



# Full wwPDB X-ray Structure Validation Report ⓘ

Nov 16, 2023 – 09:52 AM JST

PDB ID : 6LLA  
Title : Crystal structure of *Providencia alcalifaciens* 3-dehydroquinase synthase (DHQS) in complex with Mg<sup>2+</sup> and NAD  
Authors : Neetu, N.; Katiki, M.; Kumar, P.  
Deposited on : 2019-12-22  
Resolution : 1.88 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 : 4.02b-467  
Mogul : 1.8.5 (274361), CSD as541be (2020)  
Xtriage (Phenix) : 1.13  
EDS : 2.36  
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

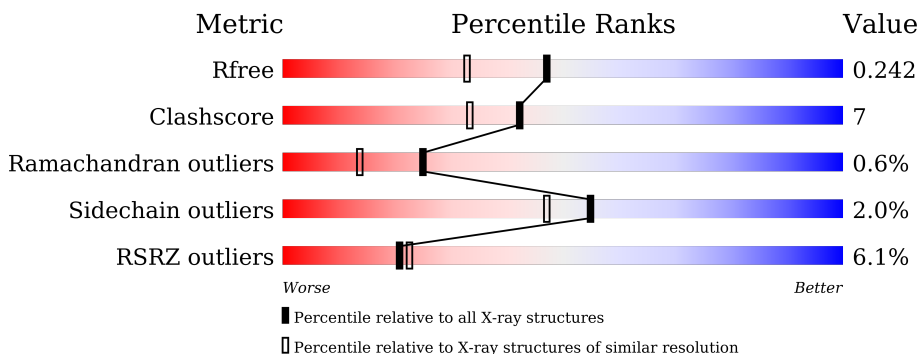
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 1.88 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	9470 (1.90-1.86)
Clashscore	141614	10282 (1.90-1.86)
Ramachandran outliers	138981	10152 (1.90-1.86)
Sidechain outliers	138945	10152 (1.90-1.86)
RSRZ outliers	127900	9303 (1.90-1.86)

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	375	 4% (red), 84% (green), 10% (yellow), 2% (grey)
1	B	375	 3% (red), 84% (green), 11% (yellow), 2% (grey)
1	C	375	 11% (red), 78% (green), 15% (yellow), 6% (grey)
1	D	375	 6% (red), 83% (green), 11% (yellow), 6% (grey)

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit crite-

ria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
4	EDO	B	411	-	-	X	-
4	EDO	C	403	-	-	-	X
5	PEG	A	415	-	-	X	-
5	PEG	B	413	-	-	X	-

## 2 Entry composition i

There are 6 unique types of molecules in this entry. The entry contains 12432 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-dehydroquinate synthase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	362	2786	1769	470	528	19	0	3	0
1	B	363	2787	1769	470	529	19	0	3	0
1	C	354	2697	1717	453	510	17	0	0	0
1	D	353	2701	1719	454	511	17	0	1	0

There are 52 discrepancies between the modelled and reference sequences:

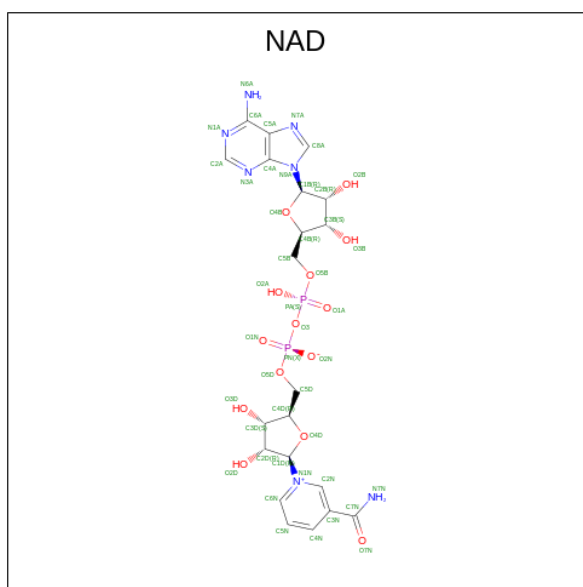
Chain	Residue	Modelled	Actual	Comment	Reference
A	-12	HIS	-	expression tag	UNP X6Q997
A	-11	HIS	-	expression tag	UNP X6Q997
A	-10	HIS	-	expression tag	UNP X6Q997
A	-9	HIS	-	expression tag	UNP X6Q997
A	-8	HIS	-	expression tag	UNP X6Q997
A	-7	HIS	-	expression tag	UNP X6Q997
A	-6	GLU	-	expression tag	UNP X6Q997
A	-5	ASN	-	expression tag	UNP X6Q997
A	-4	LEU	-	expression tag	UNP X6Q997
A	-3	TYR	-	expression tag	UNP X6Q997
A	-2	PHE	-	expression tag	UNP X6Q997
A	-1	GLN	-	expression tag	UNP X6Q997
A	0	GLY	-	expression tag	UNP X6Q997
B	-12	HIS	-	expression tag	UNP X6Q997
B	-11	HIS	-	expression tag	UNP X6Q997
B	-10	HIS	-	expression tag	UNP X6Q997
B	-9	HIS	-	expression tag	UNP X6Q997
B	-8	HIS	-	expression tag	UNP X6Q997
B	-7	HIS	-	expression tag	UNP X6Q997
B	-6	GLU	-	expression tag	UNP X6Q997
B	-5	ASN	-	expression tag	UNP X6Q997

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-4	LEU	-	expression tag	UNP X6Q997
B	-3	TYR	-	expression tag	UNP X6Q997
B	-2	PHE	-	expression tag	UNP X6Q997
B	-1	GLN	-	expression tag	UNP X6Q997
B	0	GLY	-	expression tag	UNP X6Q997
C	-12	HIS	-	expression tag	UNP X6Q997
C	-11	HIS	-	expression tag	UNP X6Q997
C	-10	HIS	-	expression tag	UNP X6Q997
C	-9	HIS	-	expression tag	UNP X6Q997
C	-8	HIS	-	expression tag	UNP X6Q997
C	-7	HIS	-	expression tag	UNP X6Q997
C	-6	GLU	-	expression tag	UNP X6Q997
C	-5	ASN	-	expression tag	UNP X6Q997
C	-4	LEU	-	expression tag	UNP X6Q997
C	-3	TYR	-	expression tag	UNP X6Q997
C	-2	PHE	-	expression tag	UNP X6Q997
C	-1	GLN	-	expression tag	UNP X6Q997
C	0	GLY	-	expression tag	UNP X6Q997
D	-12	HIS	-	expression tag	UNP X6Q997
D	-11	HIS	-	expression tag	UNP X6Q997
D	-10	HIS	-	expression tag	UNP X6Q997
D	-9	HIS	-	expression tag	UNP X6Q997
D	-8	HIS	-	expression tag	UNP X6Q997
D	-7	HIS	-	expression tag	UNP X6Q997
D	-6	GLU	-	expression tag	UNP X6Q997
D	-5	ASN	-	expression tag	UNP X6Q997
D	-4	LEU	-	expression tag	UNP X6Q997
D	-3	TYR	-	expression tag	UNP X6Q997
D	-2	PHE	-	expression tag	UNP X6Q997
D	-1	GLN	-	expression tag	UNP X6Q997
D	0	GLY	-	expression tag	UNP X6Q997

- Molecule 2 is NICOTINAMIDE-ADENINE-DINUCLEOTIDE (three-letter code: NAD) (formula: C<sub>21</sub>H<sub>27</sub>N<sub>7</sub>O<sub>14</sub>P<sub>2</sub>) (labeled as "Ligand of Interest" by depositor).



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

- Molecule 3 is MAGNESIUM ION (three-letter code: MG) (formula: Mg) (labeled as "Ligand of Interest" by depositor).

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

- Molecule 4 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: C<sub>2</sub>H<sub>6</sub>O<sub>2</sub>).



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

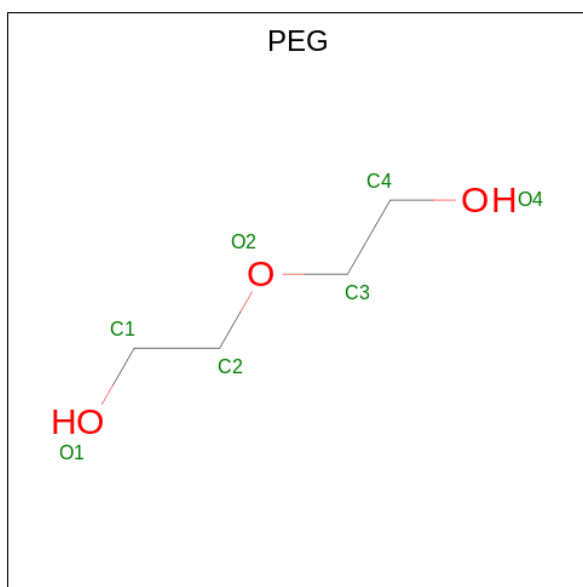
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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	B	1	Total C O 4 2 2	0	0
4	B	1	Total C O 4 2 2	0	0
4	B	1	Total C O 4 2 2	0	0
4	B	1	Total C O 4 2 2	0	0
4	B	1	Total C O 4 2 2	0	0
4	B	1	Total C O 4 2 2	0	0
4	B	1	Total C O 4 2 2	0	0
4	C	1	Total C O 4 2 2	0	0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

- Molecule 5 is DI(HYDROXYETHYL)ETHER (three-letter code: PEG) (formula: C<sub>4</sub>H<sub>10</sub>O<sub>3</sub>).





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

- Molecule 6 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	A	342	Total O 342 342	0	0
6	B	341	Total O 341 341	0	0
6	C	178	Total O 178 178	0	0

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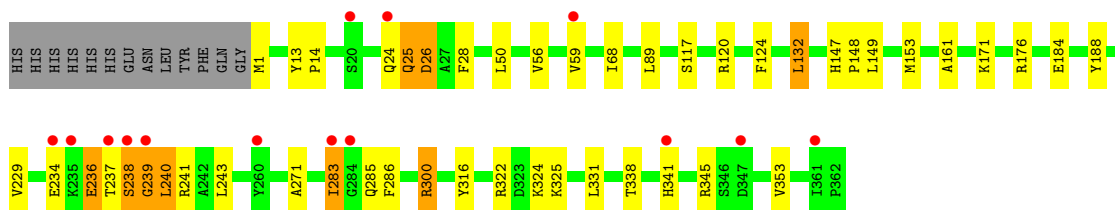
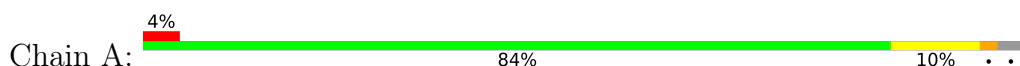
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<b>Mol</b>	<b>Chain</b>	<b>Residues</b>	<b>Atoms</b>		<b>ZeroOcc</b>	<b>AltConf</b>
6	D	233	Total 233	O 233	0	0

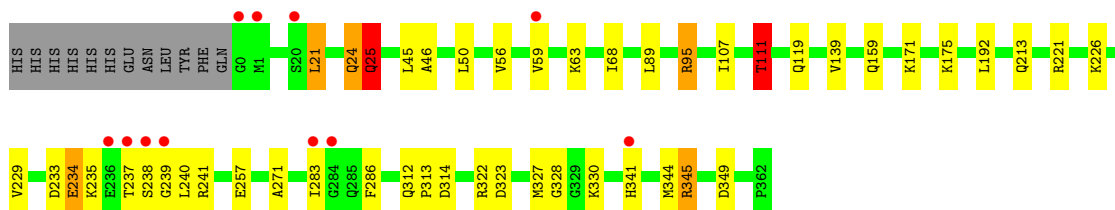
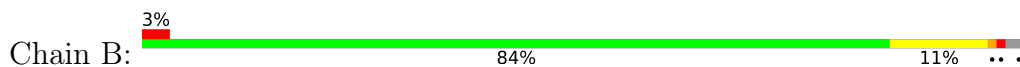
### 3 Residue-property plots [i](#)

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.

