



# wwPDB X-ray Structure Validation Summary Report ⓘ

Dec 14, 2020 – 09:54 PM EST

PDB ID : 6VC6  
Title : 2.1 Angstrom Resolution Crystal Structure of 6-phospho-alpha-glucosidase from Gut Microorganisms in Complex with NAD and Mn<sup>2+</sup>  
Authors : Wu, R.; Kim, Y.; Endres, M.; Joachimiak, J.  
Deposited on : 2019-12-20  
Resolution : 2.13 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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.

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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 : 2.15.1  
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) : 2.15.1

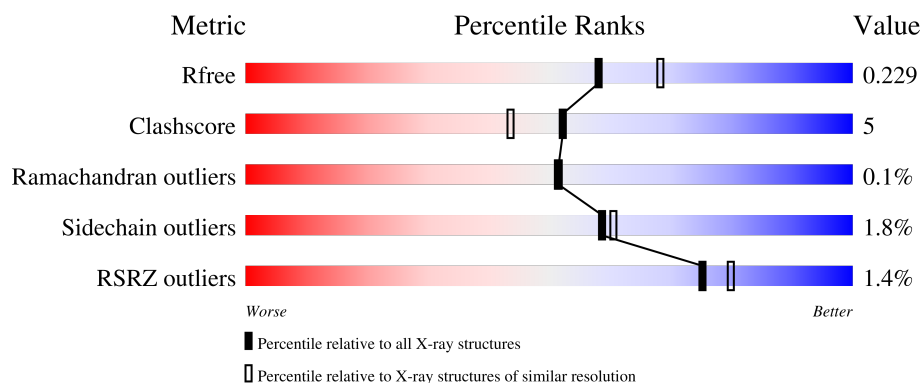
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.13 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	2523 (2.16-2.12)
Clashscore	141614	2653 (2.16-2.12)
Ramachandran outliers	138981	2618 (2.16-2.12)
Sidechain outliers	138945	2617 (2.16-2.12)
RSRZ outliers	127900	2485 (2.16-2.12)

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	441	<div> <div>88%</div> <div>9%</div> <div>.</div> </div>
1	B	441	<div> <div>87%</div> <div>11%</div> <div>.</div> </div>
1	C	441	<div> <div>86%</div> <div>12%</div> <div>..</div> </div>
1	D	441	<div> <div>82%</div> <div>15%</div> <div>.</div> </div>

## 2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 14988 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.

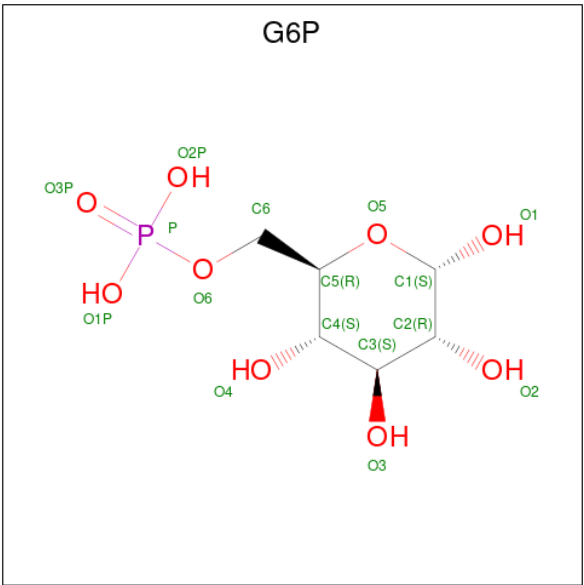
- Molecule 1 is a protein called 6-phospho-alpha-glucosidase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	432	Total	C	N	O	S	Se	0	0	0
			3504	2234	591	656	6	17			
1	B	440	Total	C	N	O	S	Se	0	0	0
			3567	2272	602	670	6	17			
1	C	438	Total	C	N	O	S	Se	0	0	0
			3552	2262	600	667	6	17			
1	D	433	Total	C	N	O	S	Se	1	0	0
			3508	2236	592	657	6	17			

- Molecule 2 is MANGANESE (II) ION (three-letter code: MN) (formula: Mn) (labeled as "Ligand of Interest" by depositor).

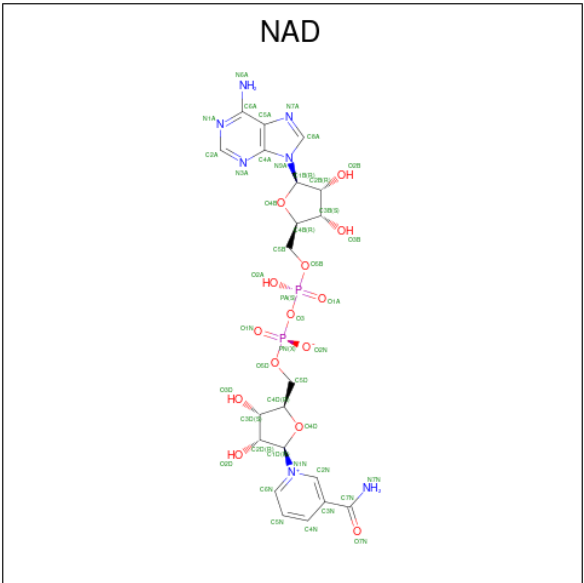
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	1	Total	Mn	0	0
			1	1		
2	A	1	Total	Mn	0	0
			1	1		
2	D	1	Total	Mn	0	0
			1	1		
2	C	1	Total	Mn	0	0
			1	1		

