



# Full wwPDB X-ray Structure Validation Report ⓘ

Jul 25, 2022 – 12:41 PM EDT

PDB ID : 7MYA  
Title : Structure of proline utilization A with the FAD covalently-modified by 1,3-dithiolane  
Authors : Tanner, J.J.; Campbell, A.C.  
Deposited on : 2021-05-20  
Resolution : 1.56 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	:	2.29
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

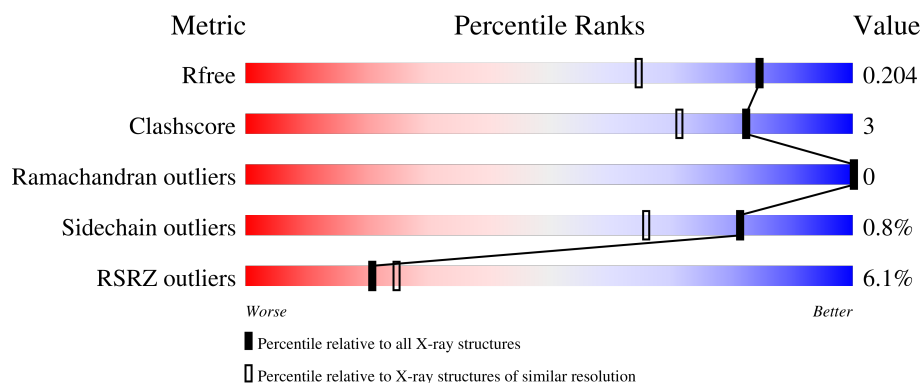
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 1.56 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	1483 (1.56-1.56)
Clashscore	141614	1529 (1.56-1.56)
Ramachandran outliers	138981	1498 (1.56-1.56)
Sidechain outliers	138945	1495 (1.56-1.56)
RSRZ outliers	127900	1465 (1.56-1.56)

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	1235	<div> <div>6%</div> <div> <div></div> <div>92%</div> <div>6%</div> </div> </div>
1	B	1235	<div> <div>6%</div> <div> <div></div> <div>92%</div> <div>7%</div> </div> </div>

## 2 Entry composition [i](#)

There are 9 unique types of molecules in this entry. The entry contains 20344 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.

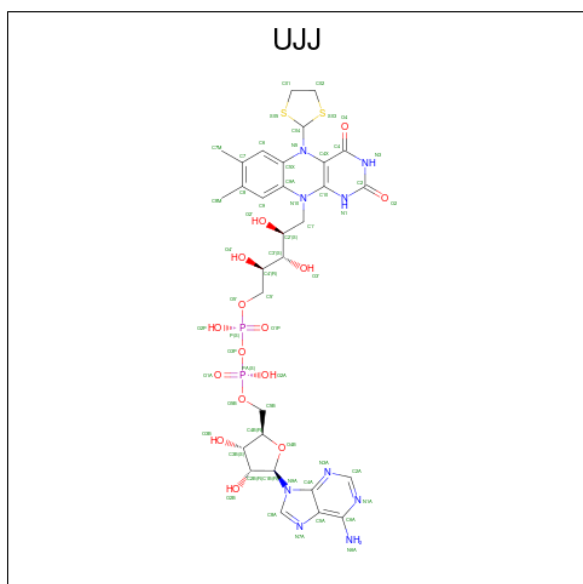
- Molecule 1 is a protein called Bifunctional protein PutA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	1217	Total	C	N	O	S	0	17	0
			9063	5713	1615	1700	35			
1	B	1217	Total	C	N	O	S	0	15	0
			9042	5698	1616	1695	33			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	SER	-	expression tag	UNP F7X6I3
A	0	MET	-	expression tag	UNP F7X6I3
B	-1	SER	-	expression tag	UNP F7X6I3
B	0	MET	-	expression tag	UNP F7X6I3

- Molecule 2 is [(2R,3S,4R,5R)-5-(6-amino-9H-purin-9-yl)-3,4-dihydroxytetrahydrofuran-2-yl] methyl (2R,3S,4S)-5-[5-(1,3-dithiolan-2-yl)-7,8-dimethyl-2,4-dioxo-1,3,4,5-tetrahydrobenzo[g]pteridin-10(2H)-yl]-2,3,4-trihydroxypentyl dihydrogen diphosphate (three-letter code: UJJ) (formula: C<sub>30</sub>H<sub>39</sub>N<sub>9</sub>O<sub>15</sub>P<sub>2</sub>S<sub>2</sub>) (labeled as "Ligand of Interest" by depositor).

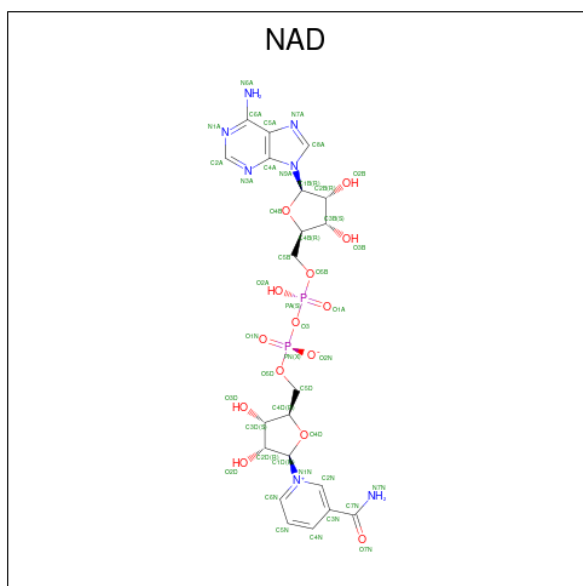


Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	S	0	0
			58	30	9	15	2	2		
2	B	1	Total	C	N	O	P	S	0	0
			58	30	9	15	2	2		

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

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

- Molecule 4 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
4	A	1	Total	C	N	O	P	0	0
			44	21	7	14	2		
4	B	1	Total	C	N	O	P	0	0
			44	21	7	14	2		

- Molecule 5 is SULFATE ION (three-letter code: SO4) (formula: O<sub>4</sub>S).



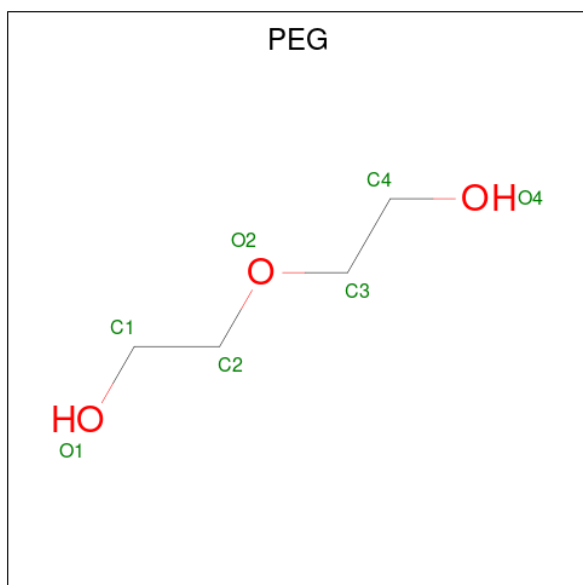
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	O	S	0	0
			5	4	1		
5	A	1	Total	O	S	0	0
			5	4	1		
5	A	1	Total	O	S	0	0
			5	4	1		
5	A	1	Total	O	S	0	0
			5	4	1		
5	B	1	Total	O	S	0	0
			5	4	1		
5	B	1	Total	O	S	0	0
			5	4	1		

- Molecule 6 is TRIETHYLENE GLYCOL (three-letter code: PGE) (formula: C<sub>6</sub>H<sub>14</sub>O<sub>4</sub>).



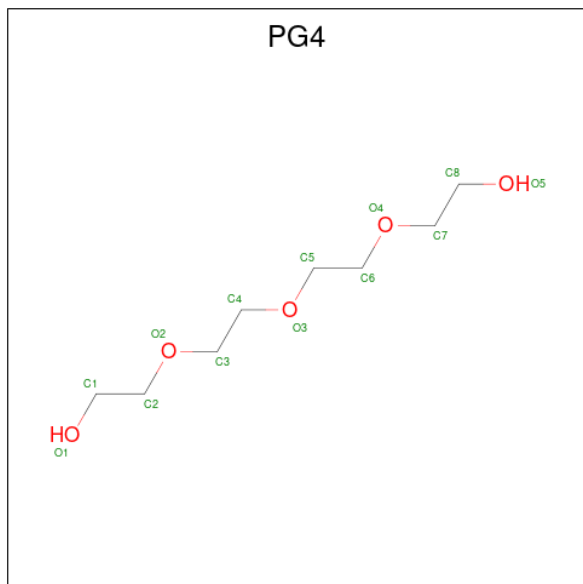
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	A	1	Total	C	O	0	0
			10	6	4		
6	B	1	Total	C	O	0	0
			10	6	4		
6	B	1	Total	C	O	0	0
			10	6	4		
6	B	1	Total	C	O	0	0
			10	6	4		

- Molecule 7 is DI(HYDROXYETHYL)ETHER (three-letter code: PEG) (formula:  $C_4H_{10}O_3$ ).