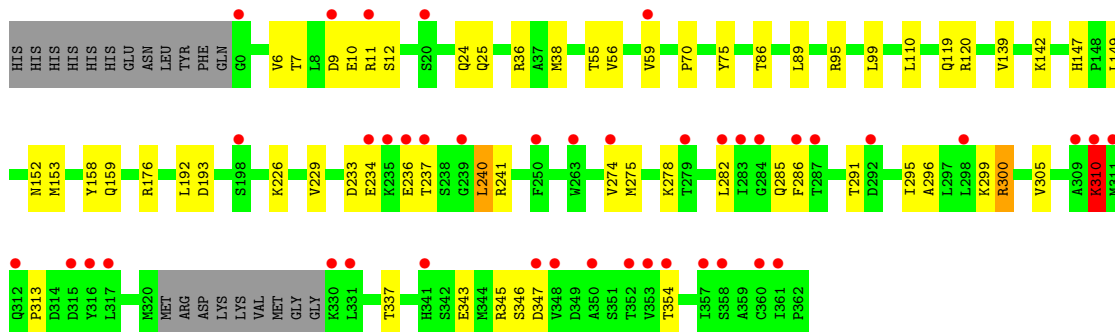
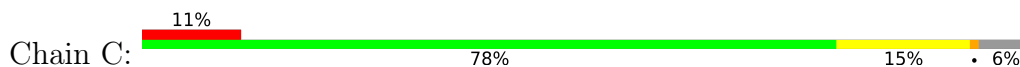
- Molecule 1: 3-dehydroquinate synthase




- Molecule 1: 3-dehydroquinate synthase

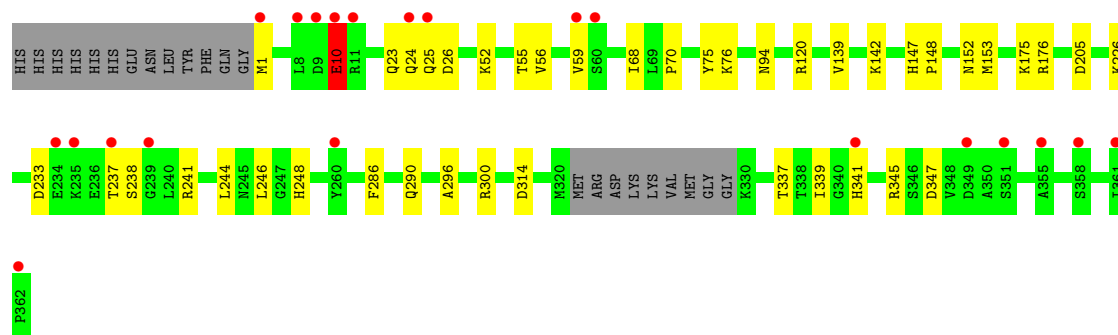


- Molecule 1: 3-dehydroquinate synthase



- Molecule 1: 3-dehydroquinate synthase

Chain D: 



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	89.39Å 59.92Å 143.80Å 90.00° 93.95° 90.00°	Depositor
Resolution (Å)	143.46 – 1.88 89.18 – 1.88	Depositor EDS
% Data completeness (in resolution range)	98.9 (143.46-1.88) 98.9 (89.18-1.88)	Depositor EDS
$R_{merge}$	0.05	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.98 (at 1.88Å)	Xtrriage
Refinement program	REFMAC 5.8.0158	Depositor
R, $R_{free}$	0.173 , 0.238 0.181 , 0.242	Depositor DCC
$R_{free}$ test set	6015 reflections (4.88%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	28.9	Xtrriage
Anisotropy	0.435	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.35 , 59.3	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	12432	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	42.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 8.23% 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: EDO, PEG, NAD, MG

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	1.05	6/2837 (0.2%)	0.91	4/3851 (0.1%)
1	B	1.09	3/2838 (0.1%)	0.96	7/3851 (0.2%)
1	C	0.81	0/2747	0.91	11/3731 (0.3%)
1	D	0.82	0/2751	0.88	7/3737 (0.2%)
All	All	0.96	9/11173 (0.1%)	0.92	29/15170 (0.2%)

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 (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	111	THR	CB-CG2	-6.69	1.30	1.52
1	B	271	ALA	CA-CB	5.93	1.65	1.52
1	A	117	SER	CB-OG	-5.72	1.34	1.42
1	A	188	TYR	CG-CD2	-5.65	1.31	1.39
1	A	271	ALA	CA-CB	5.54	1.64	1.52
1	A	13	TYR	CE2-CZ	-5.27	1.31	1.38
1	A	316	TYR	CE1-CZ	5.25	1.45	1.38
1	A	124	PHE	CG-CD1	-5.21	1.30	1.38
1	B	221	ARG	CG-CD	5.09	1.64	1.51

All (29) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	310	LYS	CA-CB-CG	10.71	136.95	113.40
1	C	310	LYS	N-CA-CB	-10.06	92.50	110.60
1	C	310	LYS	CB-CG-CD	-9.33	87.33	111.60
1	D	314	ASP	CB-CA-C	-7.91	94.58	110.40
1	D	10	GLU	CA-CB-CG	-7.37	97.19	113.40
1	C	310	LYS	CD-CE-NZ	-7.11	95.34	111.70
1	C	120	ARG	NE-CZ-NH1	6.87	123.73	120.30
1	C	300	ARG	NE-CZ-NH2	-6.76	116.92	120.30
1	D	120	ARG	NE-CZ-NH1	6.72	123.66	120.30
1	D	241	ARG	NE-CZ-NH2	6.50	123.55	120.30
1	D	241	ARG	NE-CZ-NH1	-6.48	117.06	120.30
1	B	95	ARG	CA-CB-CG	6.32	127.30	113.40
1	A	300	ARG	NE-CZ-NH2	-5.98	117.31	120.30
1	B	345	ARG	NE-CZ-NH2	-5.96	117.32	120.30
1	A	132	LEU	CB-CG-CD1	5.94	121.09	111.00
1	A	283	ILE	CG1-CB-CG2	-5.89	98.44	111.40
1	B	349	ASP	CB-CG-OD1	5.87	123.58	118.30
1	D	205	ASP	CB-CG-OD1	5.84	123.56	118.30
1	A	26	ASP	CB-CG-OD2	5.79	123.51	118.30
1	C	275	MET	CB-CG-SD	5.67	129.40	112.40
1	D	26	ASP	N-CA-C	-5.59	95.91	111.00
1	B	327	MET	CG-SD-CE	5.46	108.93	100.20
1	C	310	LYS	CG-CD-CE	5.46	128.27	111.90
1	C	110	LEU	CB-CG-CD2	-5.44	101.76	111.00
1	B	221	ARG	CG-CD-NE	-5.37	100.52	111.80
1	C	193	ASP	CB-CG-OD2	-5.33	113.50	118.30
1	C	36	ARG	NE-CZ-NH1	-5.32	117.64	120.30
1	B	45	LEU	CA-CB-CG	5.02	126.85	115.30
1	B	21	LEU	CA-CB-CG	5.00	126.81	115.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	234	GLU	Peptide

## 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	2786	0	2816	48	0
1	B	2787	0	2815	42	0
1	C	2697	0	2728	48	0
1	D	2701	0	2730	28	0
2	A	44	0	26	0	0
2	B	44	0	26	0	0
2	C	44	0	26	0	0
2	D	44	0	26	0	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	44	0	66	4	0
4	B	40	0	60	5	0
4	C	16	0	24	5	0
4	D	24	0	36	5	0
5	A	42	0	60	12	0
5	B	14	0	20	6	0
5	C	7	0	10	2	0
6	A	342	0	0	10	0
6	B	341	0	0	13	1
6	C	178	0	0	5	0
6	D	233	0	0	4	1
All	All	12432	0	11469	158	1

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 7.

All (158) 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:89:LEU:HD21	1:C:153:MET:HE2	1.13	1.09
1:A:322:ARG:NH2	6:A:501:HOH:O	1.87	1.05
1:A:89:LEU:HD21	1:C:153:MET:CE	1.93	0.98
1:B:313:PRO:HB2	5:B:413:PEG:H42	1.54	0.88
1:A:120:ARG:HH12	4:C:406:EDO:H22	1.39	0.88
1:D:94[A]:ASN:ND2	6:D:501:HOH:O	2.06	0.87
1:C:147:HIS:HD2	1:C:149:LEU:H	1.22	0.86
1:B:24:GLN:HE21	1:B:25:GLN:HE21	1.21	0.86
1:A:147:HIS:HD2	1:A:149:LEU:H	1.24	0.86
1:B:233:ASP:OD2	1:B:240:LEU:N	2.09	0.84
1:C:345:ARG:NH1	1:C:346:SER:O	2.11	0.82