- Molecule 3 is 6-O-phosphono-alpha-D-glucopyranose (three-letter code: G6P) (formula: C<sub>6</sub>H<sub>13</sub>O<sub>9</sub>P) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	O	P	0	0
			16	6	9	1		
3	B	1	Total	C	O	P	0	0
			16	6	9	1		
3	C	1	Total	C	O	P	0	0
			16	6	9	1		
3	D	1	Total	C	O	P	0	0
			16	6	9	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		
4	C	1	Total	C	N	O	P	0	0
			44	21	7	14	2		
4	D	1	Total	C	N	O	P	0	0
			44	21	7	14	2		

- Molecule 5 is GLYCEROL (three-letter code: GOL) (formula:  $C_3H_8O_3$ ).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	B	1	Total	C	O	0	0
			6	3	3		
5	C	1	Total	C	O	0	0
			6	3	3		
5	C	1	Total	C	O	0	0
			6	3	3		

- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	176	Total	O	0	0
			176	176		
6	B	156	Total	O	0	0
			156	156		
6	C	146	Total	O	0	0
			146	146		

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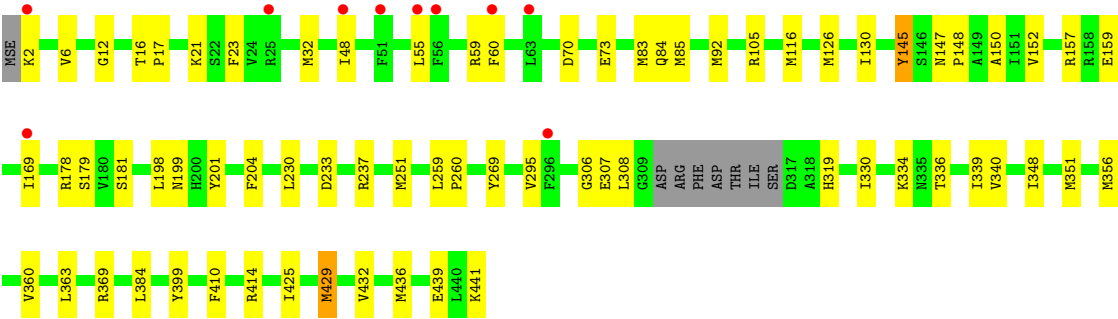
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	D	117	Total	O	0	0
			117	117		



- Molecule 1: 6-phospho-alpha-glucosidase







## 4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	194.46Å 62.26Å 160.71Å 90.00° 90.35° 90.00°	Depositor
Resolution (Å)	48.61 – 2.13 48.61 – 2.13	Depositor EDS
% Data completeness (in resolution range)	98.9 (48.61-2.13) 99.0 (48.61-2.13)	Depositor EDS
$R_{merge}$	0.12	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.38 (at 2.14Å)	Xtriage
Refinement program	PHENIX 1.11.1-2575_1309	Depositor
R, $R_{free}$	0.196 , 0.229 0.197 , 0.229	Depositor DCC
$R_{free}$ test set	5316 reflections (4.99%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	41.6	Xtriage
Anisotropy	0.300	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.33 , 37.1	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.011 for -h,-k,l	Xtriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	14988	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	49.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 10.31% of the height of the origin peak. No significant pseudotranslation is detected.*

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

<sup>2</sup>Theoretical values of  $\langle |L| \rangle$ ,  $\langle L^2 \rangle$  for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

## 5 Model quality

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: GOL, G6P, MN, NAD

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

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# $ Z  > 5$	RMSZ	# $ Z  > 5$
1	A	0.30	1/3567 (0.0%)	0.45	2/4791 (0.0%)
1	B	0.32	2/3632 (0.1%)	0.46	2/4880 (0.0%)
1	C	0.28	0/3616	0.44	0/4856
1	D	0.30	0/3571	0.43	0/4796
All	All	0.30	3/14386 (0.0%)	0.44	4/19323 (0.0%)

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	A	0	1

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	61	PRO	N-CD	5.25	1.55	1.47
1	A	138	PRO	N-CD	5.23	1.55	1.47
1	B	231	PRO	N-CD	5.16	1.55	1.47

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	60	PHE	C-N-CD	5.71	140.39	128.40
1	B	230	LEU	C-N-CD	5.54	140.03	128.40
1	A	137	SER	C-N-CD	5.49	139.93	128.40
1	B	60	PHE	C-N-CD	5.47	139.89	128.40

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	298	GLU	Mainchain

## 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	3504	0	3411	29	0
1	B	3567	0	3468	36	0
1	C	3552	0	3449	41	0
1	D	3508	0	3414	50	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
3	A	16	0	10	1	0
3	B	16	0	9	1	0
3	C	16	0	10	0	0
3	D	16	0	10	1	0
4	A	44	0	26	2	0
4	B	44	0	26	1	0
4	C	44	0	26	1	0
4	D	44	0	26	2	0
5	B	6	0	8	0	0
5	C	12	0	16	0	0
6	A	176	0	0	0	0
6	B	156	0	0	0	0
6	C	146	0	0	2	0
6	D	117	0	0	1	0
All	All	14988	0	13909	155	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 5.

