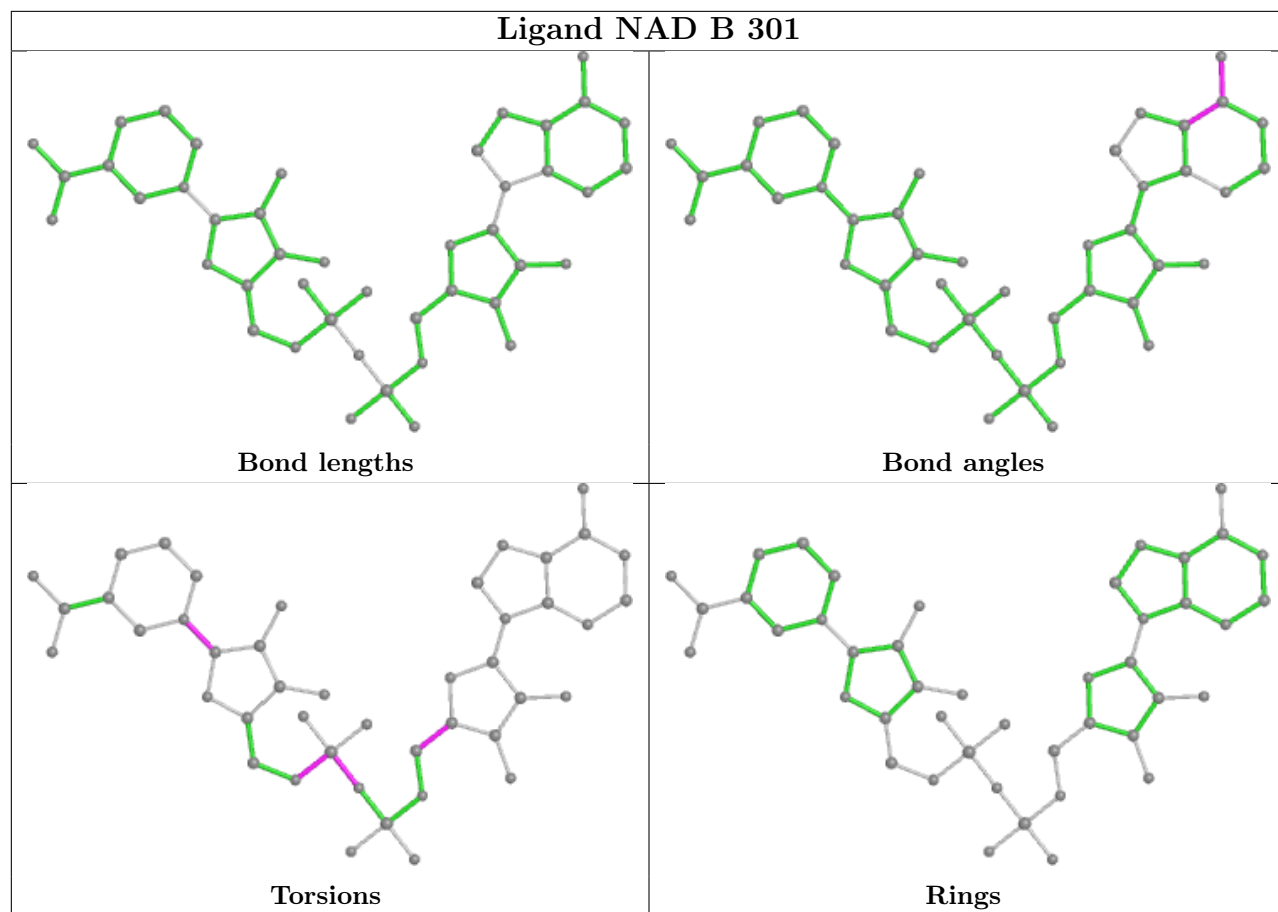


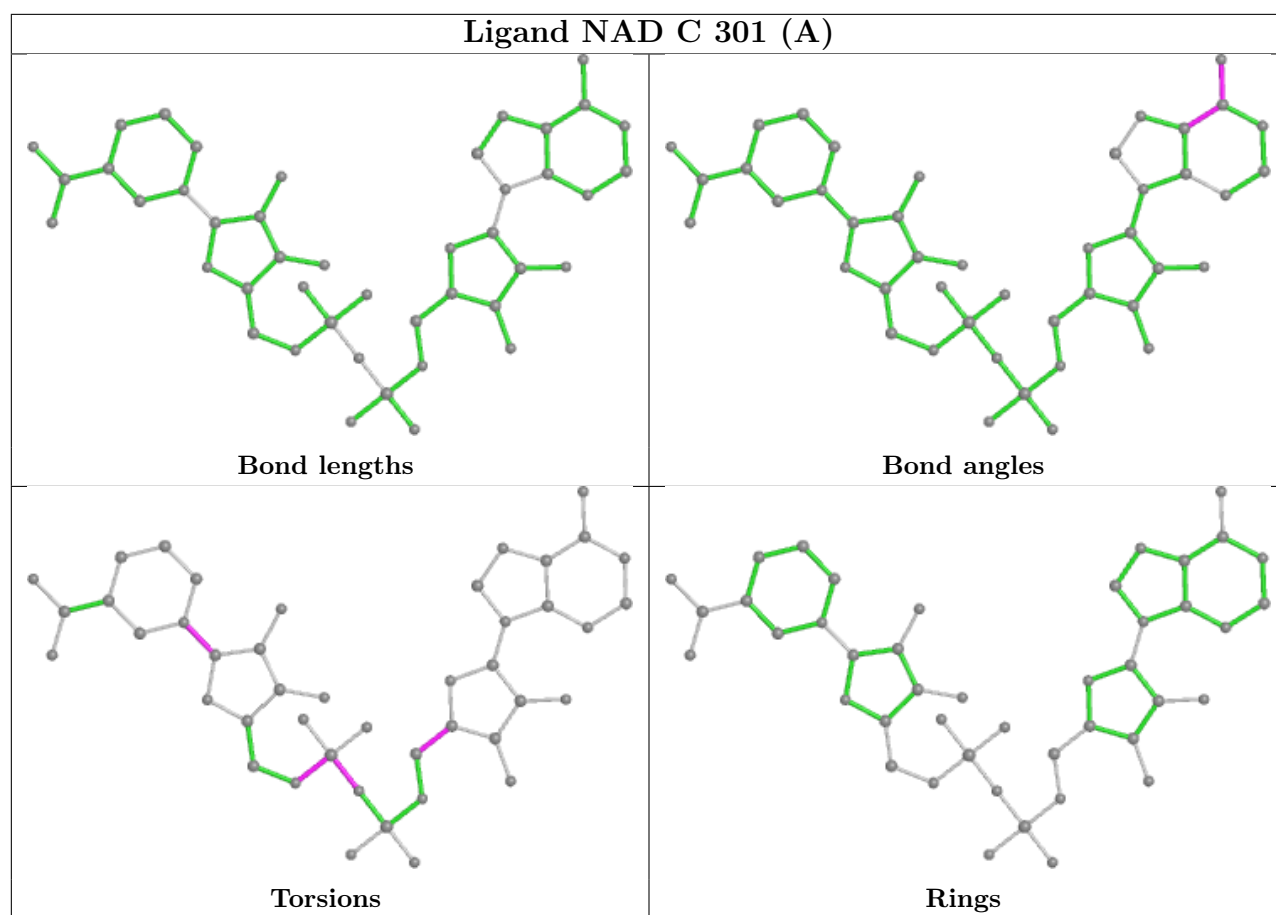












## 5.7 Other polymers [i](#)

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	251/269 (93%)	-0.35	13 (5%) 27 22	13, 20, 55, 78	0
1	B	251/269 (93%)	-0.41	5 (1%) 65 61	13, 20, 47, 81	0
1	C	259/269 (96%)	-0.38	2 (0%) 86 84	13, 25, 49, 63	0
1	D	251/269 (93%)	-0.20	13 (5%) 27 22	13, 25, 57, 88	0
All	All	1012/1076 (94%)	-0.34	33 (3%) 46 40	13, 22, 54, 88	0