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:56:VAL:O	1:A:59:VAL:HG12	1.80	0.82
1:C:159:GLN:NE2	6:C:502:HOH:O	2.17	0.78
1:C:274:VAL:CG1	1:C:295:ILE:HD11	2.13	0.77
1:B:119:GLN:HG2	1:D:153:MET:HE3	1.65	0.76
5:C:407:PEG:O4	6:C:501:HOH:O	2.04	0.74
1:B:56:VAL:O	1:B:59:VAL:HG12	1.91	0.71
1:A:341:HIS:HE1	6:B:676:HOH:O	1.73	0.70
1:C:274:VAL:HG11	1:C:295:ILE:HD11	1.71	0.70
1:A:153:MET:HE3	1:C:119:GLN:HG2	1.72	0.70
1:A:285:GLN:OE1	1:A:345:ARG:NH1	2.26	0.68
1:A:341:HIS:CD2	5:A:419:PEG:O4	2.47	0.67
1:D:76:LYS:NZ	1:D:152:ASN:HD21	1.93	0.67
1:C:147:HIS:CD2	1:C:149:LEU:H	2.10	0.67
1:D:76:LYS:HZ1	1:D:152:ASN:HD21	1.41	0.67
1:A:26:ASP:HA	4:A:406:EDO:H21	1.78	0.66
1:B:257:GLU:OE1	5:B:414:PEG:O4	2.14	0.66
1:A:234:GLU:O	6:A:503:HOH:O	2.14	0.66
1:A:24:GLN:NE2	4:A:413:EDO:H11	2.13	0.64
1:B:107:ILE:O	1:B:111:THR:HG23	1.98	0.64
5:A:418:PEG:O4	6:A:502:HOH:O	2.09	0.63
1:C:296:ALA:O	1:C:300:ARG:HG2	1.98	0.63
1:B:313:PRO:CB	5:B:413:PEG:H42	2.28	0.63
1:A:25:GLN:NE2	1:A:26:ASP:OD1	2.33	0.62
1:B:241:ARG:NH2	4:B:411:EDO:O1	2.26	0.61
1:A:147:HIS:CD2	1:A:149:LEU:H	2.11	0.60
1:C:274:VAL:HG13	1:C:295:ILE:HD11	1.81	0.60
1:A:300:ARG:HD2	6:A:557:HOH:O	2.02	0.60
1:B:175:LYS:HE2	6:B:559:HOH:O	2.00	0.60
1:B:241:ARG:HH21	4:B:411:EDO:HO1	1.49	0.59
1:D:142:LYS:HD2	4:D:403:EDO:H11	1.83	0.59
1:B:89:LEU:HD21	1:D:153:MET:HE2	1.83	0.58
1:B:233:ASP:OD2	1:B:239:GLY:N	2.35	0.58
1:A:229:VAL:HG23	1:A:240:LEU:HD22	1.85	0.57
1:A:341:HIS:HD2	5:A:419:PEG:O4	1.85	0.57
1:A:50:LEU:HD22	1:A:68:ILE:HD11	1.85	0.57
1:D:246:LEU:O	1:D:248:HIS:HD2	1.86	0.57
1:B:24:GLN:HB3	1:B:25:GLN:HG2	1.86	0.56
1:B:56:VAL:HG11	6:B:506:HOH:O	2.03	0.56
1:C:337:THR:CG2	1:C:343:GLU:HG2	2.35	0.56
1:B:313:PRO:HB2	5:B:413:PEG:C4	2.33	0.55
1:C:55:THR:O	1:C:59:VAL:HG13	2.07	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:310:LYS:NZ	6:C:506:HOH:O	2.40	0.54
1:D:55:THR:O	1:D:59:VAL:HG13	2.07	0.54
1:B:314:ASP:HB2	5:B:413:PEG:H22	1.89	0.54
1:C:9:ASP:HB2	1:C:234:GLU:OE2	2.08	0.54
1:B:312:GLN:NE2	6:B:504:HOH:O	2.33	0.54
1:B:89:LEU:HD21	1:D:153:MET:CE	2.38	0.54
1:C:233:ASP:OD1	1:C:236:GLU:HA	2.07	0.54
1:C:158:TYR:CD1	4:C:404:EDO:H22	2.44	0.52
1:B:50:LEU:HD22	1:B:68:ILE:HD11	1.89	0.52
1:A:171:LYS:NZ	6:A:513:HOH:O	2.43	0.52
1:C:278:LYS:O	1:C:282:LEU:HD13	2.08	0.52
1:B:56:VAL:CG1	6:B:506:HOH:O	2.57	0.52
1:D:337:THR:OG1	1:D:341:HIS:CG	2.64	0.51
1:C:299:LYS:HE3	1:C:305:VAL:CG2	2.40	0.51
1:C:299:LYS:HE3	1:C:305:VAL:HG21	1.93	0.51
1:C:299:LYS:NZ	6:C:511:HOH:O	2.44	0.51
1:C:38:MET:HB3	1:C:99:LEU:HD23	1.93	0.50
1:D:345:ARG:HD2	1:D:347:ASP:OD2	2.11	0.50
1:A:14:PRO:HG3	5:A:414:PEG:H32	1.92	0.50
1:C:192:LEU:HD11	1:C:229:VAL:HG11	1.94	0.50
1:A:28:PHE:H	4:A:406:EDO:H11	1.75	0.50
1:B:192:LEU:HD11	1:B:229:VAL:HG11	1.94	0.50
1:A:1:MET:HB3	5:A:418:PEG:H22	1.94	0.50
1:D:175:LYS:O	1:D:175:LYS:HD2	2.12	0.49
1:B:119:GLN:CG	1:D:153:MET:HE3	2.38	0.49
1:A:24:GLN:HE22	4:A:413:EDO:H11	1.77	0.49
1:B:323:ASP:OD1	1:B:328:GLY:HA3	2.13	0.49
1:A:240:LEU:HD23	1:A:243:LEU:HD12	1.95	0.48
1:C:337:THR:HG23	1:C:343:GLU:HG2	1.94	0.48
1:C:142:LYS:NZ	4:C:406:EDO:H21	2.29	0.48
1:A:237:THR:O	1:A:239:GLY:N	2.45	0.48
1:A:338:THR:OG1	1:A:341:HIS:HB2	2.12	0.48
1:D:70:PRO:HB2	1:D:75:TYR:CE2	2.49	0.47
1:A:241:ARG:HE	5:A:415:PEG:H41	1.79	0.47
1:A:153:MET:CE	1:C:89:LEU:HD21	2.44	0.47
1:C:11:ARG:HE	4:C:404:EDO:H11	1.79	0.47
1:B:213:GLN:HG3	6:B:645:HOH:O	2.15	0.47
1:A:238:SER:O	6:A:504:HOH:O	2.20	0.47
1:B:63:LYS:HG2	6:B:608:HOH:O	2.14	0.47
1:D:23:GLN:HE21	1:D:52:LYS:NZ	2.12	0.47
1:A:241:ARG:HE	5:A:415:PEG:H32	1.79	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:341:HIS:HD2	6:B:733:HOH:O	1.97	0.47
1:C:274:VAL:HG11	1:C:295:ILE:CD1	2.41	0.47
1:D:296:ALA:O	1:D:300:ARG:HG3	2.15	0.47
5:A:416:PEG:H21	1:C:152:ASN:O	2.15	0.47
1:B:226:LYS:HZ3	4:B:411:EDO:H12	1.80	0.47
1:C:236:GLU:HG2	1:C:241:ARG:HD3	1.97	0.46
1:A:325:LYS:HE3	6:A:521:HOH:O	2.16	0.46
1:B:171:LYS:NZ	6:B:512:HOH:O	2.48	0.46
1:B:322:ARG:HH11	1:B:330:LYS:HE2	1.80	0.46
1:B:226:LYS:NZ	4:B:411:EDO:H12	2.31	0.45
5:A:417:PEG:H21	6:C:648:HOH:O	2.16	0.45
1:C:70:PRO:HB2	1:C:75:TYR:CE2	2.50	0.45
1:A:56:VAL:CG1	6:A:510:HOH:O	2.64	0.45
1:B:314:ASP:OD2	5:B:413:PEG:O4	2.34	0.45
1:C:241:ARG:HH21	5:C:407:PEG:H12	1.81	0.45
1:D:233:ASP:OD1	1:D:238:SER:HB3	2.16	0.45
1:A:331:LEU:HD13	1:A:353:VAL:HG21	1.99	0.45
1:B:241:ARG:NH1	4:B:407:EDO:O1	2.35	0.45
1:C:10:GLU:HA	1:C:10:GLU:OE1	2.17	0.45
1:C:274:VAL:CG1	1:C:295:ILE:CD1	2.92	0.45
1:C:313:PRO:HB3	1:C:354:THR:HG23	1.99	0.44
1:A:56:VAL:HG11	6:A:510:HOH:O	2.17	0.44
1:B:25:GLN:O	6:B:501:HOH:O	2.20	0.44
1:A:283:ILE:HD13	1:A:345:ARG:NH2	2.33	0.44
1:C:345:ARG:HH12	1:C:347:ASP:C	2.20	0.44
1:B:283:ILE:HD12	1:B:345:ARG:CZ	2.48	0.44
1:B:159:GLN:NE2	6:B:508:HOH:O	2.49	0.43
1:A:147:HIS:CG	1:A:148:PRO:HD2	2.53	0.43
1:D:56:VAL:O	1:D:59:VAL:HG22	2.18	0.43
1:A:241:ARG:HE	5:A:415:PEG:C3	2.32	0.43
1:C:24:GLN:OE1	1:C:25:GLN:O	2.37	0.43
1:C:7:THR:HG22	1:C:12:SER:HB2	2.00	0.43
1:A:147:HIS:HE1	1:C:86:THR:OG1	2.02	0.42
1:B:46:ALA:HB2	1:B:68:ILE:HD12	2.01	0.42
1:D:139:VAL:HB	1:D:226:LYS:HD3	2.01	0.42
1:D:341:HIS:ND1	6:D:504:HOH:O	2.36	0.42
1:A:153:MET:HE3	1:C:119:GLN:CG	2.44	0.42
1:A:229:VAL:CG2	1:A:240:LEU:HD22	2.50	0.42
1:B:95:ARG:HG3	6:B:628:HOH:O	2.20	0.42
1:D:1:MET:N	6:D:511:HOH:O	2.52	0.42
1:A:241:ARG:HH21	5:A:415:PEG:H41	1.85	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:341:HIS:CD2	6:B:733:HOH:O	2.72	0.42
1:D:68:ILE:HD13	4:D:407:EDO:H21	2.00	0.42
1:D:341:HIS:CD2	4:D:404:EDO:O1	2.73	0.42
1:C:24:GLN:CD	1:C:25:GLN:O	2.58	0.42
1:A:283:ILE:HD13	1:A:283:ILE:HG21	1.82	0.42
1:B:139:VAL:HB	1:B:226:LYS:HD3	2.01	0.41
1:D:290:GLN:HB3	1:D:339:ILE:HD11	2.02	0.41
1:A:132:LEU:CD2	1:A:184:GLU:HG3	2.50	0.41
1:D:10:GLU:HA	6:D:546:HOH:O	2.19	0.41
1:C:6:VAL:O	1:C:12:SER:HA	2.19	0.41
1:C:56:VAL:O	1:C:59:VAL:HG22	2.20	0.41
1:C:139:VAL:HB	1:C:226:LYS:HD3	2.03	0.41
1:A:161:ALA:HA	6:A:736:HOH:O	2.20	0.41
1:B:233:ASP:O	1:B:237:THR:HA	2.20	0.41
1:D:244:LEU:HD12	4:D:403:EDO:H22	2.02	0.41
1:A:283:ILE:HG21	1:A:345:ARG:HH22	1.85	0.41
1:C:233:ASP:OD2	1:C:240:LEU:HB3	2.20	0.41
1:D:341:HIS:HD2	4:D:404:EDO:O1	2.03	0.40
1:C:142:LYS:HZ3	4:C:406:EDO:H21	1.86	0.40
1:B:233:ASP:HB3	1:B:241:ARG:HG3	2.02	0.40
1:A:153:MET:HE2	1:C:89:LEU:HD21	2.03	0.40
1:A:241:ARG:O	5:A:415:PEG:H11	2.22	0.40
1:C:278:LYS:HD3	1:C:291:THR:HG21	2.03	0.40
1:D:147:HIS:CG	1:D:148:PRO:HD2	2.57	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:B:593:HOH:O	6:D:659:HOH:O[2_545]	2.15	0.05

## 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	363/375 (97%)	350 (96%)	9 (2%)	4 (1%)	14	5
1	B	364/375 (97%)	355 (98%)	8 (2%)	1 (0%)	41	30
1	C	350/375 (93%)	343 (98%)	6 (2%)	1 (0%)	41	30
1	D	350/375 (93%)	339 (97%)	9 (3%)	2 (1%)	25	14
All	All	1427/1500 (95%)	1387 (97%)	32 (2%)	8 (1%)	25	14

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	25	GLN
1	A	238	SER
1	C	237	THR
1	A	236	GLU
1	B	25	GLN
1	D	25	GLN
1	D	10	GLU
1	A	239	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	300/309 (97%)	295 (98%)	5 (2%)	60	54
1	B	300/309 (97%)	291 (97%)	9 (3%)	41	30
1	C	290/309 (94%)	284 (98%)	6 (2%)	53	45
1	D	291/309 (94%)	287 (99%)	4 (1%)	67	62
All	All	1181/1236 (96%)	1157 (98%)	24 (2%)	55	47

All (24) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	176	ARG

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Mol	Chain	Res	Type
1	A	236	GLU
1	A	240	LEU
1	A	286	PHE
1	A	324	LYS
1	B	21	LEU
1	B	24	GLN
1	B	25	GLN
1	B	111	THR
1	B	234	GLU
1	B	235	LYS
1	B	238	SER
1	B	286	PHE
1	B	344	MET
1	C	95	ARG
1	C	176	ARG
1	C	240	LEU
1	C	285	GLN
1	C	286	PHE
1	C	310	LYS
1	D	24	GLN
1	D	176	ARG
1	D	237	THR
1	D	286	PHE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (15) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	24	GLN
1	A	147	HIS
1	A	341	HIS
1	B	24	GLN
1	B	159	GLN
1	B	341	HIS
1	C	82	ASN
1	C	119	GLN
1	C	147	HIS
1	C	213	GLN
1	C	285	GLN
1	D	23	GLN
1	D	152	ASN
1	D	248	HIS
1	D	341	HIS

### 5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

## 5.5 Carbohydrates [i](#)

There are no monosaccharides in this entry.