The worst 5 of 155 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:229:PHE:CD2	1:A:248:VAL:HG21	2.05	0.90
1:B:25:ARG:HH12	1:B:59:ARG:NH1	1.77	0.83
1:B:25:ARG:HH12	1:B:59:ARG:HH11	1.29	0.80
1:D:306:GLY:O	1:D:307:GLU:HG3	1.81	0.80
1:A:415:LEU:C	1:A:416:ILE:HD12	2.03	0.78

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	428/441 (97%)	419 (98%)	9 (2%)	0	100	100
1	B	438/441 (99%)	429 (98%)	8 (2%)	1 (0%)	47	45
1	C	434/441 (98%)	427 (98%)	6 (1%)	1 (0%)	47	45
1	D	429/441 (97%)	422 (98%)	7 (2%)	0	100	100
All	All	1729/1764 (98%)	1697 (98%)	30 (2%)	2 (0%)	51	51

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	201	TYR
1	C	201	TYR

### 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	374/364 (103%)	368 (98%)	6 (2%)	62	65
1	B	381/364 (105%)	372 (98%)	9 (2%)	49	49
1	C	379/364 (104%)	373 (98%)	6 (2%)	62	65
1	D	374/364 (103%)	368 (98%)	6 (2%)	62	65
All	All	1508/1456 (104%)	1481 (98%)	27 (2%)	59	60

5 of 27 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	211	LYS
1	C	60	PHE
1	D	145	TYR
1	B	371	MSE
1	A	251	MSE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	C	28	ASN
1	D	84	GLN
1	D	102	HIS
1	D	319	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 15 ligands modelled in this entry, 4 are monoatomic - leaving 11 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	GOL	C	504	-	5,5,5	0.37	0	5,5,5	0.28	0
3	G6P	A	502	2	16,16,16	0.58	0	24,24,24	1.61	3 (12%)
3	G6P	C	502	2	16,16,16	0.56	0	24,24,24	0.89	0
5	GOL	C	505	-	5,5,5	0.20	0	5,5,5	0.10	0
4	NAD	A	503	2	42,48,48	0.82	1 (2%)	50,73,73	1.11	3 (6%)
4	NAD	B	503	2	42,48,48	0.81	1 (2%)	50,73,73	1.12	3 (6%)
3	G6P	B	502	2	16,16,16	0.57	0	24,24,24	0.92	0
4	NAD	D	503	2	42,48,48	0.81	1 (2%)	50,73,73	1.12	3 (6%)
3	G6P	D	502	2	16,16,16	0.54	0	24,24,24	0.78	0
5	GOL	B	504	-	5,5,5	0.37	0	5,5,5	0.27	0
4	NAD	C	503	2	42,48,48	0.81	1 (2%)	50,73,73	1.13	3 (6%)

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
5	GOL	C	504	-	-	2/4/4/4	-
3	G6P	A	502	2	-	2/6/26/26	0/1/1/1
3	G6P	C	502	2	-	5/6/26/26	0/1/1/1
5	GOL	C	505	-	-	3/4/4/4	-
4	NAD	A	503	2	-	4/26/62/62	0/5/5/5
4	NAD	B	503	2	-	7/26/62/62	0/5/5/5
3	G6P	B	502	2	-	5/6/26/26	0/1/1/1
4	NAD	D	503	2	-	4/26/62/62	0/5/5/5
3	G6P	D	502	2	-	5/6/26/26	0/1/1/1
5	GOL	B	504	-	-	4/4/4/4	-
4	NAD	C	503	2	-	4/26/62/62	0/5/5/5

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	503	NAD	C5A-C4A	2.45	1.47	1.40
4	C	503	NAD	C5A-C4A	2.45	1.47	1.40
4	A	503	NAD	C5A-C4A	2.44	1.47	1.40
4	D	503	NAD	C5A-C4A	2.39	1.47	1.40

The worst 5 of 15 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	502	G6P	C1-O5-C5	4.39	121.94	113.66
3	A	502	G6P	O5-C1-C2	4.37	118.09	110.28
4	D	503	NAD	N3A-C2A-N1A	-3.30	123.53	128.68
4	A	503	NAD	N3A-C2A-N1A	-3.29	123.54	128.68
4	C	503	NAD	N3A-C2A-N1A	-3.27	123.56	128.68

There are no chirality outliers.

5 of 45 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	C	504	GOL	O1-C1-C2-C3
3	C	502	G6P	C4-C5-C6-O6
3	C	502	G6P	O5-C5-C6-O6
3	C	502	G6P	C6-O6-P-O1P
3	C	502	G6P	C6-O6-P-O2P