All (33) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	198	ILE	9.9
1	A	199	PRO	7.8
1	D	199	PRO	7.8
1	D	198	ILE	7.2
1	A	202	ALA	5.4
1	A	212	VAL	4.0
1	A	201	GLN	3.8
1	D	6	GLY	3.8
1	A	200	GLU	3.5
1	A	213	ILE	3.5
1	B	217	MET	3.2
1	B	210	GLU	3.1
1	D	3	ASN	3.1
1	D	194	VAL	3.1
1	B	212	VAL	3.1
1	D	211	GLU	3.0
1	D	48	ASP	3.0
1	D	212	VAL	3.0
1	A	211	GLU	2.9
1	D	210	GLU	2.8
1	A	216	VAL	2.7

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type	RSRZ
1	A	193	LEU	2.7
1	A	196	LYS	2.6
1	B	199	PRO	2.6
1	A	3	ASN	2.6
1	D	213	ILE	2.4
1	D	193	LEU	2.3
1	B	211	GLU	2.3
1	C	132	ASP	2.2
1	D	200	GLU	2.1
1	D	192	PRO	2.1
1	C	6	GLY	2.1
1	A	192	PRO	2.1

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

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

## 6.3 Carbohydrates [i](#)

There are no monosaccharides in this entry.

## 6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 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.

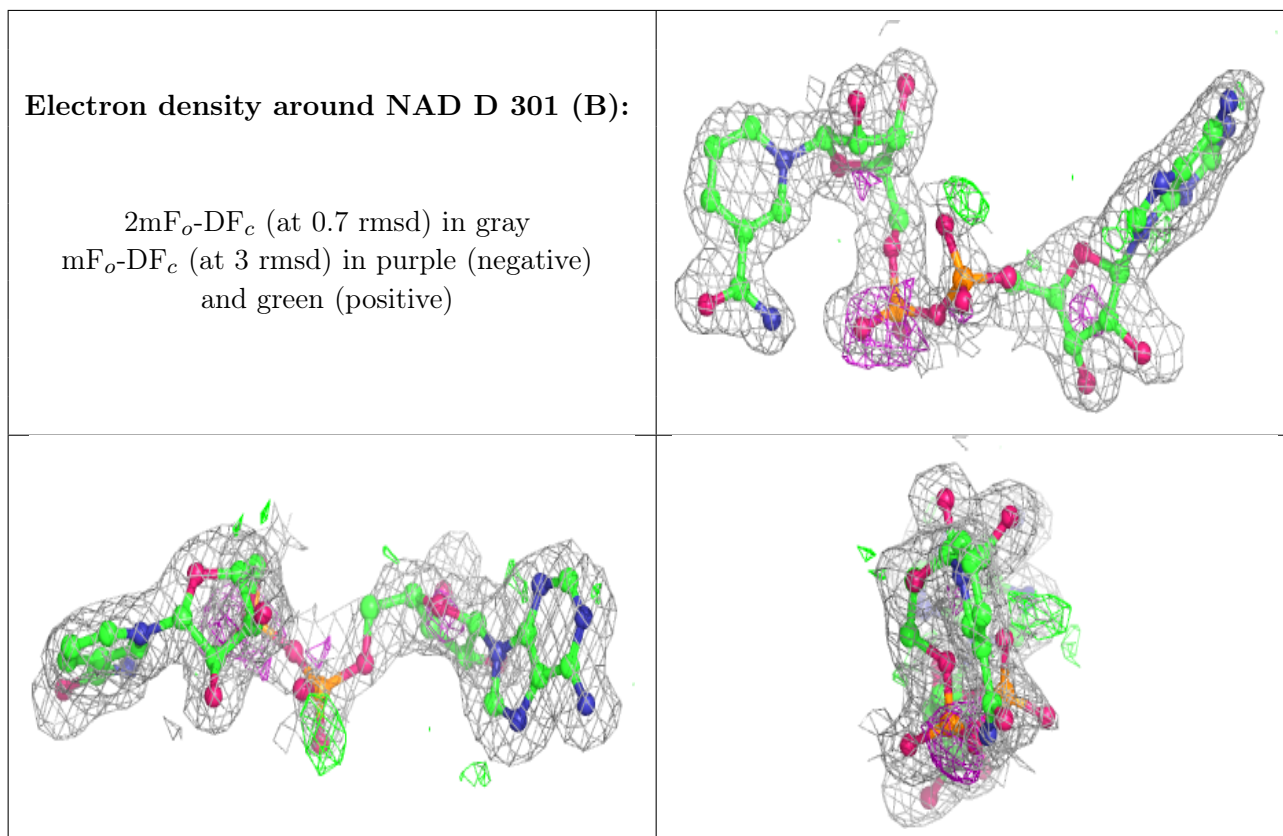
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å <sup>2</sup> )	Q<0.9
3	CIT	C	302	13/13	0.79	0.30	34,45,61,64	0
5	PG4	C	303	10/13	0.81	0.14	36,47,53,61	0
5	PG4	D	304	7/13	0.81	0.16	40,47,52,54	0
5	PG4	D	303	10/13	0.82	0.27	40,48,55,59	0
3	CIT	D	302	13/13	0.83	0.20	30,40,56,60	0
5	PG4	C	304	7/13	0.84	0.18	39,45,56,58	0
3	CIT	A	302	13/13	0.85	0.29	28,47,62,63	0
5	PG4	B	303	10/13	0.85	0.14	29,42,49,53	0
5	PG4	A	304	10/13	0.88	0.22	35,42,48,52	0
3	CIT	B	302	13/13	0.90	0.20	21,33,39,40	0
5	PG4	A	305	7/13	0.94	0.11	32,40,47,59	0