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

- Molecule 8 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
8	B	1	Total	C	O	0	0
			13	8	5		

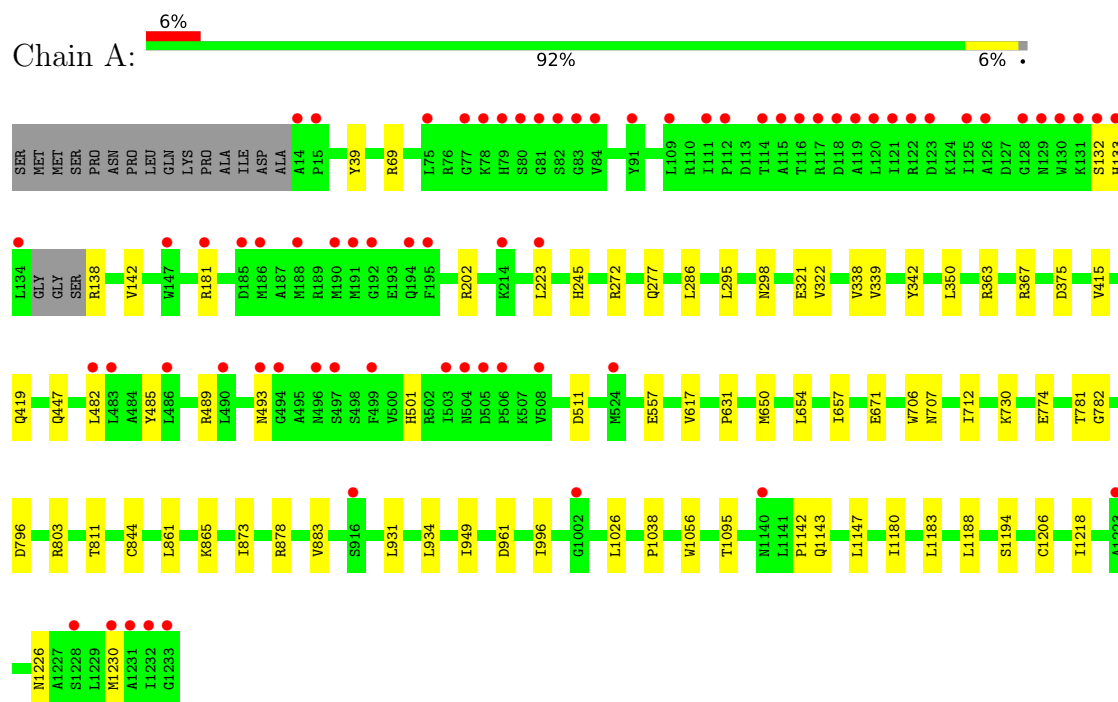
- Molecule 9 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
9	A	983	Total	O	0	0
			983	983		
9	B	941	Total	O	0	1
			941	941		

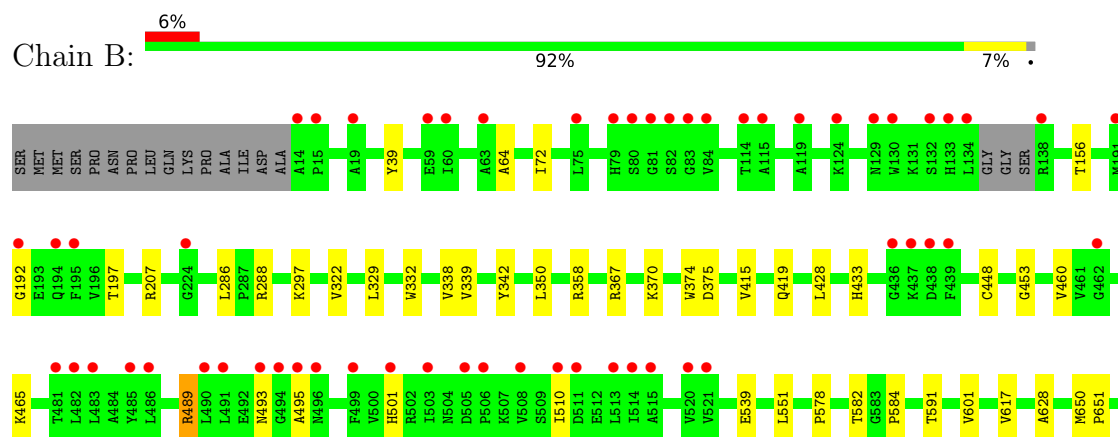
### 3 Residue-property plots [i](#)

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.

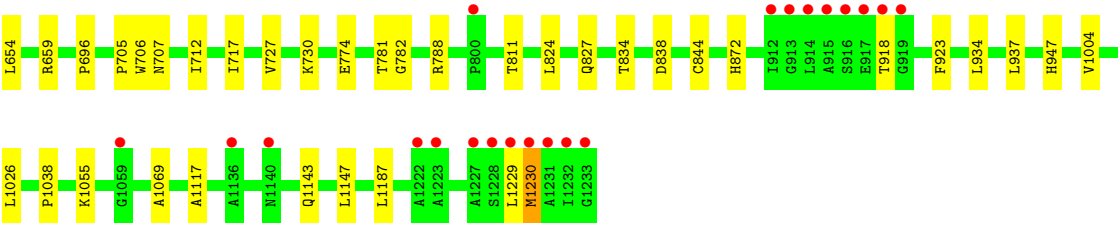
#### • Molecule 1: Bifunctional protein PutA



#### • Molecule 1: Bifunctional protein PutA







## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	101.18Å 101.88Å 125.93Å 90.00° 106.49° 90.00°	Depositor
Resolution (Å)	46.94 – 1.56 46.94 – 1.56	Depositor EDS
% Data completeness (in resolution range)	98.5 (46.94-1.56) 98.5 (46.94-1.56)	Depositor EDS
$R_{merge}$	0.09	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.48 (at 1.56Å)	Xtriage
Refinement program	PHENIX 1.19.2	Depositor
R, $R_{free}$	0.176 , 0.205 0.175 , 0.204	Depositor DCC
$R_{free}$ test set	17139 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	18.7	Xtriage
Anisotropy	0.153	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.35 , 47.0	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.51$ , $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.97	EDS
Total number of atoms	20344	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	25.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.11% 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

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: PGE, UJJ, PG4, MG, PEG, NAD, SO4

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	0.33	0/9260	0.58	0/12604
1	B	0.33	0/9251	0.59	0/12595
All	All	0.33	0/18511	0.58	0/25199