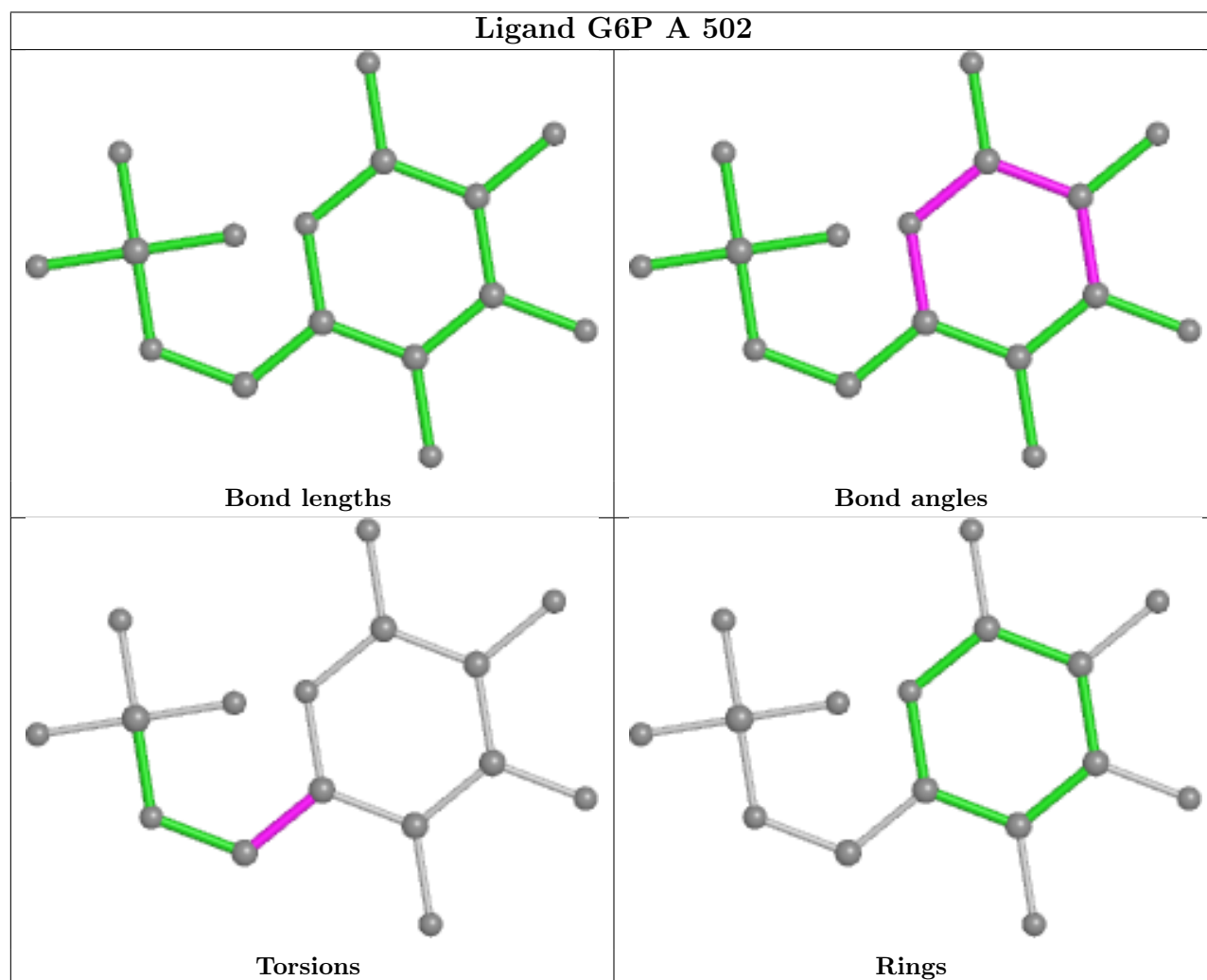
There are no ring outliers.

7 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	502	G6P	1	0
4	A	503	NAD	2	0
4	B	503	NAD	1	0
3	B	502	G6P	1	0
4	D	503	NAD	2	0
3	D	502	G6P	1	0
4	C	503	NAD	1	0