*Continued on next page...*

Continued from previous page...

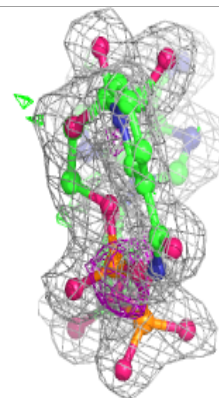
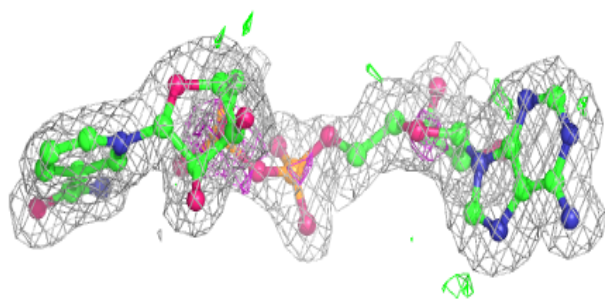
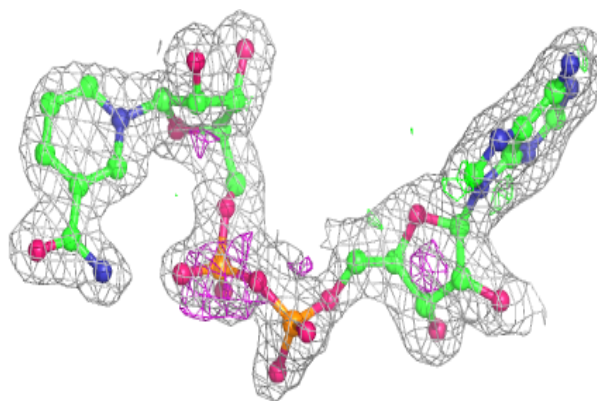
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
2	NAD	D	301[B]	44/44	0.94	0.10	24,27,30,31	44
2	NAD	D	301[A]	44/44	0.94	0.10	22,27,30,30	44
5	PG4	B	304	7/13	0.95	0.15	33,37,43,47	0
2	NAD	A	301[A]	44/44	0.96	0.10	16,22,23,25	44
2	NAD	A	301[B]	44/44	0.96	0.10	16,22,24,26	44
2	NAD	C	301[B]	44/44	0.97	0.08	17,21,25,25	44
2	NAD	B	301	44/44	0.97	0.06	14,17,20,22	0
2	NAD	C	301[A]	44/44	0.97	0.08	15,21,24,26	44
4	PO4	A	303	5/5	0.99	0.07	21,23,24,27	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.