## 5.6 Ligand geometry [i](#)

Of 48 ligands modelled in this entry, 4 are monoatomic - leaving 44 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with  $|Z| > 2$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
2	NAD	B	401	-	42,48,48	1.52	6 (14%)	50,73,73	1.60	6 (12%)
4	EDO	D	407	-	3,3,3	0.49	0	2,2,2	0.54	0
4	EDO	D	408	-	3,3,3	0.68	0	2,2,2	0.19	0
4	EDO	A	411	-	3,3,3	0.64	0	2,2,2	0.28	0
2	NAD	C	401	-	42,48,48	1.62	9 (21%)	50,73,73	0.90	2 (4%)
4	EDO	C	403	-	3,3,3	0.62	0	2,2,2	0.26	0
5	PEG	A	419	-	6,6,6	0.87	0	5,5,5	0.71	0
4	EDO	D	403	-	3,3,3	0.45	0	2,2,2	0.19	0
4	EDO	A	408	-	3,3,3	0.57	0	2,2,2	0.32	0
4	EDO	B	411	-	3,3,3	0.94	0	2,2,2	0.82	0
4	EDO	B	404	-	3,3,3	0.86	0	2,2,2	0.41	0
5	PEG	C	407	-	6,6,6	0.55	0	5,5,5	0.47	0
5	PEG	A	415	-	6,6,6	0.71	0	5,5,5	0.77	0
4	EDO	A	405	-	3,3,3	0.82	0	2,2,2	0.14	0
4	EDO	D	406	-	3,3,3	0.40	0	2,2,2	0.30	0
4	EDO	B	406	-	3,3,3	0.63	0	2,2,2	0.67	0
4	EDO	A	409	-	3,3,3	0.28	0	2,2,2	0.28	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
2	NAD	A	401	-	42,48,48	1.67	6 (14%)	50,73,73	1.46	7 (14%)
4	EDO	B	409	-	3,3,3	0.65	0	2,2,2	0.26	0
4	EDO	A	407	-	3,3,3	0.65	0	2,2,2	0.75	0
4	EDO	C	406	-	3,3,3	0.47	0	2,2,2	1.11	0
5	PEG	A	417	-	6,6,6	0.51	0	5,5,5	0.70	0
4	EDO	A	412	-	3,3,3	0.44	0	2,2,2	0.25	0
4	EDO	C	405	-	3,3,3	0.24	0	2,2,2	0.78	0
4	EDO	A	410	-	3,3,3	0.38	0	2,2,2	0.46	0
4	EDO	B	407	-	3,3,3	0.41	0	2,2,2	0.96	0
4	EDO	A	404	-	3,3,3	0.47	0	2,2,2	0.35	0
4	EDO	B	405	-	3,3,3	0.58	0	2,2,2	0.55	0
4	EDO	A	403	-	3,3,3	0.51	0	2,2,2	0.44	0
4	EDO	B	412	-	3,3,3	0.35	0	2,2,2	0.48	0
4	EDO	B	410	-	3,3,3	0.51	0	2,2,2	0.80	0
4	EDO	A	406	-	3,3,3	0.43	0	2,2,2	0.77	0
4	EDO	D	404	-	3,3,3	0.37	0	2,2,2	0.96	0
2	NAD	D	401	-	42,48,48	1.61	4 (9%)	50,73,73	1.38	4 (8%)
5	PEG	A	416	-	6,6,6	0.47	0	5,5,5	0.60	0
4	EDO	B	403	-	3,3,3	0.60	0	2,2,2	0.47	0
5	PEG	A	418	-	6,6,6	0.56	0	5,5,5	0.62	0
4	EDO	C	404	-	3,3,3	0.71	0	2,2,2	1.00	0
5	PEG	A	414	-	6,6,6	0.74	0	5,5,5	0.89	0
4	EDO	A	413	-	3,3,3	0.58	0	2,2,2	0.06	0
4	EDO	B	408	-	3,3,3	0.49	0	2,2,2	0.18	0
4	EDO	D	405	-	3,3,3	0.27	0	2,2,2	0.67	0
5	PEG	B	413	-	6,6,6	0.53	0	5,5,5	1.33	1 (20%)
5	PEG	B	414	-	6,6,6	0.68	0	5,5,5	0.51	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
2	NAD	B	401	-	-	3/26/62/62	0/5/5/5
4	EDO	D	407	-	-	1/1/1/1	-
4	EDO	D	408	-	-	0/1/1/1	-
4	EDO	A	411	-	-	1/1/1/1	-
2	NAD	C	401	-	-	2/26/62/62	0/5/5/5
4	EDO	C	403	-	-	0/1/1/1	-
5	PEG	A	419	-	-	4/4/4/4	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	EDO	D	403	-	-	0/1/1/1	-
4	EDO	A	408	-	-	1/1/1/1	-
4	EDO	B	411	-	-	1/1/1/1	-
4	EDO	B	404	-	-	0/1/1/1	-
5	PEG	C	407	-	-	3/4/4/4	-
5	PEG	A	415	-	-	2/4/4/4	-
4	EDO	A	405	-	-	0/1/1/1	-
4	EDO	D	406	-	-	1/1/1/1	-
4	EDO	B	406	-	-	1/1/1/1	-
4	EDO	A	409	-	-	1/1/1/1	-
2	NAD	A	401	-	-	1/26/62/62	0/5/5/5
4	EDO	B	409	-	-	0/1/1/1	-
4	EDO	A	407	-	-	1/1/1/1	-
4	EDO	C	406	-	-	1/1/1/1	-
5	PEG	A	417	-	-	3/4/4/4	-
4	EDO	A	412	-	-	1/1/1/1	-
4	EDO	C	405	-	-	0/1/1/1	-
4	EDO	A	410	-	-	0/1/1/1	-
4	EDO	B	407	-	-	0/1/1/1	-
4	EDO	A	404	-	-	1/1/1/1	-
4	EDO	B	405	-	-	1/1/1/1	-
4	EDO	A	403	-	-	1/1/1/1	-
4	EDO	B	412	-	-	1/1/1/1	-
4	EDO	B	410	-	-	0/1/1/1	-
4	EDO	A	406	-	-	1/1/1/1	-
4	EDO	D	404	-	-	0/1/1/1	-
2	NAD	D	401	-	-	1/26/62/62	0/5/5/5
5	PEG	A	416	-	-	1/4/4/4	-
4	EDO	B	403	-	-	1/1/1/1	-
5	PEG	A	418	-	-	1/4/4/4	-
4	EDO	C	404	-	-	1/1/1/1	-
5	PEG	A	414	-	-	2/4/4/4	-
4	EDO	A	413	-	-	1/1/1/1	-
4	EDO	B	408	-	-	1/1/1/1	-
4	EDO	D	405	-	-	1/1/1/1	-
5	PEG	B	413	-	-	1/4/4/4	-
5	PEG	B	414	-	-	2/4/4/4	-

All (25) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	401	NAD	O4B-C1B	6.46	1.50	1.41
2	D	401	NAD	C2D-C1D	-5.69	1.45	1.53
2	C	401	NAD	O4B-C1B	4.59	1.47	1.41
2	D	401	NAD	O4B-C1B	4.56	1.47	1.41
2	C	401	NAD	C2B-C1B	-4.50	1.46	1.53
2	A	401	NAD	C2A-N3A	4.33	1.39	1.32
2	B	401	NAD	C2D-C1D	-3.99	1.47	1.53
2	B	401	NAD	O4B-C1B	3.93	1.46	1.41
2	B	401	NAD	C2B-C1B	-3.41	1.48	1.53
2	C	401	NAD	C2D-C1D	-3.07	1.49	1.53
2	B	401	NAD	C2A-N3A	3.04	1.37	1.32
2	A	401	NAD	O3B-C3B	2.73	1.49	1.43
2	C	401	NAD	O4D-C1D	2.56	1.44	1.41
2	D	401	NAD	C2B-C1B	-2.39	1.50	1.53
2	C	401	NAD	C5A-N7A	-2.37	1.31	1.39
2	C	401	NAD	C5A-C4A	2.36	1.47	1.40
2	C	401	NAD	C2N-N1N	-2.36	1.32	1.35
2	B	401	NAD	C2N-N1N	2.31	1.37	1.35
2	A	401	NAD	C7N-N7N	-2.27	1.28	1.33
2	C	401	NAD	C3N-C7N	2.12	1.53	1.50
2	A	401	NAD	O3D-C3D	-2.11	1.38	1.43
2	C	401	NAD	C5N-C4N	-2.08	1.34	1.38
2	D	401	NAD	PN-O1N	-2.08	1.43	1.50
2	B	401	NAD	O4D-C4D	-2.06	1.40	1.45
2	A	401	NAD	C3N-C7N	2.06	1.53	1.50

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	401	NAD	N3A-C2A-N1A	-6.13	119.10	128.68
2	D	401	NAD	C5A-C6A-N6A	5.17	128.21	120.35
2	A	401	NAD	N3A-C2A-N1A	-4.57	121.54	128.68
2	B	401	NAD	O7N-C7N-N7N	-4.45	116.26	122.58
2	A	401	NAD	O4B-C1B-C2B	-4.44	100.44	106.93
2	B	401	NAD	O7N-C7N-C3N	3.99	124.41	119.63
2	B	401	NAD	C6N-N1N-C2N	-3.66	118.64	121.97
2	D	401	NAD	O7N-C7N-N7N	-3.37	117.79	122.58
2	D	401	NAD	C3N-C7N-N7N	3.25	121.65	117.75
2	A	401	NAD	C3N-C7N-N7N	3.06	121.43	117.75
2	A	401	NAD	C6N-N1N-C2N	-2.87	119.35	121.97
5	B	413	PEG	O2-C2-C1	2.71	121.98	110.07
2	A	401	NAD	O4D-C4D-C3D	2.63	110.31	105.11
2	A	401	NAD	O2B-C2B-C3B	2.24	119.07	111.82

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	401	NAD	C5N-C4N-C3N	-2.21	117.73	120.34
2	D	401	NAD	N6A-C6A-N1A	-2.14	114.13	118.57
2	B	401	NAD	C3N-C2N-N1N	2.11	122.49	120.43
2	B	401	NAD	C3B-C2B-C1B	2.08	104.11	100.98
2	C	401	NAD	C5A-C6A-N6A	2.08	123.51	120.35
2	A	401	NAD	O7N-C7N-N7N	-2.02	119.70	122.58

There are no chirality outliers.

All (46) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	401	NAD	O4D-C1D-N1N-C6N
2	B	401	NAD	O4D-C1D-N1N-C6N
2	C	401	NAD	O4D-C1D-N1N-C6N
2	D	401	NAD	O4D-C1D-N1N-C6N
5	B	413	PEG	C1-C2-O2-C3
5	A	418	PEG	C4-C3-O2-C2
5	A	415	PEG	O2-C3-C4-O4
4	A	407	EDO	O1-C1-C2-O2
4	B	411	EDO	O1-C1-C2-O2
4	C	406	EDO	O1-C1-C2-O2
4	D	405	EDO	O1-C1-C2-O2
5	A	419	PEG	O1-C1-C2-O2
5	A	417	PEG	O2-C3-C4-O4
5	C	407	PEG	O1-C1-C2-O2
5	B	414	PEG	O2-C3-C4-O4
4	A	411	EDO	O1-C1-C2-O2
4	B	412	EDO	O1-C1-C2-O2
4	A	406	EDO	O1-C1-C2-O2
4	A	409	EDO	O1-C1-C2-O2
4	B	408	EDO	O1-C1-C2-O2
4	D	406	EDO	O1-C1-C2-O2
5	B	414	PEG	C1-C2-O2-C3
5	A	419	PEG	C4-C3-O2-C2
5	C	407	PEG	C4-C3-O2-C2
5	A	415	PEG	C4-C3-O2-C2
5	A	417	PEG	C4-C3-O2-C2
5	A	419	PEG	O2-C3-C4-O4
4	A	408	EDO	O1-C1-C2-O2
4	B	406	EDO	O1-C1-C2-O2
5	A	419	PEG	C1-C2-O2-C3
5	A	414	PEG	C1-C2-O2-C3

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Mol	Chain	Res	Type	Atoms
4	A	404	EDO	O1-C1-C2-O2
4	D	407	EDO	O1-C1-C2-O2
5	A	417	PEG	C1-C2-O2-C3
2	B	401	NAD	C3B-C4B-C5B-O5B
5	A	414	PEG	C4-C3-O2-C2
4	A	403	EDO	O1-C1-C2-O2
4	B	403	EDO	O1-C1-C2-O2
5	A	416	PEG	O2-C3-C4-O4
5	C	407	PEG	C1-C2-O2-C3
2	B	401	NAD	O4B-C4B-C5B-O5B
4	A	412	EDO	O1-C1-C2-O2
4	A	413	EDO	O1-C1-C2-O2
4	C	404	EDO	O1-C1-C2-O2
2	C	401	NAD	C2D-C1D-N1N-C2N
4	B	405	EDO	O1-C1-C2-O2