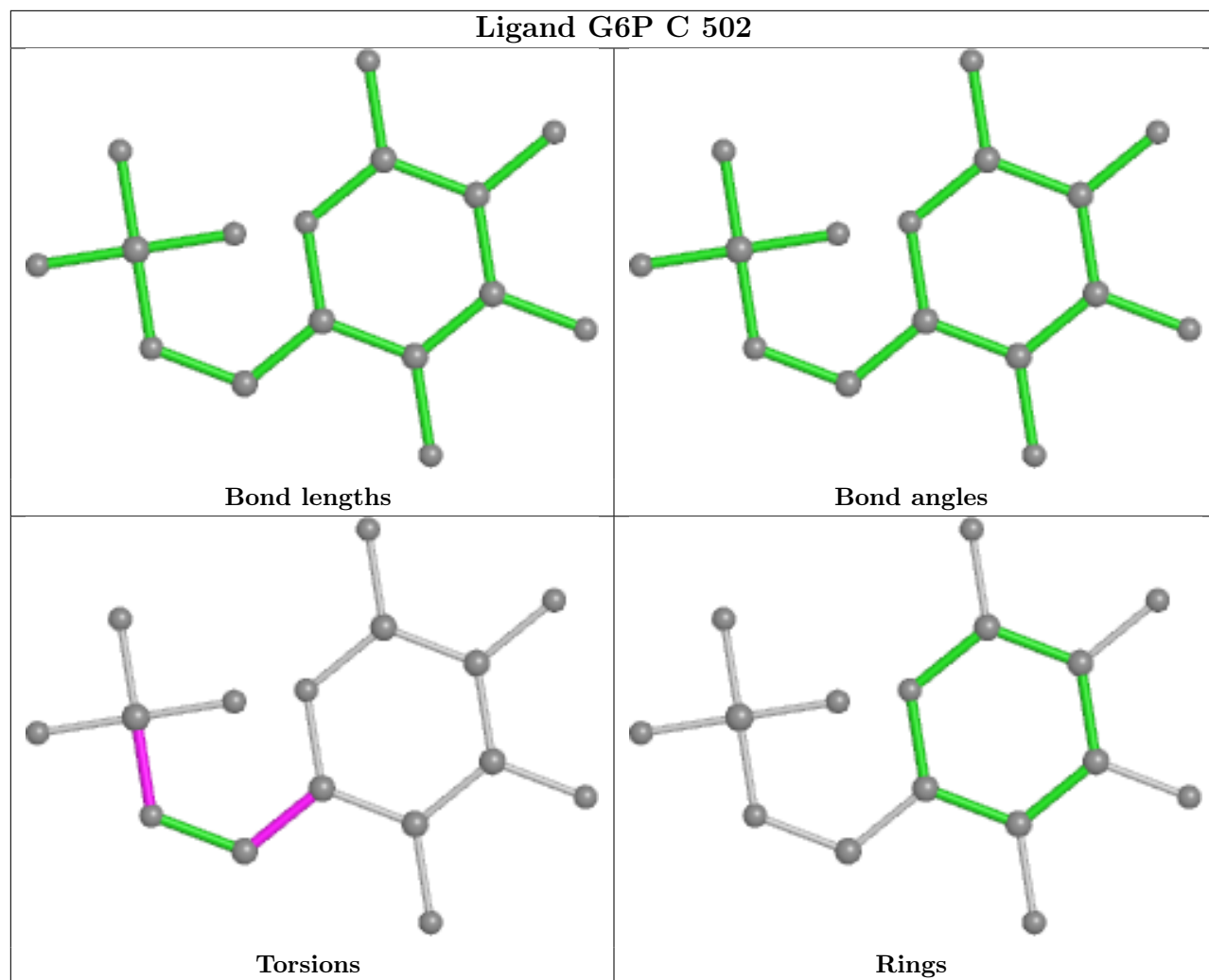
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be

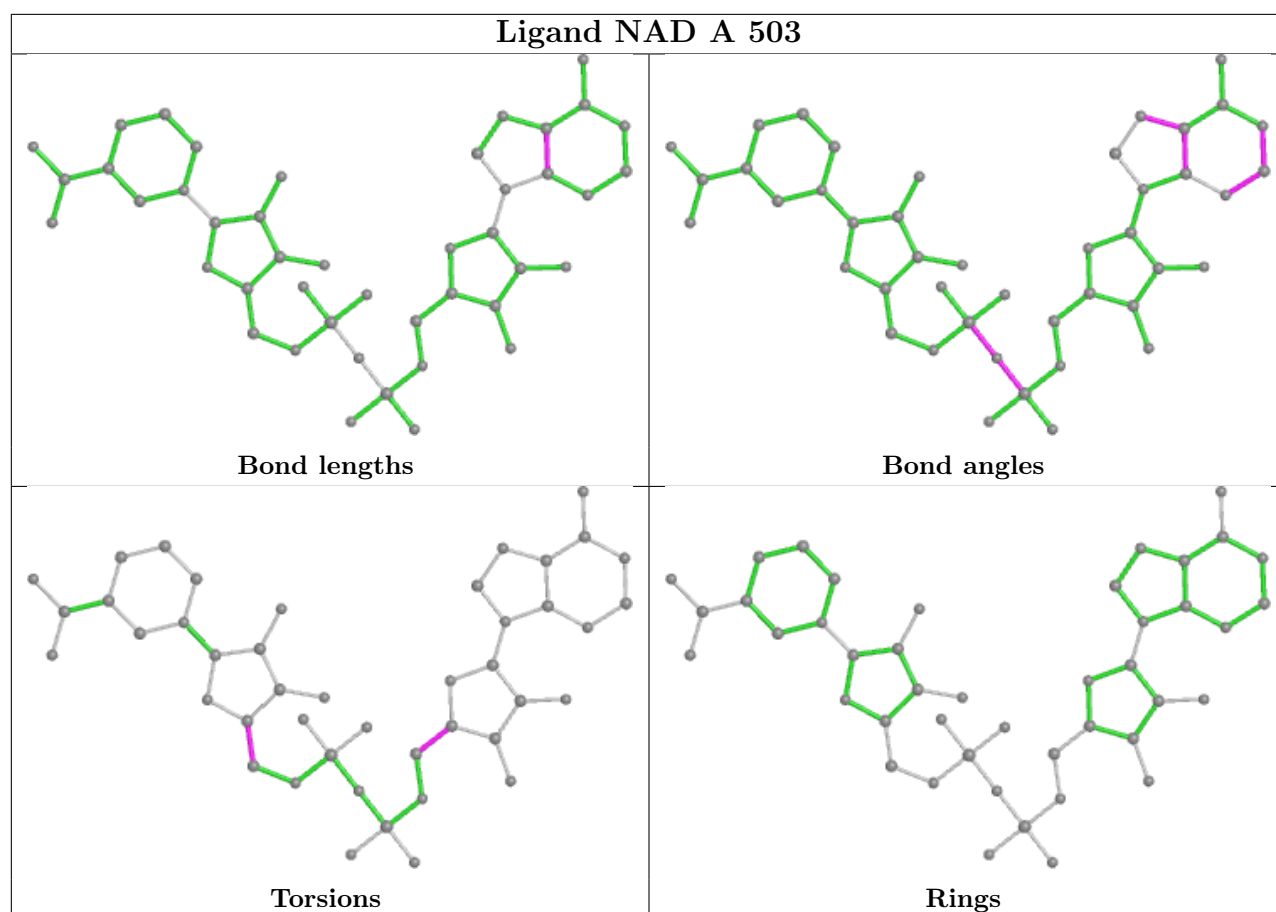
highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.