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

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	9063	0	9093	43	0
1	B	9042	0	9076	51	0
2	A	58	0	0	0	0
2	B	58	0	0	1	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
4	A	44	0	26	1	0
4	B	44	0	26	2	0
5	A	25	0	0	0	0
5	B	10	0	0	0	0
6	A	10	0	14	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
6	B	30	0	42	4	0
7	B	21	0	30	1	0
8	B	13	0	18	0	0
9	A	983	0	0	8	0
9	B	941	0	0	5	0
All	All	20344	0	18325	93	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 (93) 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:298:ASN:ND2	9:A:2302:HOH:O	2.15	0.74
1:B:582:THR:HA	6:B:1308:PGE:H2	1.71	0.73
1:A:1183:LEU:O	9:A:2301:HOH:O	2.07	0.72
1:B:539:GLU:OE1	9:B:2401:HOH:O	2.09	0.71
1:B:844:CYS:SG	4:B:1303:NAD:C4N	2.82	0.66
1:B:286:LEU:HD21	1:B:322:VAL:HG11	1.78	0.66
1:B:788[B]:ARG:HD2	1:B:1187:LEU:HD11	1.78	0.66
1:B:339[B]:VAL:HG21	1:B:350:LEU:HD21	1.77	0.65
1:A:1206:CYS:SG	9:A:3228:HOH:O	2.54	0.65
1:A:286:LEU:HD21	1:A:322:VAL:HG11	1.80	0.64
1:A:1095[B]:THR:HG22	9:A:2886:HOH:O	1.98	0.64
1:A:861:LEU:HD22	1:A:865:LYS:HE3	1.80	0.63
1:A:138:ARG:HE	1:A:142[A]:VAL:HG21	1.65	0.62
1:A:557:GLU:OE1	9:A:2303:HOH:O	2.16	0.61
1:B:650:MET:O	1:B:654:LEU:HG	2.01	0.60
1:A:873:ILE:HG13	1:A:883:VAL:HB	1.84	0.59
1:B:358:ARG:HG2	1:B:415:VAL:HG11	1.84	0.59
1:A:996[B]:ILE:HD12	1:A:1218:ILE:HG12	1.85	0.58
1:A:69:ARG:NH2	1:A:511:ASP:OD1	2.37	0.57
1:B:824:LEU:HD23	1:B:827:GLN:HG3	1.87	0.55
1:B:297:LYS:HD2	1:B:329:LEU:HA	1.87	0.55
1:A:844:CYS:SG	4:A:2203:NAD:C4N	2.95	0.55
1:B:937:LEU:HD21	1:B:947:HIS:CD2	2.42	0.55
1:A:1180:ILE:HG23	1:A:1188:LEU:HD12	1.90	0.54
1:A:650:MET:O	1:A:654:LEU:HG	2.08	0.53
1:B:551:LEU:HD22	6:B:1311:PGE:H42	1.90	0.53
1:B:288[B]:ARG:NE	9:B:2409:HOH:O	2.41	0.52
1:A:961:ASP:OD2	1:B:1055:LYS:NZ	2.34	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:156:THR:OG1	9:B:2402:HOH:O	2.17	0.51
1:A:367[A]:ARG:NE	1:A:447:GLN:OE1	2.44	0.51
1:A:245:HIS:CE1	1:A:295:LEU:HD11	2.46	0.50
1:A:712:ILE:HD13	1:A:781:THR:HG21	1.92	0.50
1:A:339:VAL:HG21	1:A:350:LEU:HD21	1.92	0.50
1:A:1226:ASN:O	1:A:1230:MET:HG2	2.11	0.50
1:A:338:VAL:HG22	1:A:367[B]:ARG:HD2	1.94	0.49
1:B:370:LYS:NZ	2:B:1301:UJJ:O1A	2.42	0.48
1:A:202:ARG:NH2	9:A:2321:HOH:O	2.45	0.48
1:B:782:GLY:O	1:B:811:THR:HA	2.13	0.48
1:B:1004:VAL:HG22	7:B:1312:PEG:H42	1.95	0.48
1:B:1026:LEU:HD23	1:B:1038:PRO:HG2	1.95	0.47
1:A:202:ARG:NH2	9:A:2324:HOH:O	2.46	0.47
1:A:803:ARG:NH2	1:A:1194:SER:H	2.13	0.47
1:B:338:VAL:HG22	1:B:367:ARG:HB3	1.96	0.47
1:B:1143:GLN:O	1:B:1147:LEU:HG	2.15	0.47
1:A:706:TRP:CE3	1:A:707:ASN:HA	2.50	0.46
1:B:197:THR:O	1:B:207:ARG:HD2	2.15	0.46
1:B:706:TRP:CE3	1:B:707:ASN:HA	2.51	0.46
1:B:465:LYS:HA	1:B:465:LYS:HD3	1.73	0.46
1:A:223[B]:LEU:HD13	1:A:482:LEU:HA	1.97	0.45
1:A:657:ILE:HD13	1:A:671:GLU:HG2	1.99	0.45
1:A:485:TYR:O	1:A:489:ARG:HG2	2.15	0.45
1:B:591[B]:THR:CG2	1:B:601:VAL:HG13	2.47	0.45
1:A:1026:LEU:HD23	1:A:1038:PRO:HG2	1.98	0.45
1:B:844:CYS:SG	4:B:1303:NAD:C3N	3.05	0.45
1:A:132:SER:HB2	1:A:133:HIS:CE1	2.51	0.44
1:B:297:LYS:HG3	1:B:332:TRP:HB2	1.99	0.44
1:A:485:TYR:CZ	1:A:489:ARG:HD2	2.52	0.44
1:B:338:VAL:HG22	1:B:367:ARG:HD2	1.99	0.44
1:A:1147:LEU:HD22	1:B:1147:LEU:HD22	2.00	0.44
1:A:782:GLY:O	1:A:811:THR:HA	2.18	0.43
1:B:448:CYS:HB2	1:B:453:GLY:HA3	2.00	0.43
1:B:617:VAL:HG12	1:B:774:GLU:HB2	2.00	0.43
1:A:931:LEU:HB2	1:A:949:ILE:HG22	2.01	0.43
1:A:1056:TRP:CD1	1:A:1142:PRO:HD3	2.54	0.43
1:B:495:ALA:HB2	1:B:1230:MET:HG3	2.01	0.43
1:B:659:ARG:HD3	9:B:2776:HOH:O	2.18	0.43
1:B:72:ILE:HG22	1:B:510:ILE:HD12	2.01	0.42
1:B:834:THR:O	1:B:838:ASP:HB3	2.18	0.42
1:B:705:PRO:HD3	1:B:781:THR:HB	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:428:LEU:HD11	1:B:460:VAL:HG21	2.00	0.42
1:B:64:ALA:HA	1:B:433:HIS:CD2	2.55	0.42
1:B:192:GLY:O	1:B:207:ARG:NH1	2.47	0.42
1:B:628:ALA:HB2	1:B:696:PRO:HG3	2.02	0.42
1:B:489:ARG:HA	1:B:489:ARG:NE	2.35	0.42
1:A:375:ASP:OD1	1:A:375:ASP:N	2.48	0.42
6:B:1311:PGE:H52	9:B:3190:HOH:O	2.19	0.41
1:A:367[B]:ARG:HA	1:A:419:GLN:HB2	2.01	0.41
1:B:717:ILE:HG12	1:B:727:VAL:HG11	2.03	0.41
1:B:651:PRO:HB3	6:B:1311:PGE:H2	2.03	0.41
1:A:1143:GLN:O	1:A:1147:LEU:HG	2.20	0.41
1:B:493:ASN:O	1:B:501:HIS:HB2	2.21	0.41
1:B:918:THR:HB	1:B:923:PHE:CD1	2.56	0.41
1:B:1069:ALA:HA	1:B:1117:ALA:HB1	2.03	0.41
1:B:712:ILE:HD13	1:B:781:THR:HG21	2.02	0.41
1:A:363:ARG:HA	1:A:415:VAL:O	2.21	0.40
1:B:578:PRO:O	1:B:584:PRO:HA	2.21	0.40
1:B:367:ARG:HA	1:B:419:GLN:HB2	2.03	0.40
1:B:374:TRP:CZ3	1:B:1229:LEU:HG	2.56	0.40
1:A:617:VAL:HG12	1:A:774:GLU:HB2	2.02	0.40
1:A:272:ARG:HB3	1:A:277:GLN:HG3	2.03	0.40
1:A:493:ASN:O	1:A:501:HIS:HB2	2.22	0.40
1:A:878:ARG:HD2	9:A:2643:HOH:O	2.21	0.40

There are no symmetry-related clashes.

## 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	1230/1235 (100%)	1204 (98%)	26 (2%)	0	100	100
1	B	1230/1235 (100%)	1204 (98%)	26 (2%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
All	All	2460/2470 (100%)	2408 (98%)	52 (2%)	0	100	100

There are no Ramachandran outliers to report.

### 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	909/951 (96%)	901 (99%)	8 (1%)	78	61
1	B	909/951 (96%)	902 (99%)	7 (1%)	81	66
All	All	1818/1902 (96%)	1803 (99%)	15 (1%)	81	66

All (15) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	39	TYR
1	A	181	ARG
1	A	321	GLU
1	A	342	TYR
1	A	631	PRO
1	A	730	LYS
1	A	796	ASP
1	A	934	LEU
1	B	39	TYR
1	B	342	TYR
1	B	489	ARG
1	B	730	LYS
1	B	872	HIS
1	B	934	LEU
1	B	1230	MET

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

Mol	Chain	Res	Type
1	B	419	GLN

### 5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

### 5.5 Carbohydrates [i](#)

There are no monosaccharides in this entry.