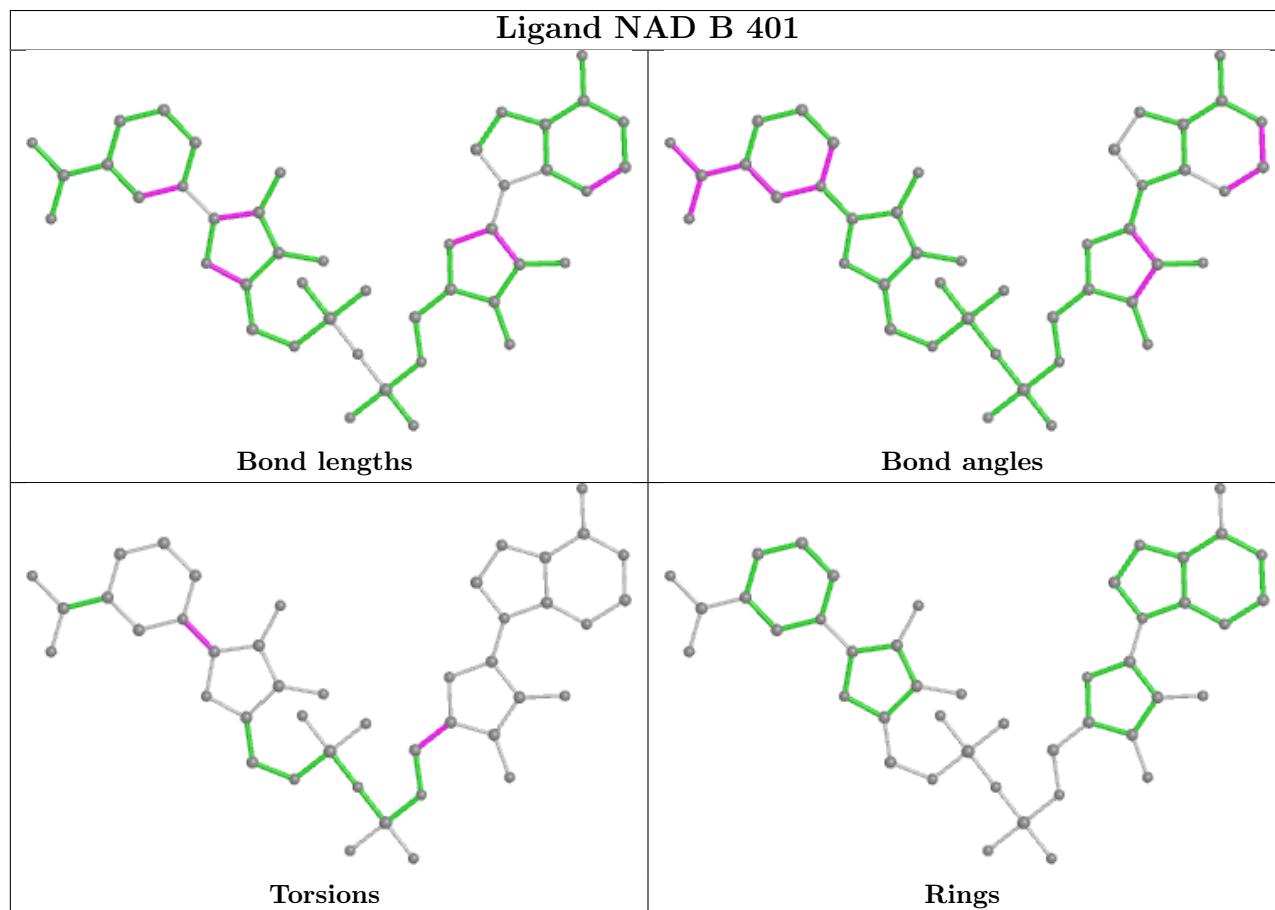
There are no ring outliers.

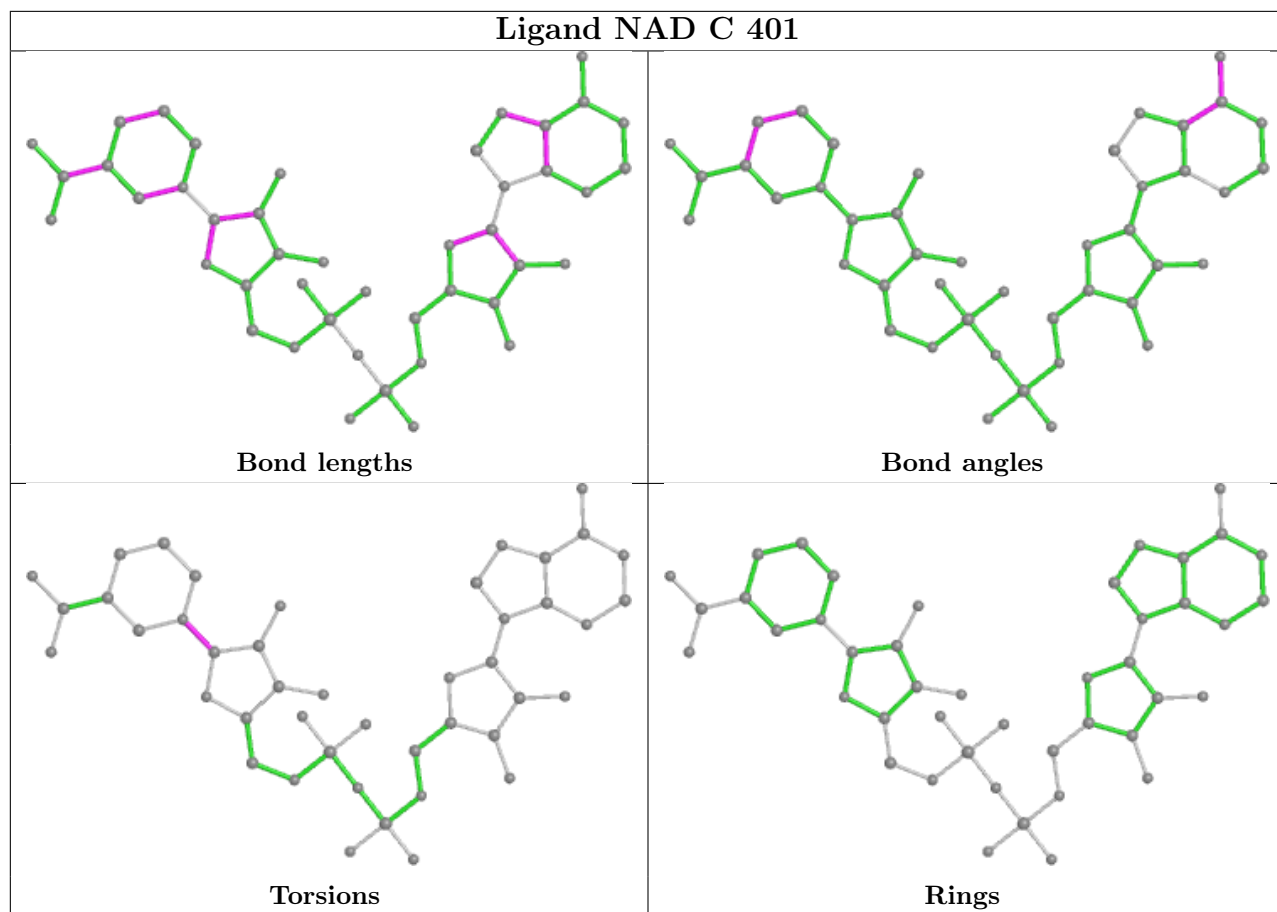
18 monomers are involved in 39 short contacts:

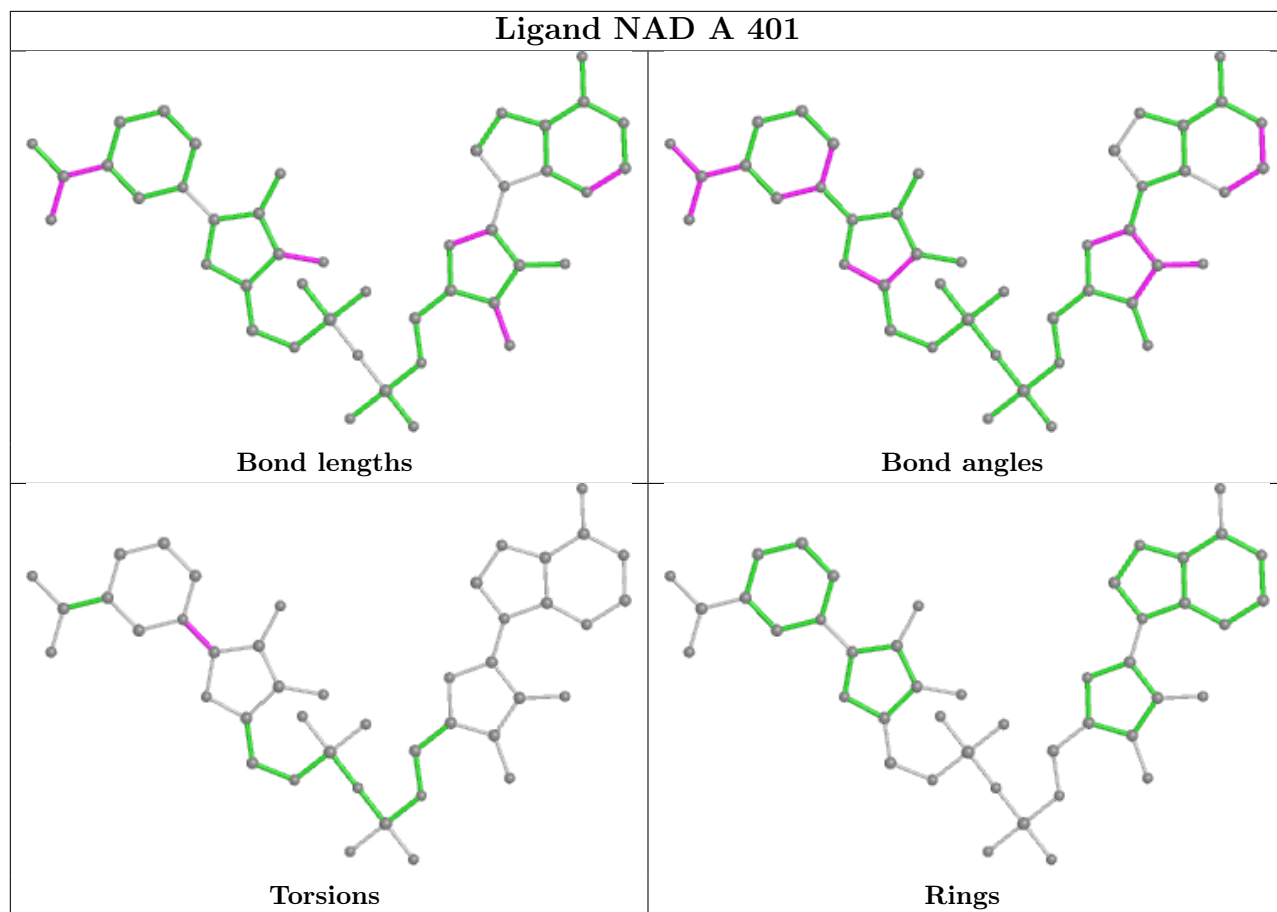
Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	D	407	EDO	1	0
5	A	419	PEG	2	0
4	D	403	EDO	2	0
4	B	411	EDO	4	0
5	C	407	PEG	2	0
5	A	415	PEG	5	0
4	C	406	EDO	3	0
5	A	417	PEG	1	0
4	B	407	EDO	1	0
4	A	406	EDO	2	0
4	D	404	EDO	2	0
5	A	416	PEG	1	0
5	A	418	PEG	2	0
4	C	404	EDO	2	0
5	A	414	PEG	1	0
4	A	413	EDO	2	0
5	B	413	PEG	5	0
5	B	414	PEG	1	0