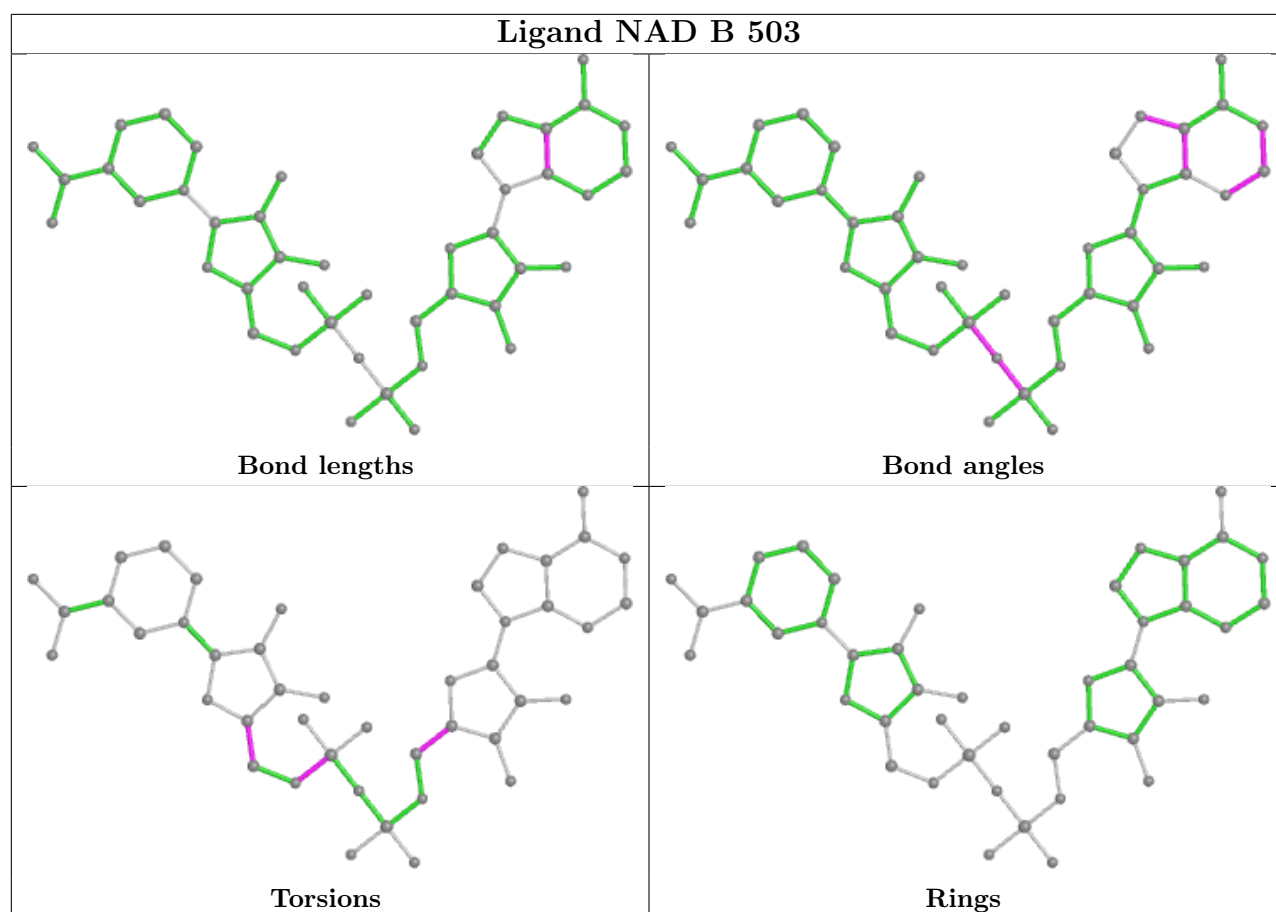


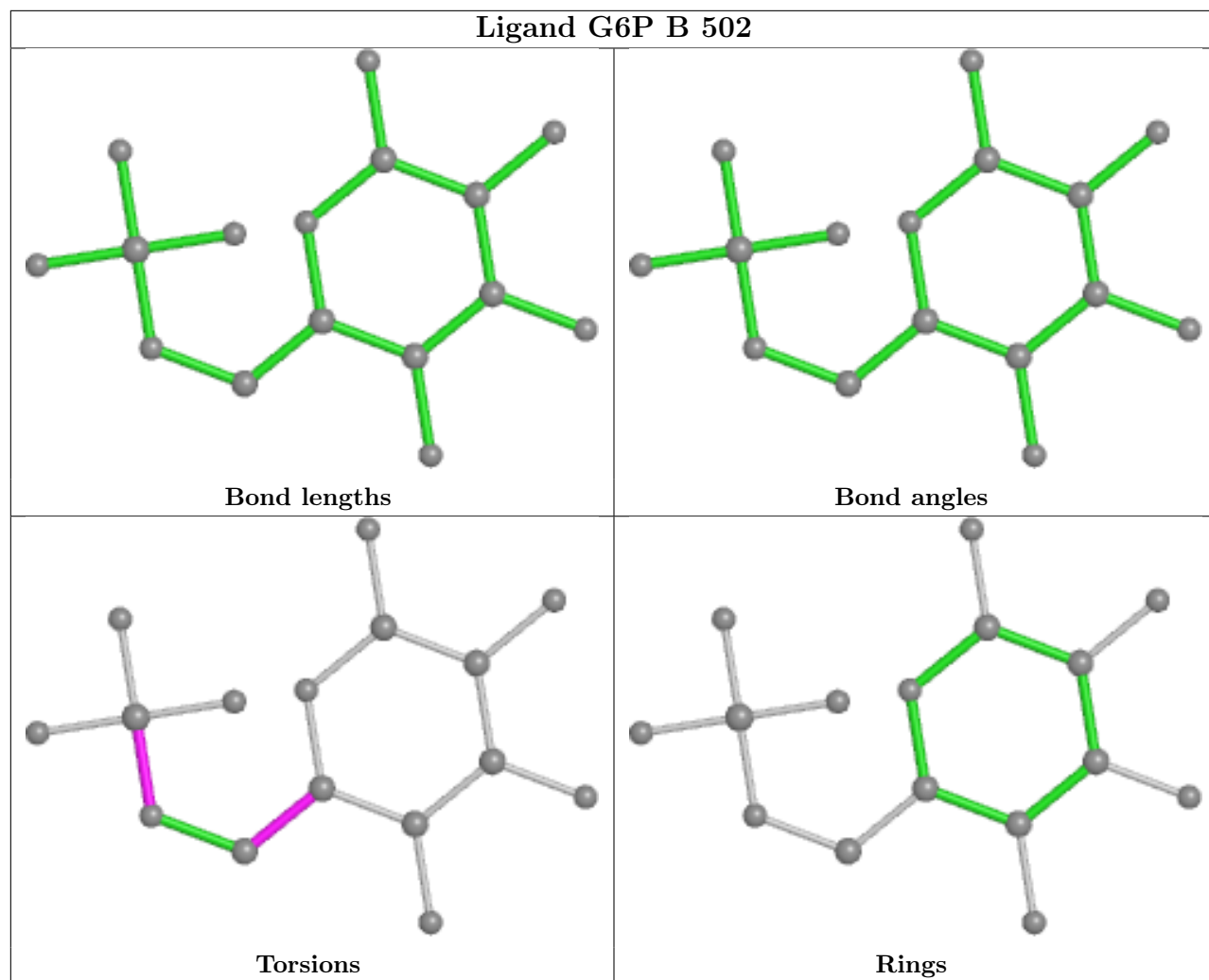