### 5.6 Ligand geometry [i](#)

Of 21 ligands modelled in this entry, 2 are monoatomic - leaving 19 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	$\# Z  > 2$
5	SO4	A	2207	-	4,4,4	0.14	0	6,6,6	0.08	0
6	PGE	A	2209	-	9,9,9	0.31	0	8,8,8	0.38	0
6	PGE	B	1311	-	9,9,9	0.31	0	8,8,8	0.32	0
5	SO4	A	2205	-	4,4,4	0.15	0	6,6,6	0.13	0
4	NAD	B	1303	3	42,48,48	3.44	16 (38%)	50,73,73	1.58	5 (10%)
4	NAD	A	2203	3	42,48,48	3.55	16 (38%)	50,73,73	1.71	6 (12%)
7	PEG	B	1312	-	6,6,6	0.48	0	5,5,5	0.35	0
5	SO4	A	2204	-	4,4,4	0.14	0	6,6,6	0.14	0
6	PGE	B	1308	-	9,9,9	0.30	0	8,8,8	0.35	0
2	UJJ	A	2201	-	56,64,64	2.27	20 (35%)	63,98,98	1.63	9 (14%)
2	UJJ	B	1301	-	56,64,64	2.29	19 (33%)	63,98,98	1.60	10 (15%)
5	SO4	A	2208	-	4,4,4	0.16	0	6,6,6	0.14	0
7	PEG	B	1309	-	6,6,6	0.30	0	5,5,5	0.25	0



Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
6	PGE	B	1306	-	9,9,9	0.30	0	8,8,8	0.26	0
5	SO4	B	1304	-	4,4,4	0.22	0	6,6,6	0.16	0
8	PG4	B	1310	-	12,12,12	0.50	0	11,11,11	0.19	0
5	SO4	B	1305	-	4,4,4	0.13	0	6,6,6	0.07	0
5	SO4	A	2206	-	4,4,4	0.14	0	6,6,6	0.12	0
7	PEG	B	1307	-	6,6,6	0.49	0	5,5,5	0.28	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
4	NAD	A	2203	3	-	4/26/62/62	0/5/5/5
8	PG4	B	1310	-	-	2/10/10/10	-
7	PEG	B	1312	-	-	2/4/4/4	-
7	PEG	B	1309	-	-	1/4/4/4	-
6	PGE	A	2209	-	-	2/7/7/7	-
6	PGE	B	1308	-	-	1/7/7/7	-
6	PGE	B	1311	-	-	6/7/7/7	-
6	PGE	B	1306	-	-	1/7/7/7	-
2	UJJ	A	2201	-	-	5/30/61/61	0/6/7/7
2	UJJ	B	1301	-	-	3/30/61/61	0/6/7/7
4	NAD	B	1303	3	-	2/26/62/62	0/5/5/5
7	PEG	B	1307	-	-	0/4/4/4	-

All (71) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	2203	NAD	O4D-C1D	-9.57	1.27	1.41
4	B	1303	NAD	O4D-C1D	-9.38	1.28	1.41
4	B	1303	NAD	C3B-C4B	-8.10	1.32	1.53
4	A	2203	NAD	C3B-C4B	-8.10	1.32	1.53
4	A	2203	NAD	C7N-N7N	7.92	1.48	1.33
4	A	2203	NAD	C3D-C4D	-7.52	1.33	1.53
4	B	1303	NAD	C7N-N7N	7.49	1.47	1.33
4	A	2203	NAD	O4D-C4D	7.08	1.60	1.45
4	B	1303	NAD	C3D-C4D	-6.94	1.35	1.53
4	A	2203	NAD	O4B-C4B	6.93	1.60	1.45
4	B	1303	NAD	O4B-C4B	6.81	1.60	1.45

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	1301	UJJ	O4-C4	6.80	1.36	1.23
4	B	1303	NAD	O4D-C4D	6.67	1.59	1.45
2	A	2201	UJJ	O4-C4	6.52	1.36	1.23
2	B	1301	UJJ	C6-C5X	5.84	1.49	1.39
2	B	1301	UJJ	O2-C2	5.69	1.35	1.23
2	A	2201	UJJ	C6-C5X	5.50	1.48	1.39
4	A	2203	NAD	O4B-C1B	-5.43	1.33	1.41
2	A	2201	UJJ	C9-C9A	5.36	1.48	1.39
4	B	1303	NAD	O4B-C1B	-5.35	1.33	1.41
2	A	2201	UJJ	O2-C2	5.22	1.34	1.23
2	B	1301	UJJ	C9-C9A	4.86	1.47	1.39
2	A	2201	UJJ	C4X-N5	4.69	1.45	1.37
4	A	2203	NAD	O3D-C3D	4.47	1.53	1.43
4	B	1303	NAD	C3N-C7N	4.29	1.57	1.50
4	B	1303	NAD	C6A-N6A	4.18	1.49	1.34
2	B	1301	UJJ	C4X-N5	4.13	1.44	1.37
4	A	2203	NAD	C3N-C7N	4.09	1.56	1.50
4	B	1303	NAD	O3D-C3D	4.04	1.52	1.43
4	A	2203	NAD	C6A-N6A	4.04	1.48	1.34
2	B	1301	UJJ	C10-N1	3.92	1.44	1.37
2	B	1301	UJJ	C2B-C1B	-3.73	1.48	1.53
2	B	1301	UJJ	C2-N1	3.64	1.43	1.37
2	A	2201	UJJ	C2-N1	3.50	1.43	1.37
2	A	2201	UJJ	C10-N1	3.50	1.43	1.37
4	B	1303	NAD	O3B-C3B	3.19	1.50	1.43
4	A	2203	NAD	O3B-C3B	3.05	1.50	1.43
4	B	1303	NAD	C2N-N1N	2.94	1.38	1.35
2	A	2201	UJJ	C2B-C1B	-2.93	1.49	1.53
2	B	1301	UJJ	C2A-N3A	2.92	1.36	1.32
2	A	2201	UJJ	C5X-N5	2.87	1.46	1.42
2	A	2201	UJJ	C6A-N6A	2.75	1.44	1.34
2	B	1301	UJJ	C6A-N6A	2.66	1.43	1.34
4	A	2203	NAD	C2N-N1N	2.65	1.38	1.35
2	A	2201	UJJ	C10-N10	2.56	1.43	1.38
2	B	1301	UJJ	C9A-N10	2.52	1.45	1.41
4	A	2203	NAD	C2A-N1A	2.51	1.38	1.33
2	B	1301	UJJ	C10-N10	2.50	1.42	1.38
4	B	1303	NAD	C2A-N1A	2.50	1.38	1.33
4	A	2203	NAD	C5A-C4A	-2.49	1.34	1.40
2	A	2201	UJJ	C2A-N3A	2.48	1.36	1.32
2	A	2201	UJJ	PA-O5B	-2.47	1.49	1.59
2	A	2201	UJJ	C9A-C5X	-2.47	1.36	1.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	1301	UJJ	C9A-C5X	-2.45	1.36	1.40
2	A	2201	UJJ	O2'-C2'	-2.37	1.38	1.43
4	A	2203	NAD	O7N-C7N	-2.35	1.19	1.24
4	A	2203	NAD	C2A-N3A	2.34	1.35	1.32
2	A	2201	UJJ	PA-O2A	-2.31	1.44	1.55
2	A	2201	UJJ	O3B-C3B	-2.30	1.37	1.43
2	B	1301	UJJ	PA-O2A	-2.28	1.44	1.55
2	A	2201	UJJ	C2-N3	2.27	1.41	1.37
2	B	1301	UJJ	O3B-C3B	-2.23	1.37	1.43
2	B	1301	UJJ	C2-N3	2.20	1.41	1.37
2	A	2201	UJJ	C4X-C4	2.19	1.49	1.43
2	B	1301	UJJ	C4X-C4	2.18	1.49	1.43
2	B	1301	UJJ	PA-O5B	-2.18	1.50	1.59
4	B	1303	NAD	C5A-C4A	-2.16	1.35	1.40
4	B	1303	NAD	O7N-C7N	-2.15	1.20	1.24
2	A	2201	UJJ	C2B-C3B	-2.14	1.47	1.53
2	B	1301	UJJ	P-O1P	2.07	1.58	1.50
4	B	1303	NAD	C2A-N3A	2.02	1.35	1.32

All (30) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	2203	NAD	C1B-N9A-C4A	-7.46	113.53	126.64
4	B	1303	NAD	C1B-N9A-C4A	-6.31	115.56	126.64
2	B	1301	UJJ	N3A-C2A-N1A	-5.98	119.32	128.68
2	A	2201	UJJ	N3A-C2A-N1A	-5.90	119.46	128.68
4	A	2203	NAD	N3A-C2A-N1A	-5.42	120.21	128.68
4	B	1303	NAD	N3A-C2A-N1A	-5.25	120.48	128.68
2	A	2201	UJJ	C4X-C4-N3	4.96	119.74	110.99
2	A	2201	UJJ	C4-N3-C2	-4.75	119.50	126.34
2	B	1301	UJJ	C4-N3-C2	-4.20	120.28	126.34
2	B	1301	UJJ	N3-C2-N1	4.20	122.55	115.80
2	B	1301	UJJ	C4X-C4-N3	4.14	118.29	110.99
2	A	2201	UJJ	N3-C2-N1	3.66	121.68	115.80
4	A	2203	NAD	C3N-C2N-N1N	-3.31	117.19	120.43
4	A	2203	NAD	O4D-C1D-C2D	-3.09	102.41	106.93
4	B	1303	NAD	C3N-C2N-N1N	-2.99	117.51	120.43
4	B	1303	NAD	O4B-C1B-C2B	-2.77	102.88	106.93
2	B	1301	UJJ	C1B-N9A-C4A	-2.76	121.80	126.64
2	B	1301	UJJ	C2A-N1A-C6A	2.66	123.30	118.75
2	A	2201	UJJ	C2A-N1A-C6A	2.59	123.18	118.75
4	A	2203	NAD	C2N-C3N-C4N	2.58	121.19	118.26

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	2203	NAD	O4B-C1B-C2B	-2.53	103.23	106.93
2	A	2201	UJJ	O4-C4-C4X	-2.53	121.33	127.54
2	A	2201	UJJ	C4A-C5A-N7A	-2.43	106.86	109.40
2	B	1301	UJJ	C6-C5X-N5	-2.36	117.50	121.58
2	B	1301	UJJ	O4-C4-C4X	-2.31	121.87	127.54
2	B	1301	UJJ	O2-C2-N1	-2.29	117.50	121.82
2	B	1301	UJJ	C1'-N10-C9A	2.27	124.30	120.51
4	B	1303	NAD	C3N-C7N-N7N	2.13	120.30	117.75
2	A	2201	UJJ	C1B-N9A-C4A	-2.06	123.02	126.64
2	A	2201	UJJ	C1'-N10-C9A	2.06	123.94	120.51

There are no chirality outliers.