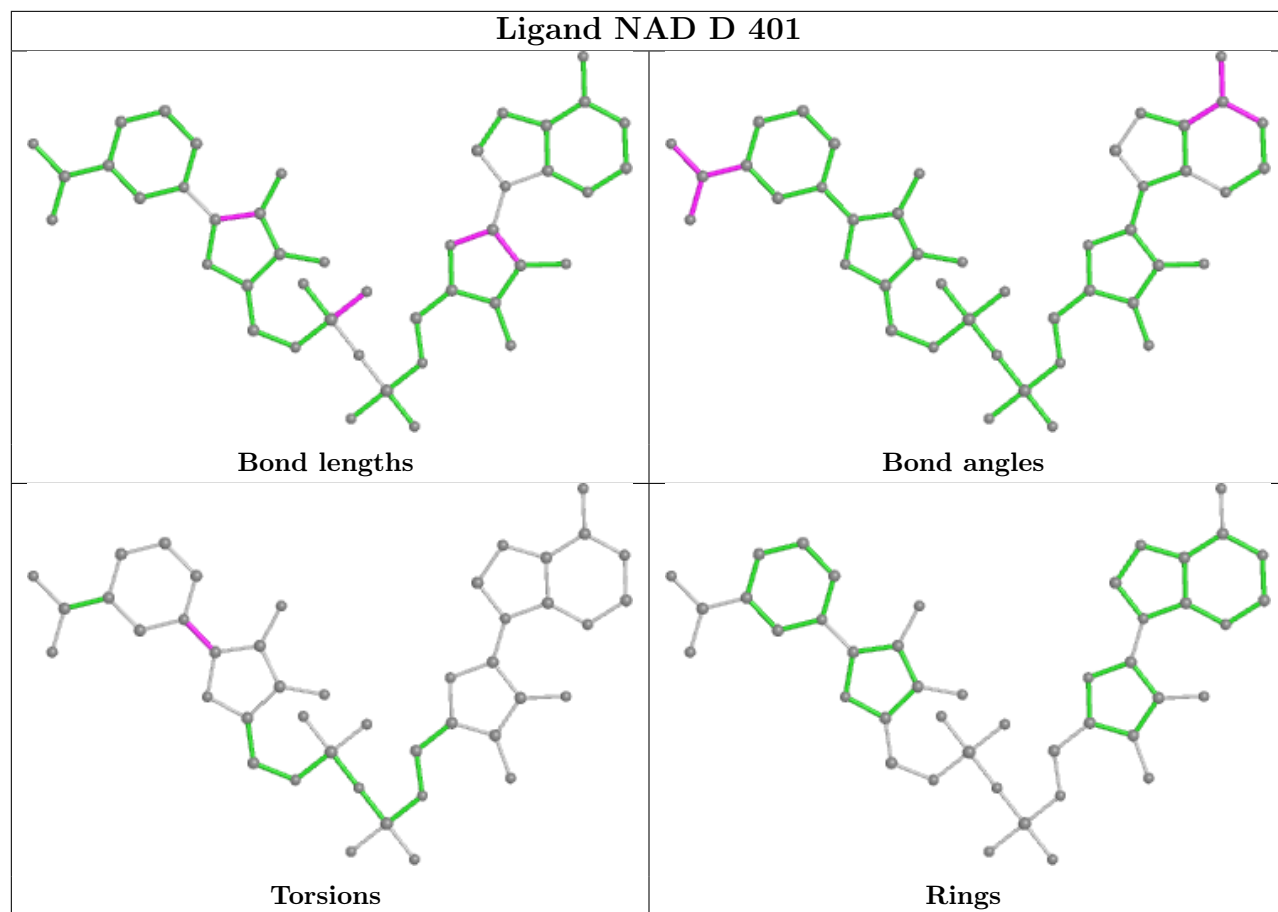
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will

also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.









## 5.7 Other polymers [i](#)

There are no such residues in this entry.

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

There are no chain breaks in this entry.



## 6 Fit of model and data [i](#)

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

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 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	362/375 (96%)	0.17	14 (3%) 39 41	20, 28, 51, 114	0
1	B	363/375 (96%)	0.14	11 (3%) 50 51	18, 27, 52, 106	0
1	C	354/375 (94%)	0.65	42 (11%) 4 4	21, 51, 101, 130	0
1	D	353/375 (94%)	0.40	21 (5%) 22 24	22, 43, 80, 110	0
All	All	1432/1500 (95%)	0.34	88 (6%) 21 22	18, 35, 87, 130	0

All (88) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	362	PRO	10.6
1	A	235	LYS	8.5
1	C	361	ILE	7.5
1	A	260	TYR	6.8
1	A	237	THR	6.3
1	D	59	VAL	6.2
1	C	353	VAL	5.4
1	C	283	ILE	5.4
1	C	237	THR	5.3
1	B	284	GLY	5.1
1	D	361	ILE	5.0
1	C	234	GLU	5.0
1	C	59	VAL	4.9
1	B	236	GLU	4.8
1	A	239	GLY	4.7
1	A	238	SER	4.4
1	C	286	PHE	4.2
1	C	239	GLY	4.1
1	B	341	HIS	4.1
1	C	358	SER	4.0
1	C	309	ALA	4.0

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>	<b>RSRZ</b>
1	C	9	ASP	4.0
1	D	25	GLN	3.9
1	A	24	GLN	3.9
1	C	347	ASP	3.8
1	B	237	THR	3.8
1	C	360	CYS	3.8
1	D	10	GLU	3.8
1	B	0	GLY	3.7
1	A	283	ILE	3.6
1	C	274	VAL	3.6
1	C	11	ARG	3.6
1	D	234	GLU	3.5
1	D	341	HIS	3.5
1	C	311	MET	3.5
1	B	238	SER	3.4
1	C	310	LYS	3.4
1	D	237	THR	3.4
1	D	9	ASP	3.4
1	C	341	HIS	3.3
1	C	235	LYS	3.3
1	B	239	GLY	3.2
1	C	250	PHE	3.2
1	A	341	HIS	3.2
1	D	60	SER	3.1
1	C	284	GLY	3.1
1	A	234	GLU	3.1
1	C	279	THR	3.0
1	C	263	TRP	3.0
1	D	11	ARG	3.0
1	A	59	VAL	2.9
1	C	236	GLU	2.9
1	D	349	ASP	2.9
1	B	20	SER	2.9
1	A	284	GLY	2.9
1	D	239	GLY	2.9
1	B	59	VAL	2.8
1	D	235	LYS	2.7
1	D	355	ALA	2.7
1	C	0	GLY	2.7
1	C	348	VAL	2.7
1	C	317	LEU	2.7
1	C	352	THR	2.6

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Mol	Chain	Res	Type	RSRZ
1	C	330	LYS	2.6
1	D	8	LEU	2.5
1	C	354	THR	2.5
1	B	283	ILE	2.5
1	D	358	SER	2.4
1	D	24	GLN	2.4
1	C	315	ASP	2.4
1	C	198	SER	2.3
1	D	1	MET	2.3
1	D	260	TYR	2.3
1	A	361	ILE	2.3
1	C	20	SER	2.3
1	C	357	ILE	2.2
1	C	331	LEU	2.2
1	C	350	ALA	2.2
1	C	298	LEU	2.2
1	B	1	MET	2.2
1	C	316	TYR	2.2
1	C	282	LEU	2.1
1	A	347[A]	ASP	2.1
1	C	287	THR	2.0
1	C	312	GLN	2.0
1	C	292	ASP	2.0
1	A	20	SER	2.0
1	D	351	SER	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, 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.

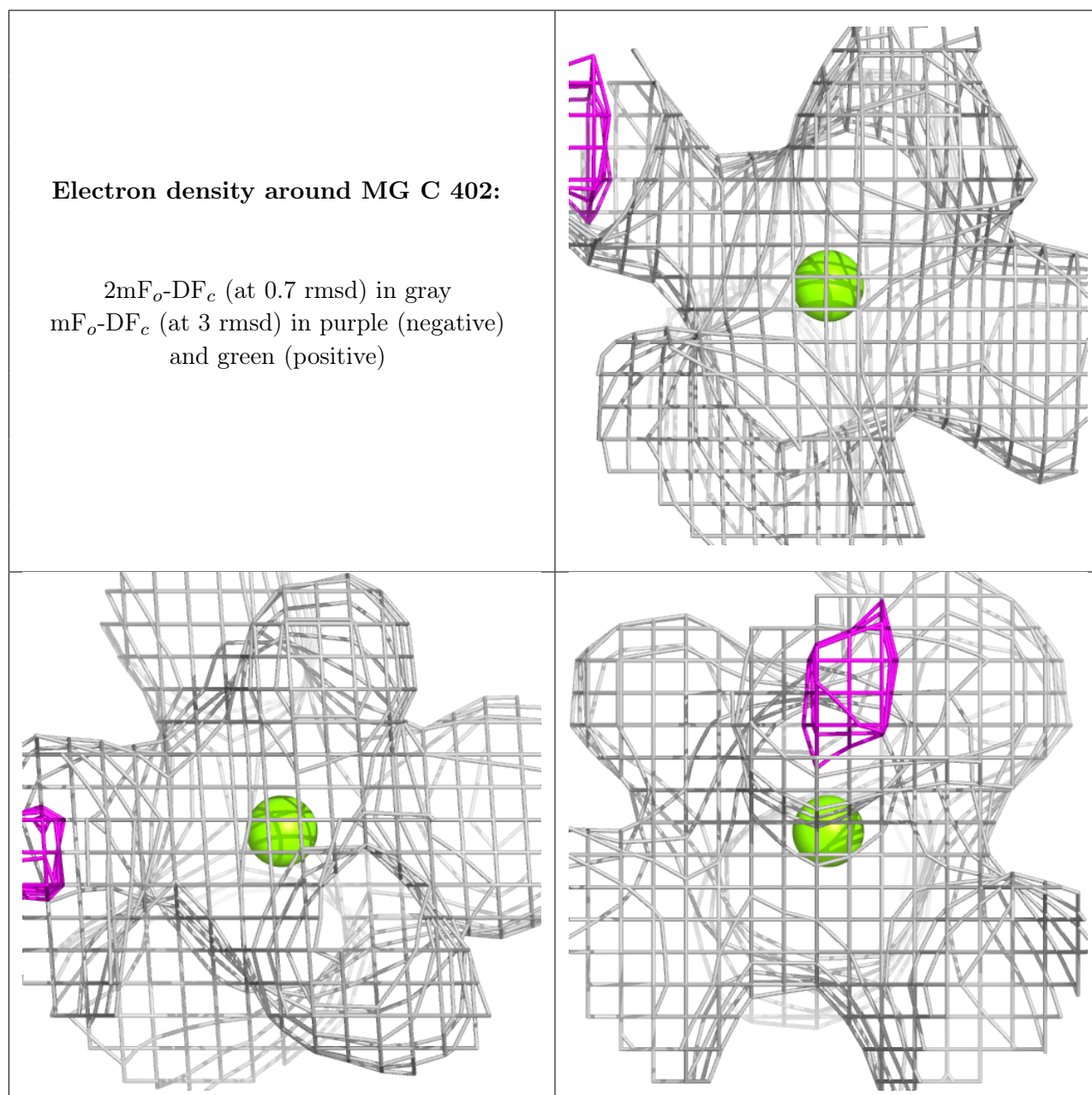
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å <sup>2</sup> )	Q<0.9
4	EDO	C	403	4/4	0.55	0.43	65,72,81,84	0
4	EDO	A	413	4/4	0.66	0.19	58,64,76,77	0
5	PEG	A	416	7/7	0.72	0.26	58,64,75,79	0
4	EDO	B	409	4/4	0.75	0.23	66,69,77,78	0
4	EDO	A	405	4/4	0.76	0.20	54,59,60,61	0
4	EDO	B	410	4/4	0.78	0.23	54,58,70,93	0
5	PEG	A	417	7/7	0.78	0.14	60,63,66,74	0
4	EDO	D	407	4/4	0.80	0.15	57,64,67,67	0
4	EDO	D	406	4/4	0.81	0.12	67,68,68,70	0
4	EDO	B	411	4/4	0.82	0.23	36,38,57,68	0
4	EDO	A	412	4/4	0.83	0.15	72,75,78,80	0
4	EDO	A	406	4/4	0.83	0.23	50,52,60,63	0
5	PEG	A	419	7/7	0.84	0.17	36,55,67,77	0
5	PEG	B	414	7/7	0.84	0.14	49,63,71,72	0
5	PEG	C	407	7/7	0.84	0.16	43,60,73,88	0
5	PEG	A	415	7/7	0.85	0.20	33,49,64,66	0
4	EDO	C	406	4/4	0.85	0.23	44,52,56,58	0
4	EDO	A	408	4/4	0.86	0.16	39,48,55,59	0
4	EDO	A	407	4/4	0.86	0.19	45,47,49,52	0
4	EDO	D	408	4/4	0.87	0.14	49,53,57,60	0
5	PEG	A	414	7/7	0.87	0.22	49,51,85,87	0
4	EDO	B	406	4/4	0.88	0.11	51,53,57,57	0
4	EDO	C	404	4/4	0.89	0.18	44,50,57,62	0
4	EDO	B	405	4/4	0.89	0.19	42,43,51,51	0
4	EDO	A	411	4/4	0.89	0.10	49,55,65,66	0
4	EDO	D	405	4/4	0.90	0.20	51,57,59,73	0
4	EDO	D	403	4/4	0.90	0.14	40,45,46,50	0
5	PEG	A	418	7/7	0.90	0.10	39,51,56,57	0
4	EDO	B	404	4/4	0.91	0.12	40,45,50,50	0
4	EDO	A	410	4/4	0.91	0.09	57,57,60,63	0
4	EDO	B	408	4/4	0.92	0.12	40,47,52,65	0
4	EDO	A	403	4/4	0.92	0.16	29,32,41,45	0
4	EDO	A	404	4/4	0.93	0.14	33,44,48,49	0
4	EDO	B	403	4/4	0.94	0.12	44,59,63,68	0
4	EDO	B	412	4/4	0.94	0.15	38,38,47,54	0
4	EDO	D	404	4/4	0.94	0.17	43,43,52,54	0
4	EDO	A	409	4/4	0.94	0.14	36,43,52,59	0
4	EDO	B	407	4/4	0.94	0.16	44,47,64,69	0
5	PEG	B	413	7/7	0.95	0.19	28,39,43,50	0
3	MG	C	402	1/1	0.96	0.05	41,41,41,41	0
2	NAD	A	401	44/44	0.96	0.11	17,26,33,37	0
2	NAD	B	401	44/44	0.96	0.10	18,25,29,33	0
4	EDO	C	405	4/4	0.96	0.11	58,62,69,71	0