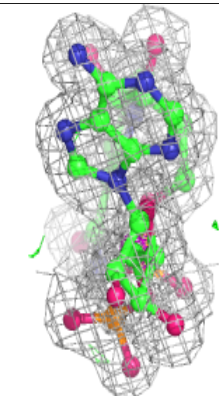
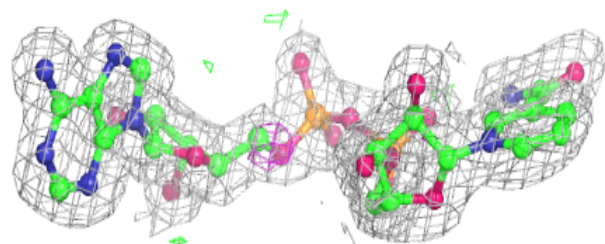
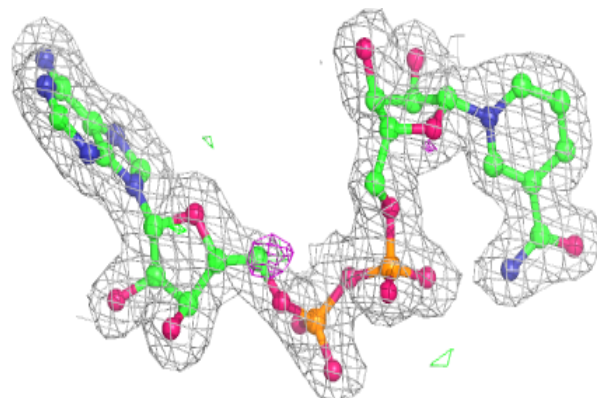


**Electron density around NAD D 301 (A):**

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

**Electron density around NAD A 301 (A):**

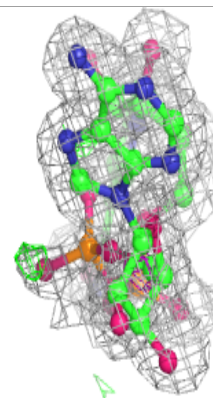
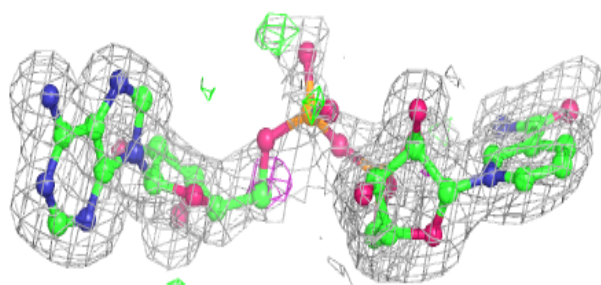
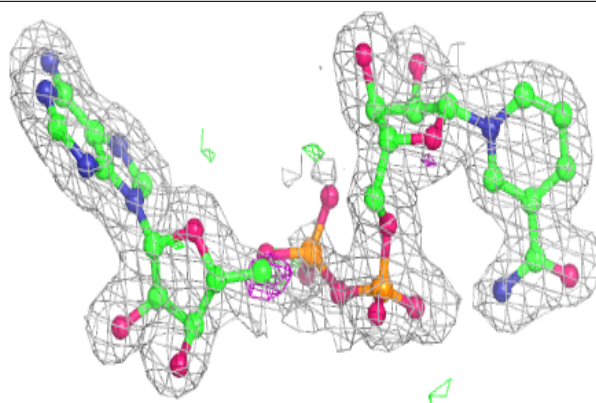
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)