All (29) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	2203	NAD	C5B-O5B-PA-O1A
6	A	2209	PGE	O1-C1-C2-O2
7	B	1312	PEG	O2-C3-C4-O4
6	B	1311	PGE	O1-C1-C2-O2
6	B	1306	PGE	O1-C1-C2-O2
8	B	1310	PG4	C8-C7-O4-C6
4	B	1303	NAD	C4D-C5D-O5D-PN
6	B	1311	PGE	C1-C2-O2-C3
4	A	2203	NAD	C5B-O5B-PA-O3
6	B	1311	PGE	C4-C3-O2-C2
4	A	2203	NAD	C4D-C5D-O5D-PN
2	A	2201	UJJ	C2'-C3'-C4'-O4'
7	B	1309	PEG	C1-C2-O2-C3
8	B	1310	PG4	C6-C5-O3-C4
7	B	1312	PEG	C4-C3-O2-C2
6	B	1311	PGE	C3-C4-O3-C5
6	B	1311	PGE	O3-C5-C6-O4
2	A	2201	UJJ	PA-O3P-P-O1P
2	B	1301	UJJ	PA-O3P-P-O1P
2	A	2201	UJJ	C4'-C5'-O5'-P
4	A	2203	NAD	C3D-C4D-C5D-O5D
6	A	2209	PGE	C4-C3-O2-C2
2	A	2201	UJJ	C5'-O5'-P-O3P
2	A	2201	UJJ	PA-O3P-P-O2P
2	B	1301	UJJ	PA-O3P-P-O2P
2	B	1301	UJJ	C4'-C5'-O5'-P
4	B	1303	NAD	C5B-O5B-PA-O1A

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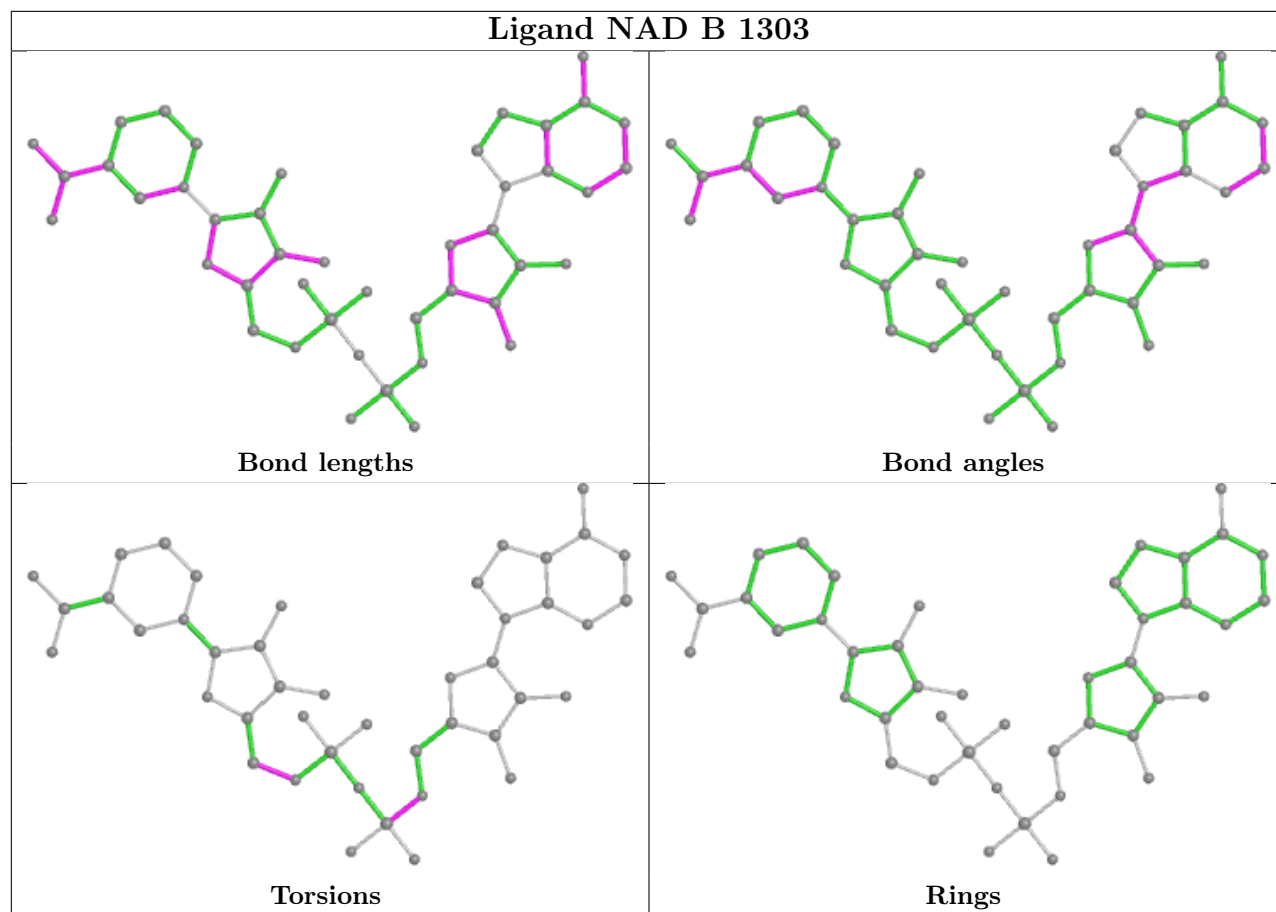
Mol	Chain	Res	Type	Atoms
6	B	1308	PGE	O3-C5-C6-O4
6	B	1311	PGE	O2-C3-C4-O3