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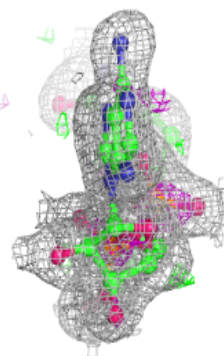
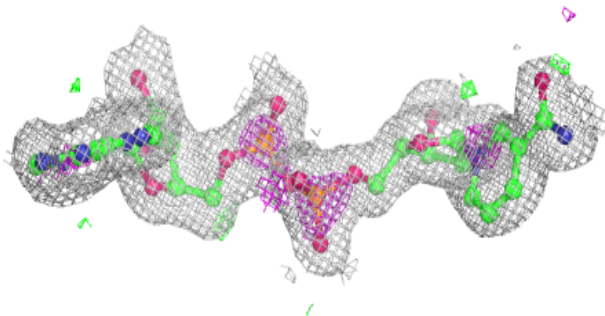
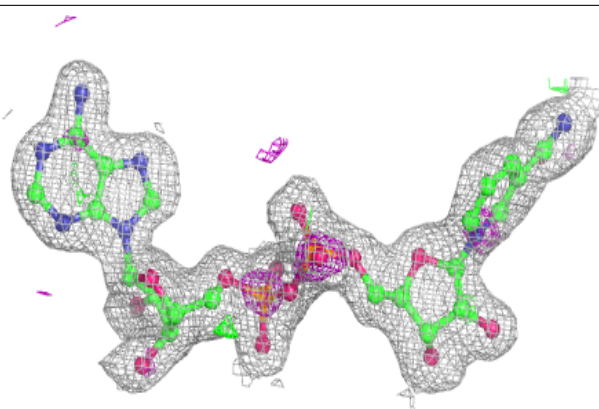
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
2	NAD	C	401	44/44	0.97	0.10	25,32,40,44	0
3	MG	D	402	1/1	0.97	0.06	28,28,28,28	0
2	NAD	D	401	44/44	0.98	0.10	25,28,33,39	0
3	MG	A	402	1/1	0.98	0.10	20,20,20,20	0
3	MG	B	402	1/1	0.98	0.09	21,21,21,21	0

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.

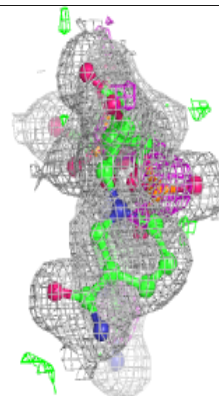
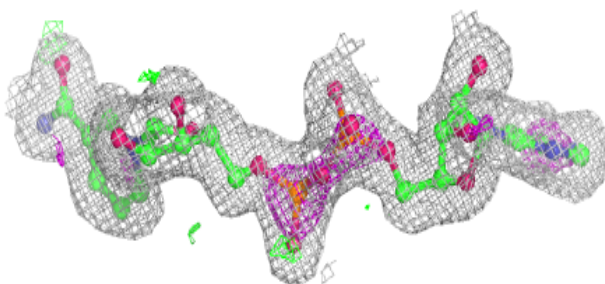
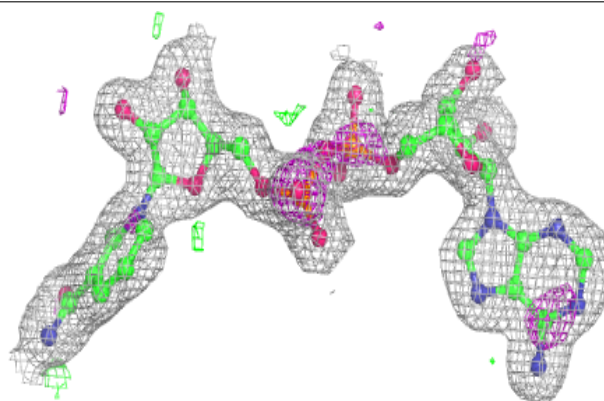


**Electron density around NAD A 401:**

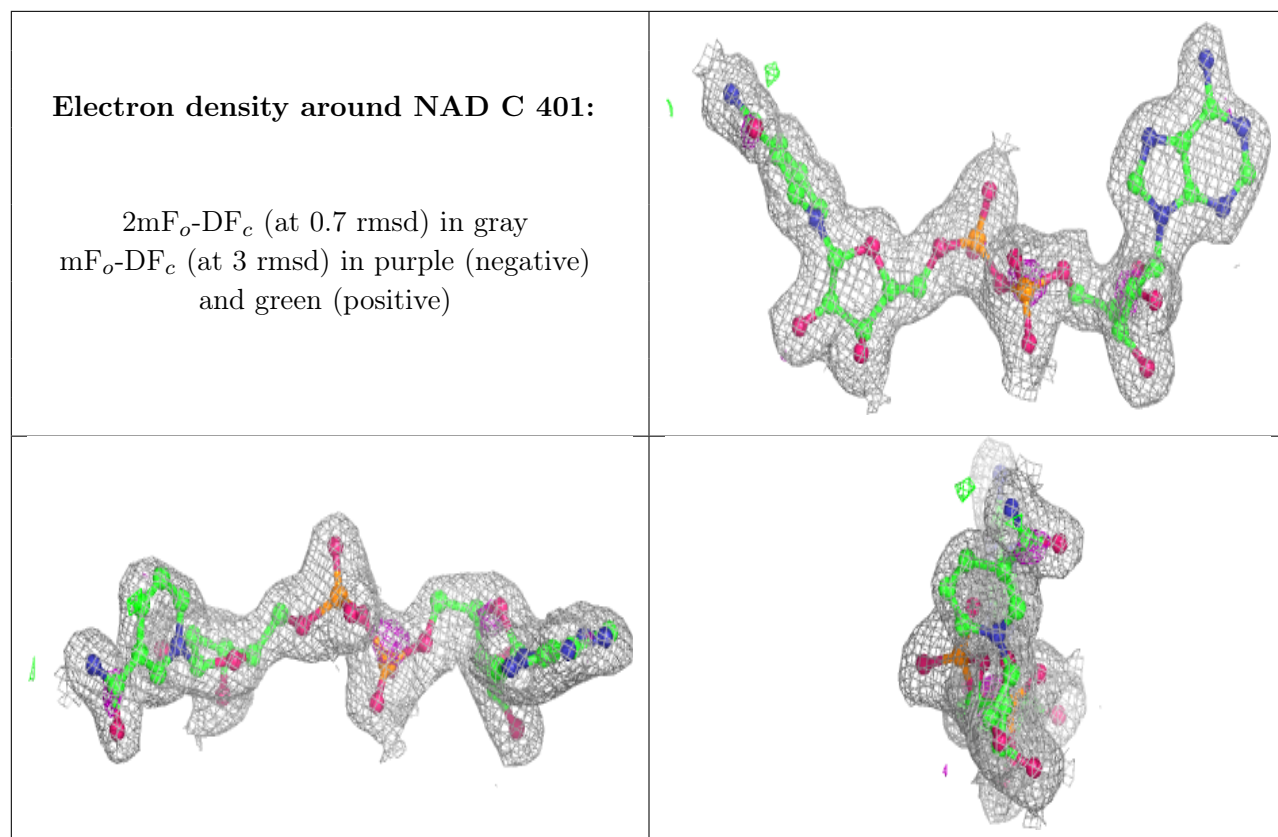
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

**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)

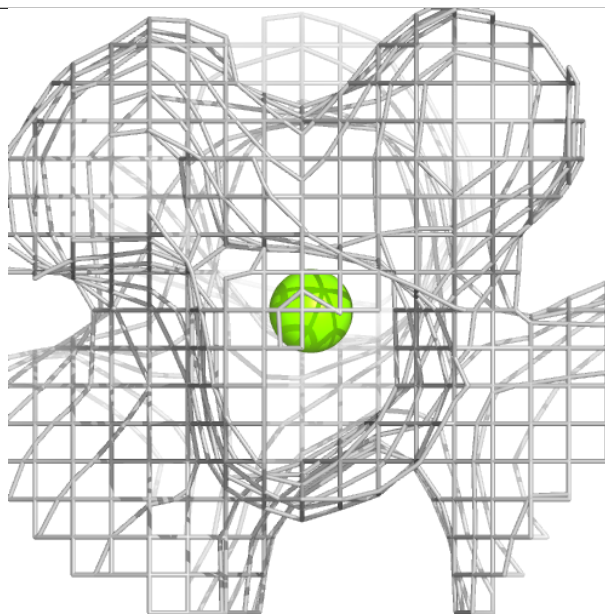
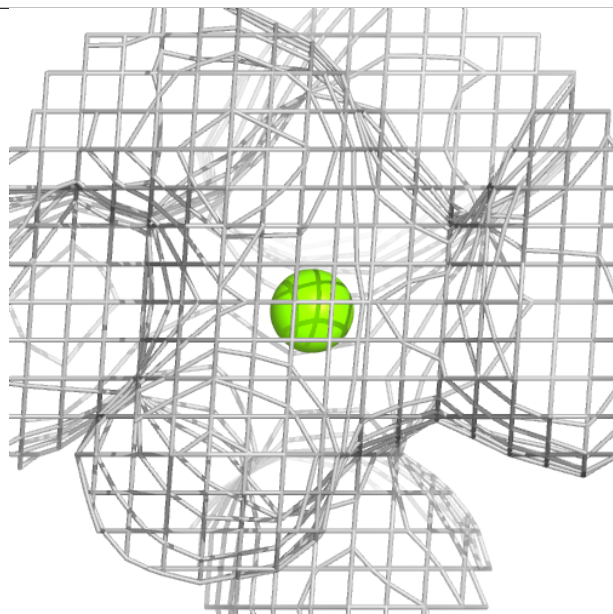
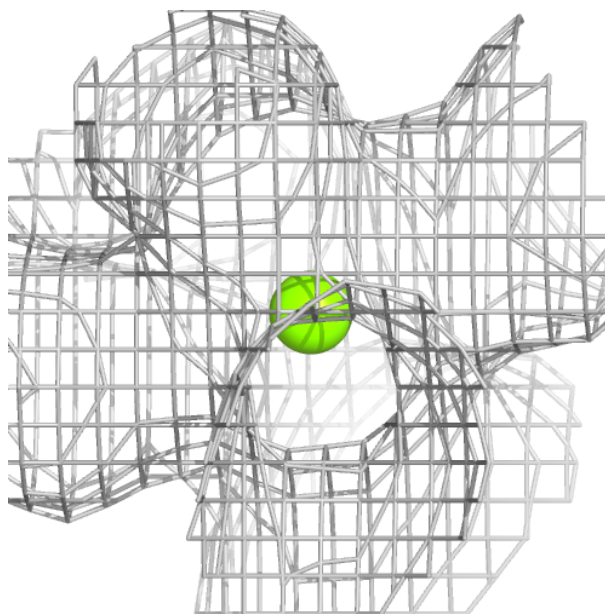