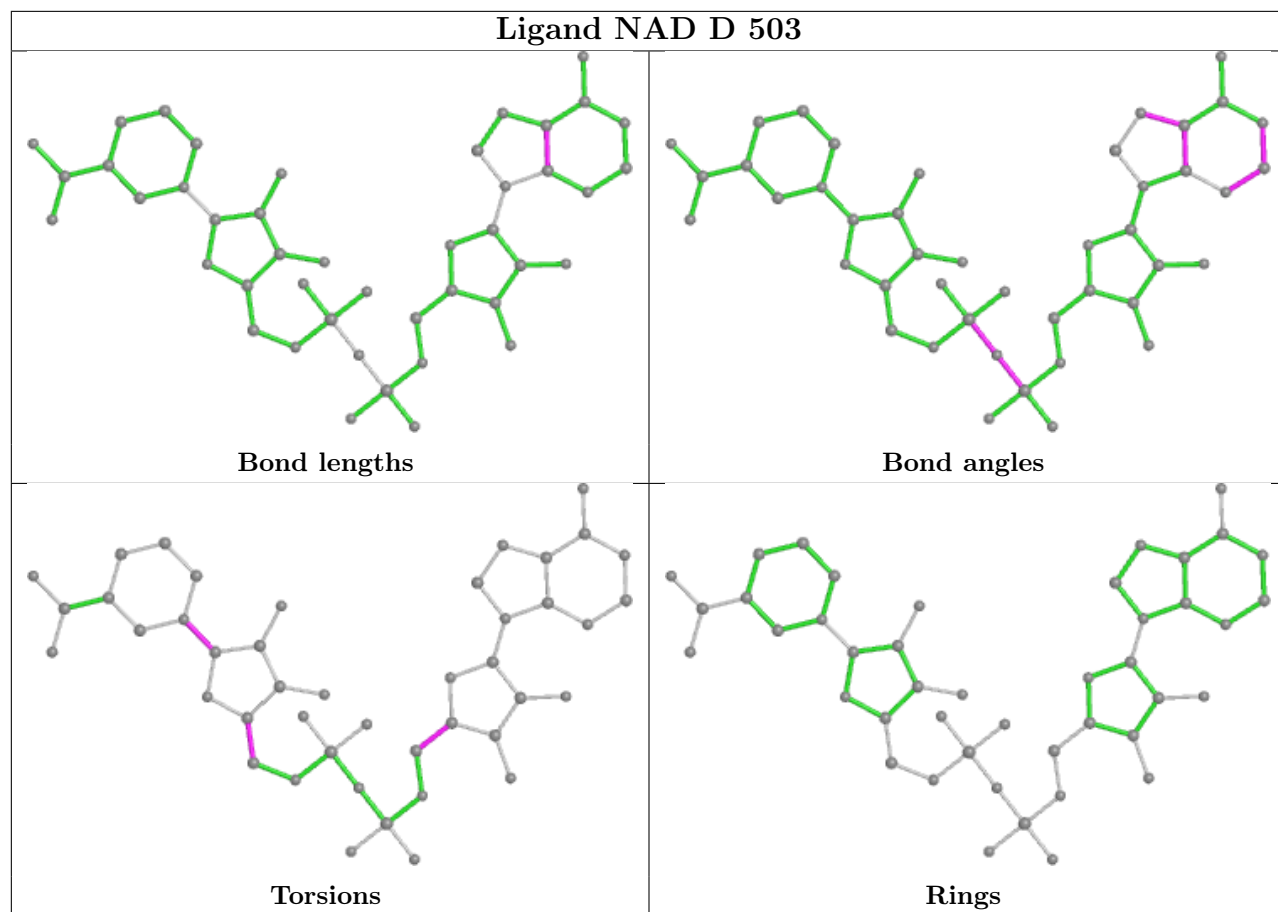