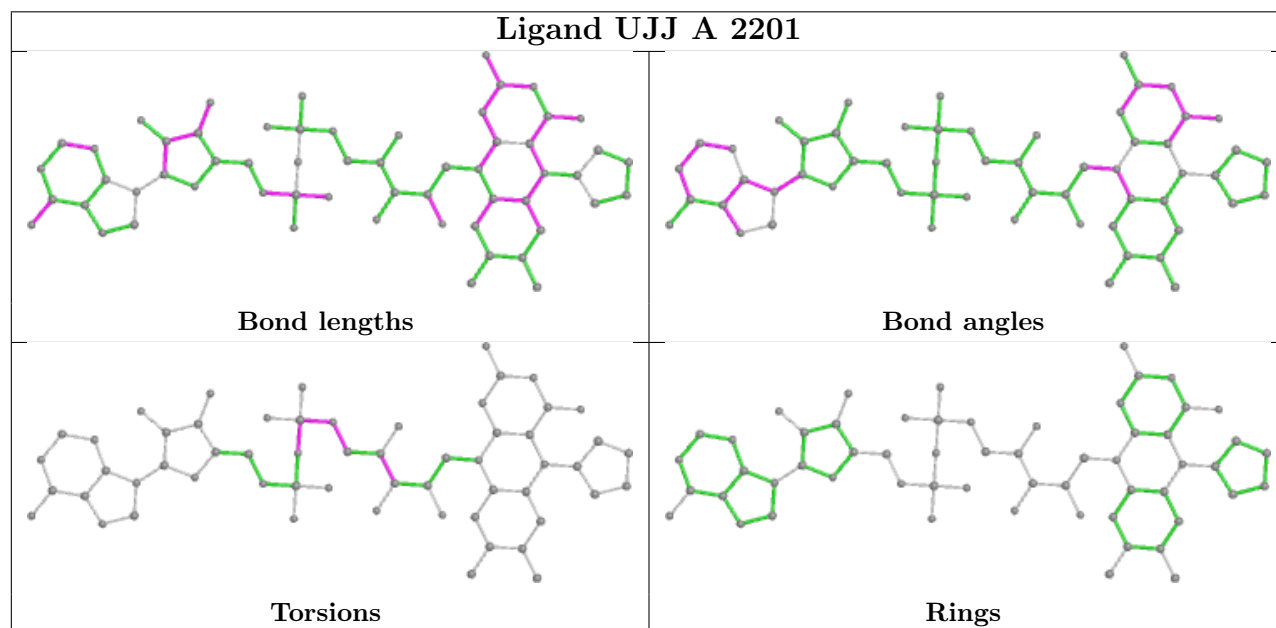
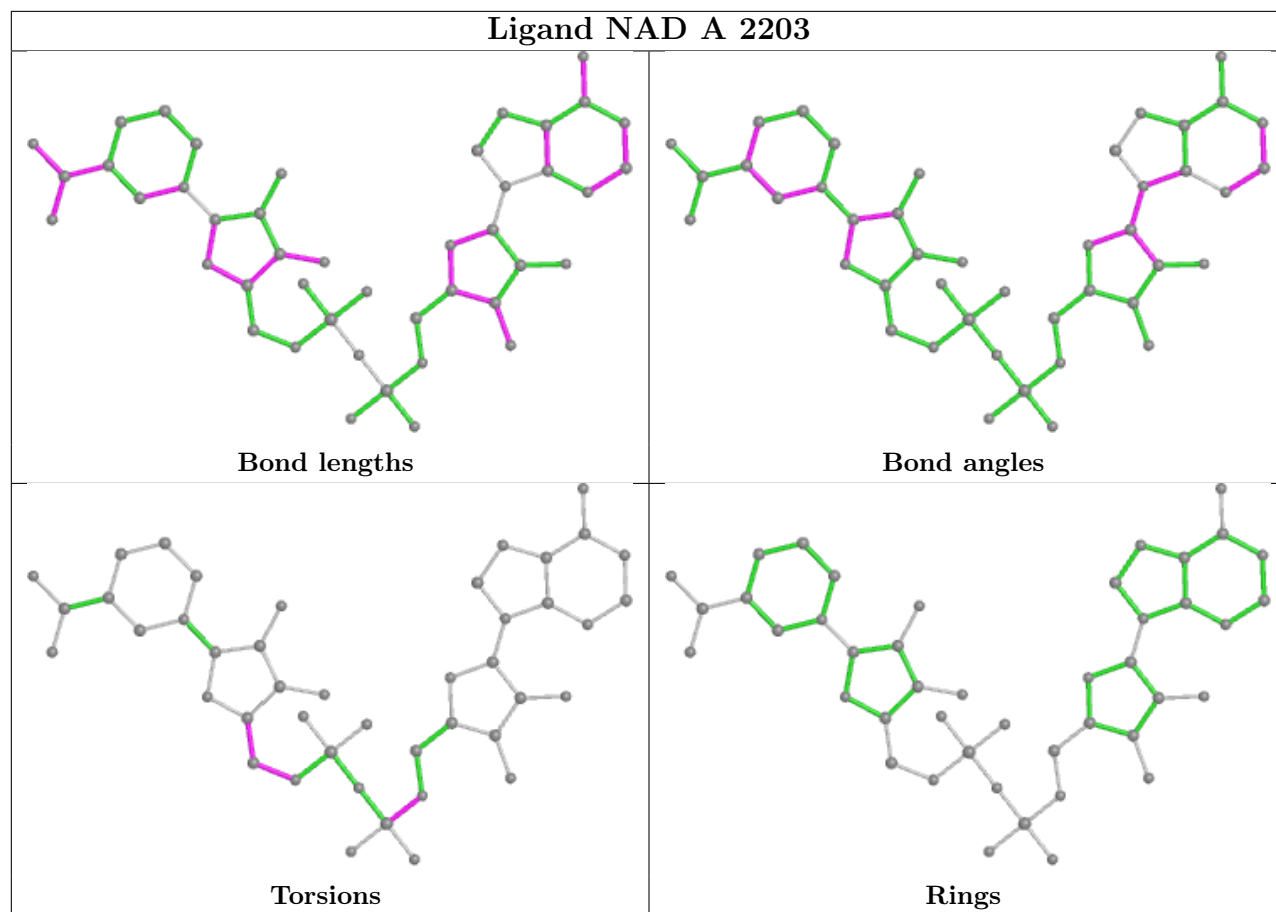
There are no ring outliers.

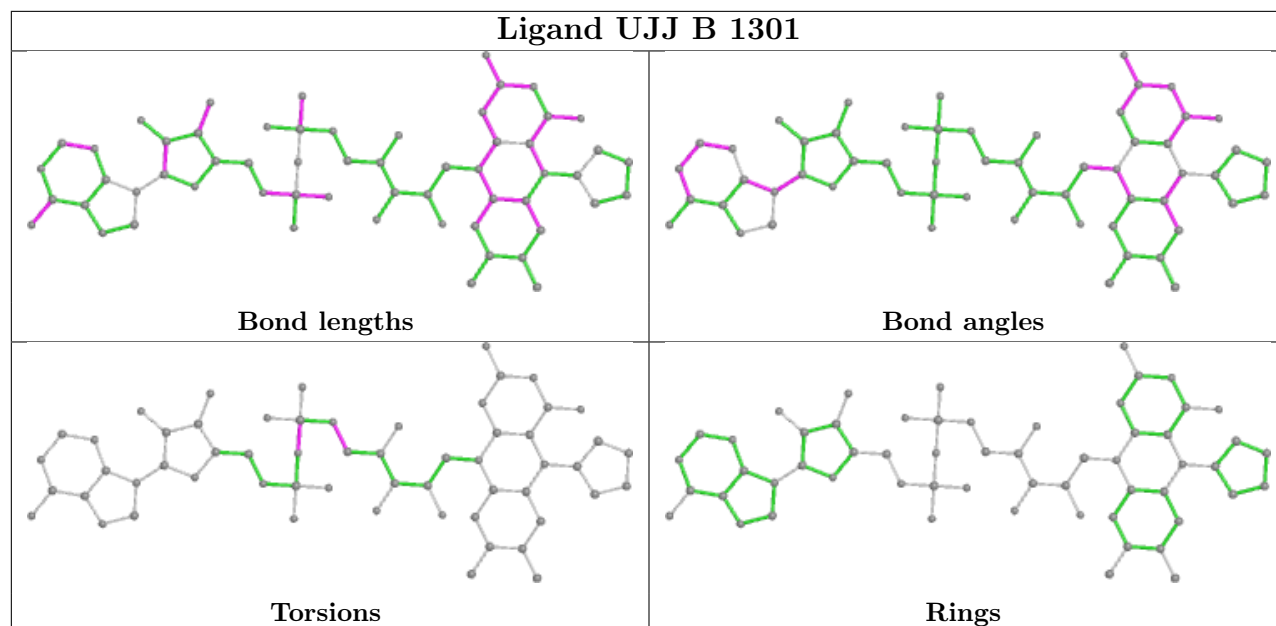
6 monomers are involved in 9 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	B	1311	PGE	3	0
4	B	1303	NAD	2	0
4	A	2203	NAD	1	0
7	B	1312	PEG	1	0
6	B	1308	PGE	1	0
2	B	1301	UJJ	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	1217/1235 (98%)	0.23	70 (5%) 23 26	13, 21, 48, 86	0
1	B	1217/1235 (98%)	0.26	78 (6%) 19 23	12, 22, 48, 83	0
All	All	2434/2470 (98%)	0.24	148 (6%) 21 25	12, 21, 48, 86	0

All (148) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	81	GLY	9.8
1	A	80	SER	7.2
1	B	81	GLY	6.8
1	B	14	ALA	6.7
1	A	133	HIS	6.7
1	A	132	SER	6.4
1	A	82	SER	6.1
1	A	83	GLY	6.0
1	A	1231	ALA	5.9
1	A	84	VAL	5.7
1	A	482	LEU	5.6
1	A	490	LEU	5.6
1	B	1223	ALA	5.6
1	A	79	HIS	5.6
1	B	195	PHE	5.5
1	B	1231	ALA	5.3
1	A	192	GLY	5.3
1	B	1222	ALA	5.2
1	A	134	LEU	5.2
1	A	111	ILE	5.2
1	A	120	LEU	5.2
1	A	14	ALA	5.1
1	B	496	ASN	5.0
1	B	1232	ILE	5.0

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Mol	Chain	Res	Type	RSRZ
1	B	918	THR	5.0
1	A	486	LEU	5.0
1	B	506	PRO	5.0
1	A	191	MET	5.0
1	A	506	PRO	4.9
1	A	195	PHE	4.8
1	A	130	TRP	4.8
1	B	1233	GLY	4.8
1	A	116	THR	4.7
1	B	490	LEU	4.7
1	A	115	ALA	4.6
1	A	496	ASN	4.5
1	A	223[A]	LEU	4.4
1	B	912	ILE	4.3
1	A	119	ALA	4.3
1	B	15	PRO	4.2
1	B	60	ILE	4.2
1	B	80	SER	4.2
1	A	508	VAL	4.2
1	A	75	LEU	4.1
1	B	438	ASP	4.0
1	A	125	ILE	4.0
1	B	84	VAL	4.0
1	B	486	LEU	4.0
1	B	192	GLY	4.0
1	A	123	ASP	4.0
1	B	482	LEU	3.9
1	B	916	SER	3.9
1	B	508	VAL	3.9
1	A	112	PRO	3.8
1	A	129	ASN	3.8
1	B	914	LEU	3.8
1	B	82	SER	3.8
1	B	505	ASP	3.7
1	A	77	GLY	3.6
1	B	915	ALA	3.6
1	A	483	LEU	3.6
1	B	1140	ASN	3.6
1	B	1230	MET	3.6
1	B	917	GLU	3.5
1	B	439	PHE	3.5
1	B	79	HIS	3.4