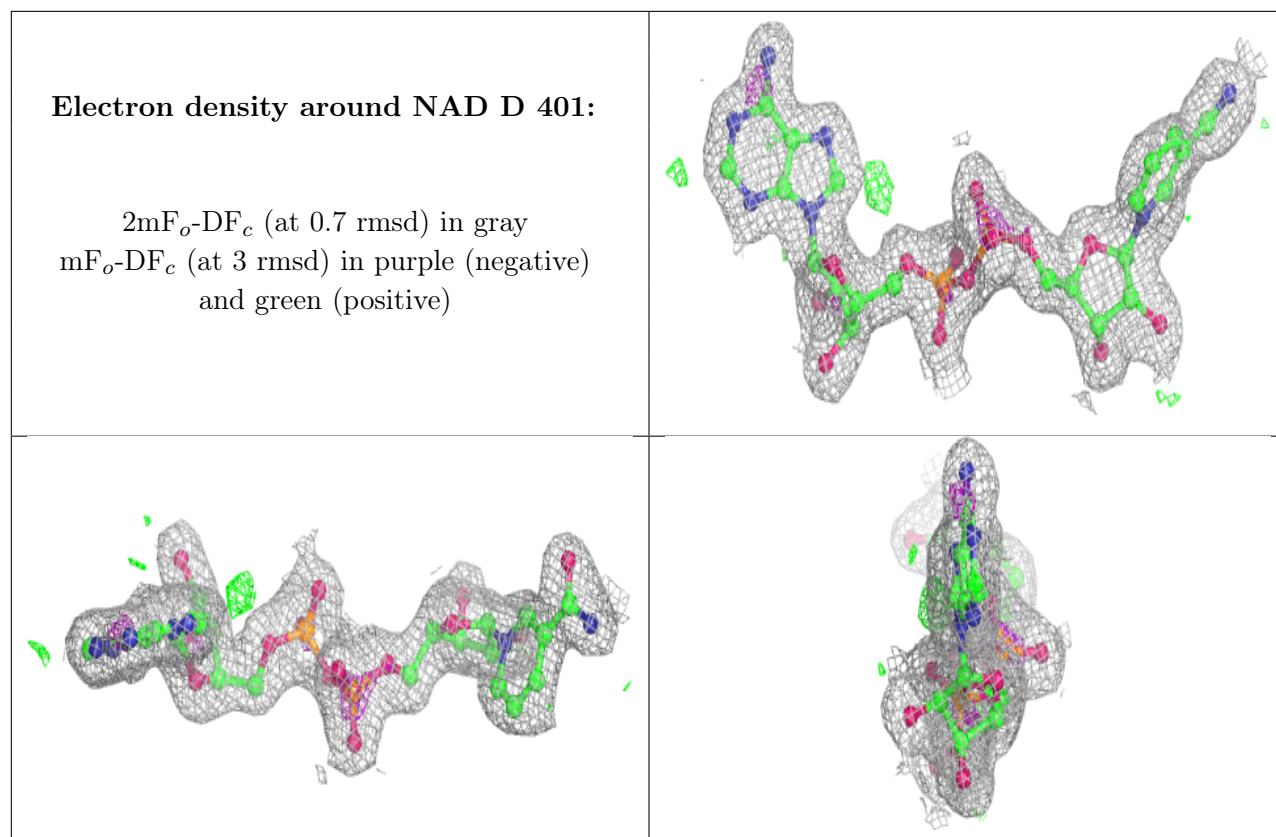


**Electron density around MG D 402:**

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

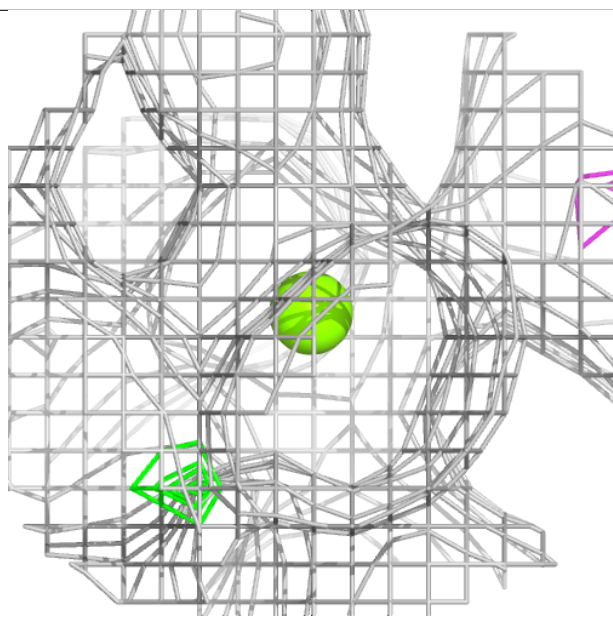
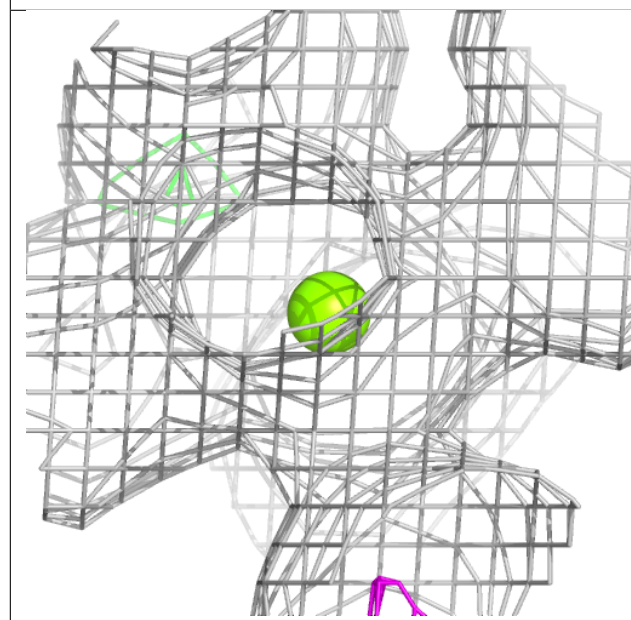
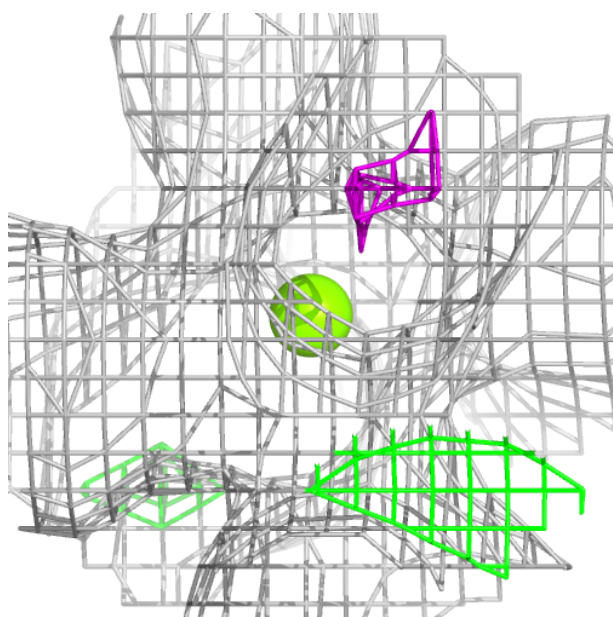


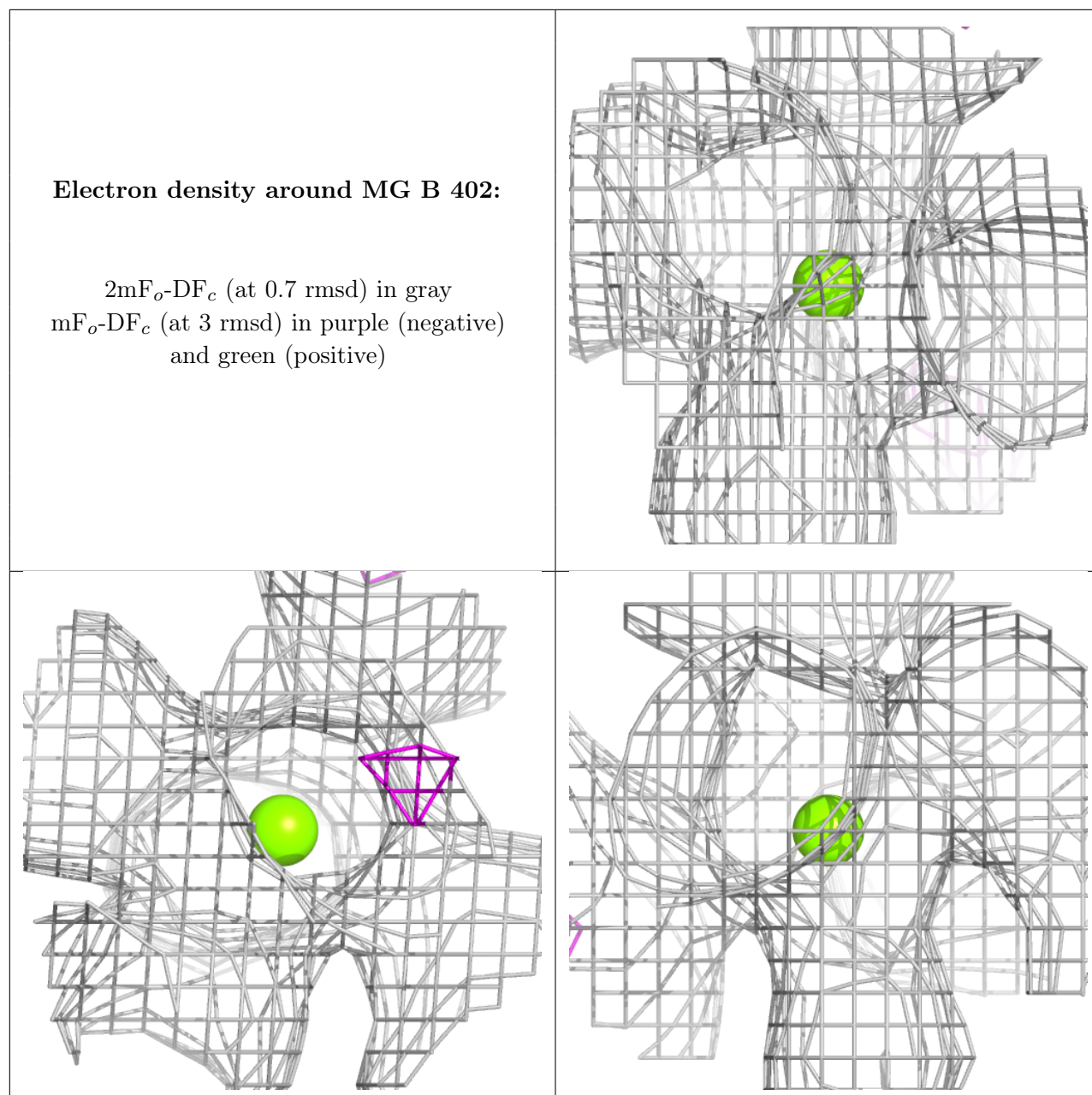




**Electron density around MG A 402:**

$2mF_o-DF_c$  (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.