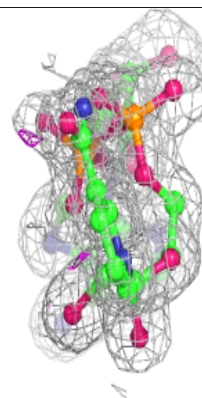
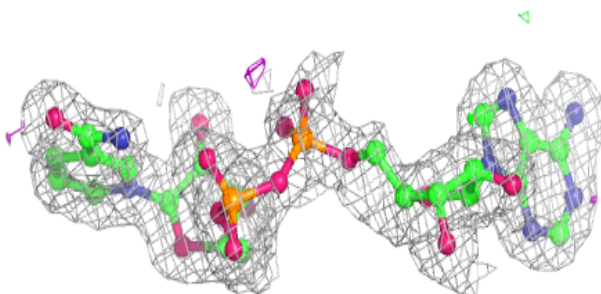
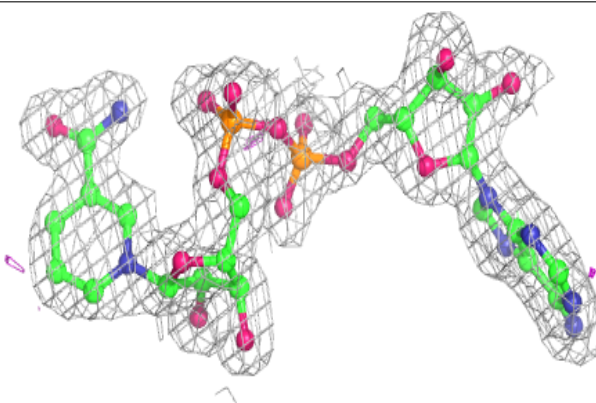


**Electron density around NAD A 301 (B):**

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

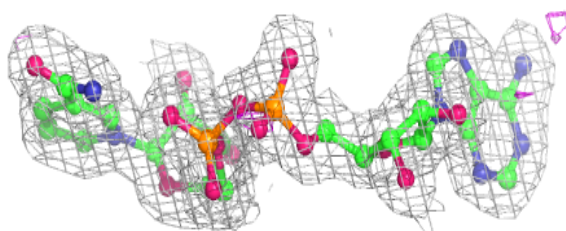
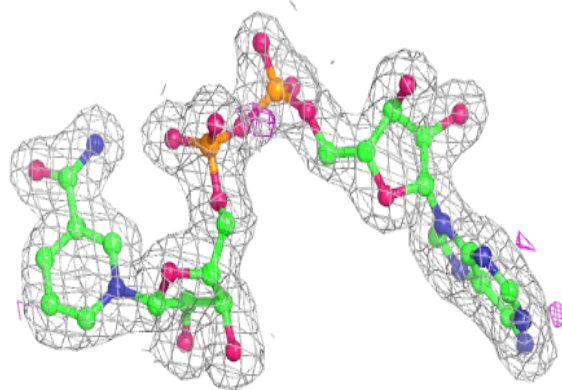
**Electron density around NAD C 301 (B):**

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

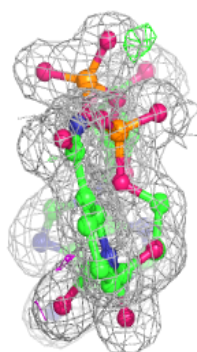
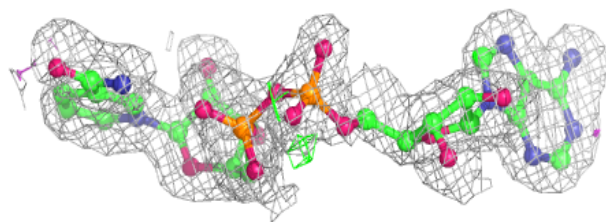
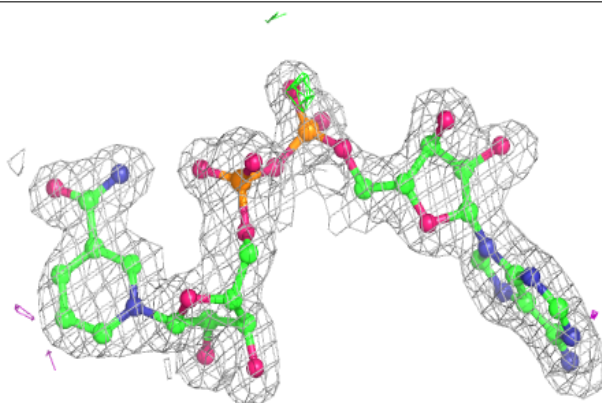


**Electron density around NAD B 301:**

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

**Electron density around NAD C 301 (A):**

$2mF_o-DF_c$  (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.