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Mol	Chain	Res	Type	RSRZ
1	A	190	MET	3.4
1	A	503	ILE	3.4
1	A	128	GLY	3.4
1	A	1233	GLY	3.4
1	B	133	HIS	3.3
1	B	1059	GLY	3.3
1	B	129	ASN	3.3
1	B	491	LEU	3.2
1	B	913	GLY	3.2
1	B	503	ILE	3.1
1	A	194	GLN	3.1
1	A	185	ASP	3.1
1	B	132	SER	3.1
1	B	495	ALA	3.1
1	B	514	ILE	3.1
1	B	1227	ALA	3.1
1	A	1002	GLY	3.1
1	A	114	THR	3.0
1	B	115	ALA	3.0
1	B	481	THR	3.0
1	B	483	LEU	3.0
1	B	494	GLY	3.0
1	A	181	ARG	2.9
1	A	126	ALA	2.9
1	A	497	SER	2.9
1	A	147	TRP	2.9
1	B	493	ASN	2.9
1	A	499	PHE	2.9
1	A	121	ILE	2.9
1	A	1232	ILE	2.9
1	B	224	GLY	2.8
1	A	188	MET	2.8
1	B	919	GLY	2.8
1	B	19	ALA	2.8
1	A	214	LYS	2.8
1	A	504	ASN	2.8
1	A	186	MET	2.7
1	B	800	PRO	2.7
1	B	510	ILE	2.7
1	B	119	ALA	2.7
1	A	117	ARG	2.6
1	B	436	GLY	2.6

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Mol	Chain	Res	Type	RSRZ
1	B	138	ARG	2.6
1	B	501	HIS	2.6
1	B	134	LEU	2.5
1	A	78	LYS	2.5
1	B	515	ALA	2.4
1	A	131	LYS	2.4
1	A	1223	ALA	2.4
1	B	63	ALA	2.4
1	B	1229	LEU	2.4
1	B	521	VAL	2.4
1	A	91	TYR	2.4
1	B	499	PHE	2.4
1	A	109	LEU	2.4
1	A	493	ASN	2.3
1	A	1140	ASN	2.3
1	B	194	GLN	2.3
1	A	1230	MET	2.3
1	A	122	ARG	2.3
1	B	520	VAL	2.3
1	B	191	MET	2.3
1	A	524	MET	2.3
1	A	118	ASP	2.3
1	B	485	TYR	2.2
1	B	437	LYS	2.2
1	B	114	THR	2.2
1	A	494	GLY	2.2
1	B	130	TRP	2.2
1	A	15	PRO	2.2
1	B	1136	ALA	2.1
1	A	505	ASP	2.1
1	B	75	LEU	2.1
1	B	462	GLY	2.1
1	B	1228	SER	2.1
1	A	916	SER	2.1
1	B	511	ASP	2.1
1	B	513	LEU	2.1
1	B	59	GLU	2.0
1	B	83	GLY	2.0
1	A	1228	SER	2.0
1	B	124	LYS	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](#)

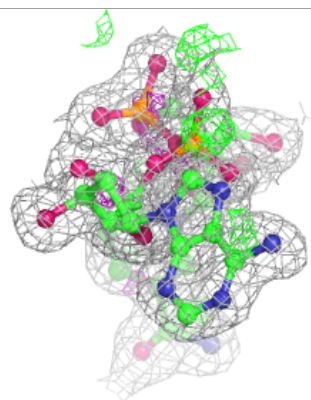
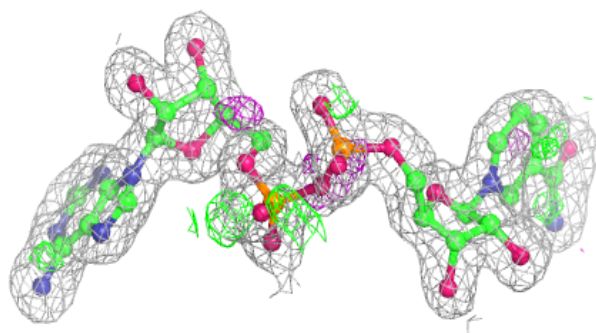
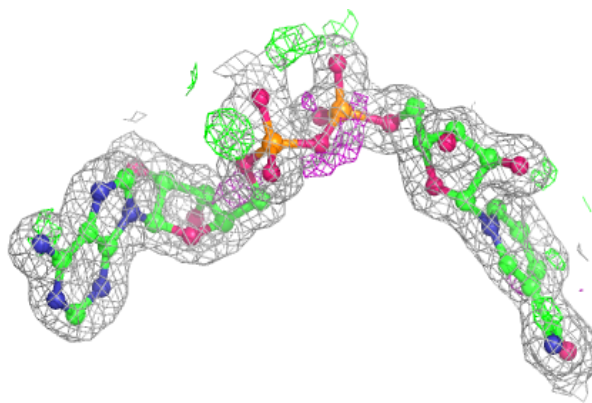
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
6	PGE	B	1311	10/10	0.75	0.26	47,56,63,67	0
5	SO4	B	1305	5/5	0.81	0.23	100,102,102,104	0
6	PGE	B	1308	10/10	0.88	0.14	33,47,62,66	0
7	PEG	B	1312	7/7	0.88	0.17	23,44,50,51	0
7	PEG	B	1307	7/7	0.89	0.10	40,42,51,53	0
5	SO4	A	2205	5/5	0.89	0.19	61,61,64,72	0
8	PG4	B	1310	13/13	0.89	0.12	40,51,57,58	0
5	SO4	A	2208	5/5	0.90	0.23	64,69,75,80	0
7	PEG	B	1309	7/7	0.92	0.15	27,28,35,42	0
4	NAD	A	2203	44/44	0.93	0.09	18,22,28,30	0
6	PGE	A	2209	10/10	0.93	0.09	23,38,47,51	0
5	SO4	A	2206	5/5	0.94	0.11	54,56,61,64	0
2	UJJ	B	1301	58/58	0.94	0.08	16,21,26,31	0
6	PGE	B	1306	10/10	0.94	0.08	27,40,45,48	0
3	MG	A	2202	1/1	0.95	0.11	32,32,32,32	0
5	SO4	A	2207	5/5	0.95	0.23	71,72,74,78	0
3	MG	B	1302	1/1	0.95	0.11	27,27,27,27	0
4	NAD	B	1303	44/44	0.96	0.09	13,16,20,32	0
2	UJJ	A	2201	58/58	0.96	0.07	13,19,24,32	0
5	SO4	B	1304	5/5	0.99	0.05	15,19,23,24	0
5	SO4	A	2204	5/5	0.99	0.05	17,19,22,23	0

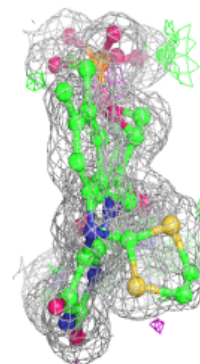
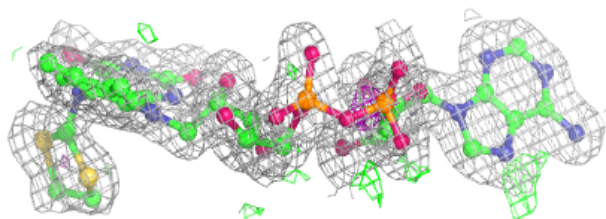
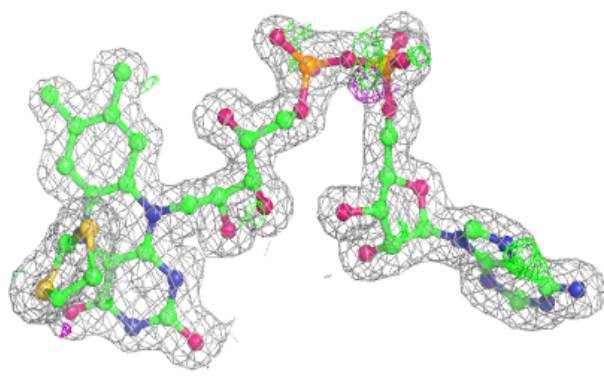
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

**Electron density around NAD A 2203:**

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

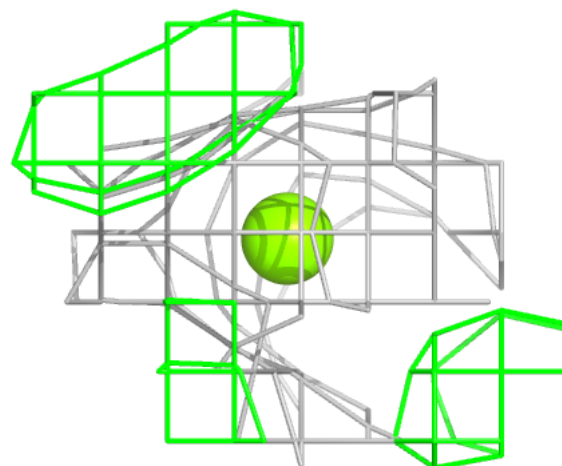
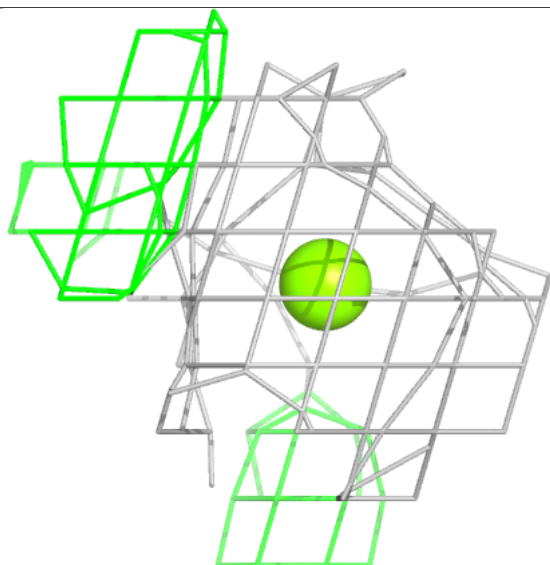
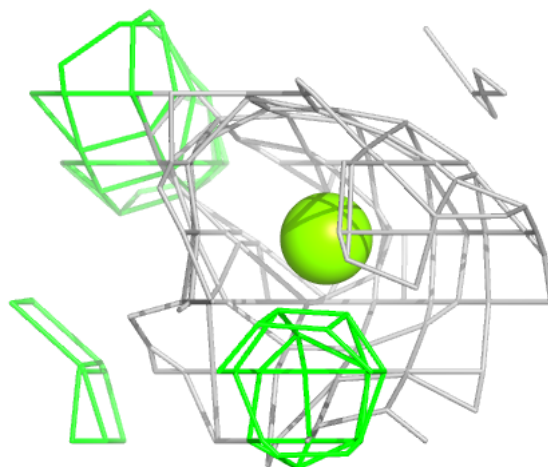
**Electron density around UJJ B 1301:**

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



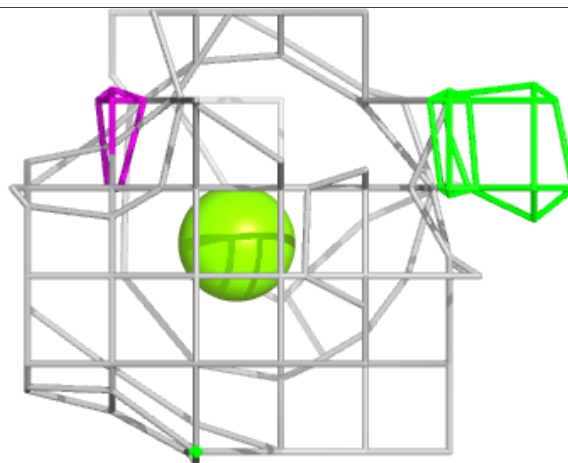
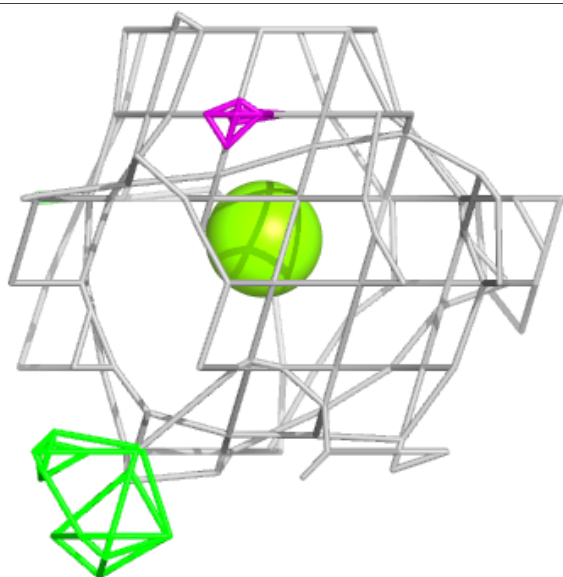
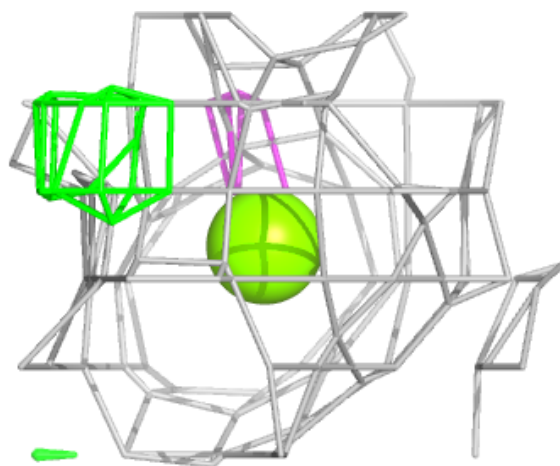
**Electron density around MG A 2202:**

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



**Electron density around MG B 1302:**

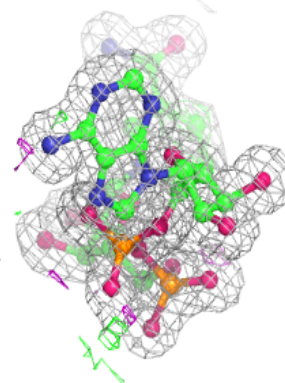
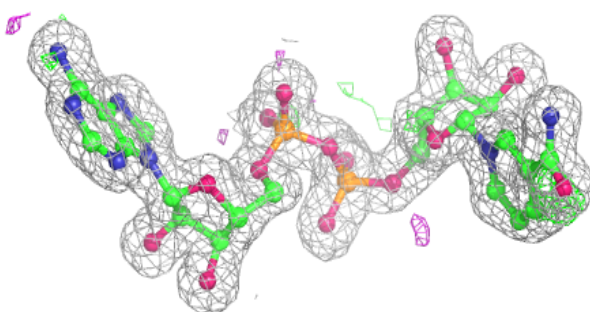
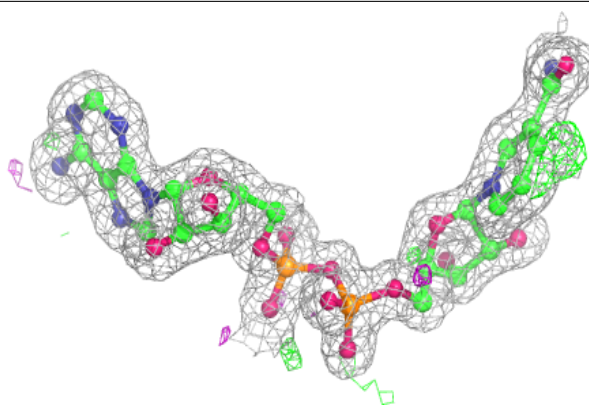
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)



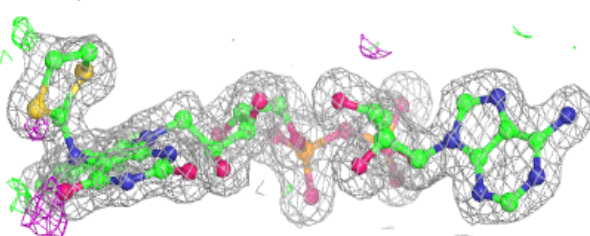
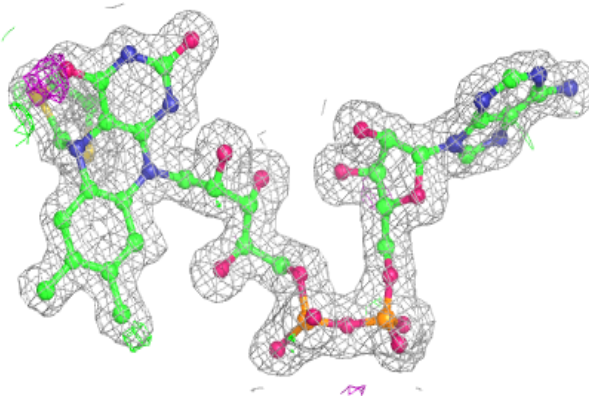


**Electron density around NAD B 1303:**

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

**Electron density around UJJ A 2201:**

$2mF_o-DF_c$  (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.