## Ligand G6P C 502



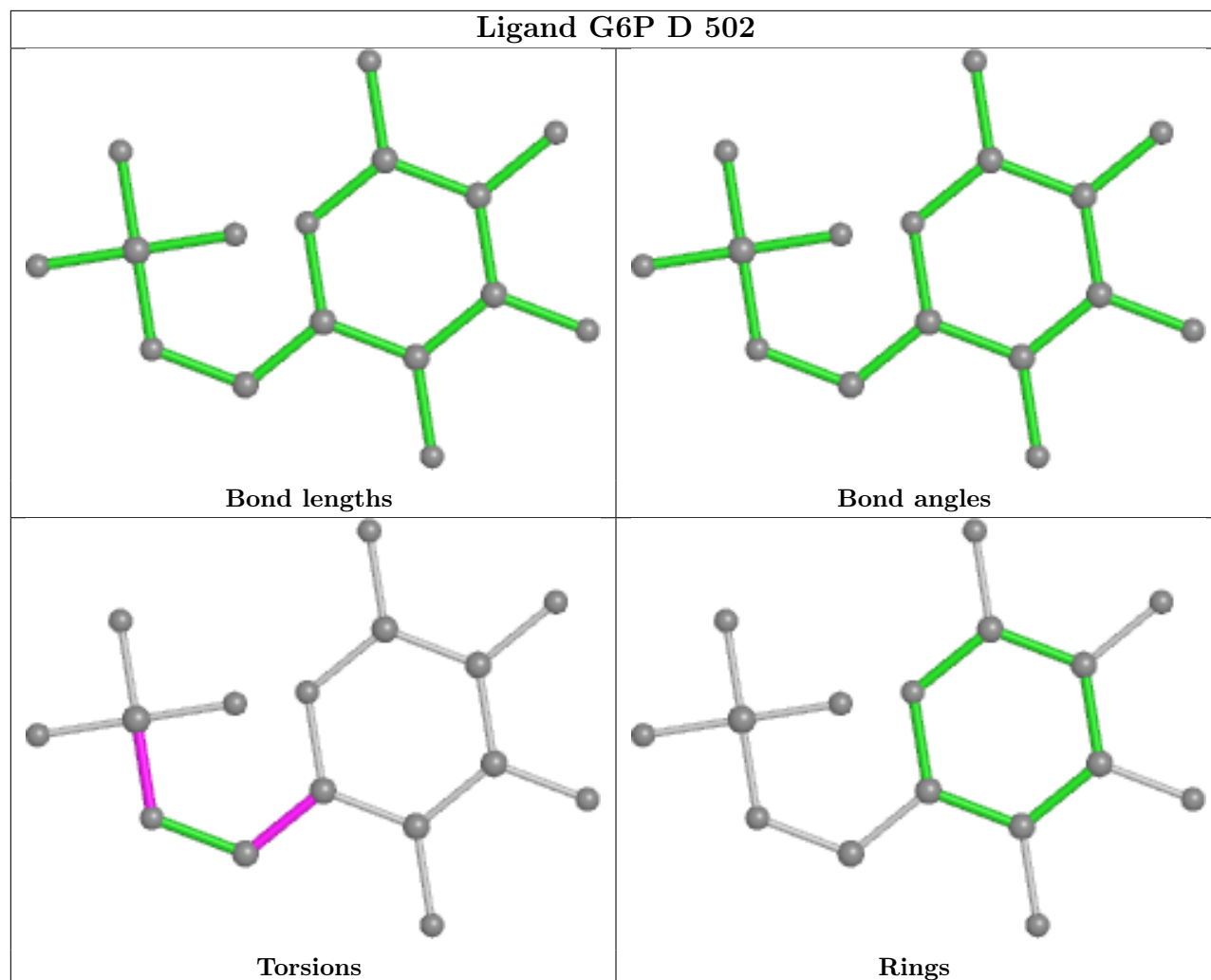


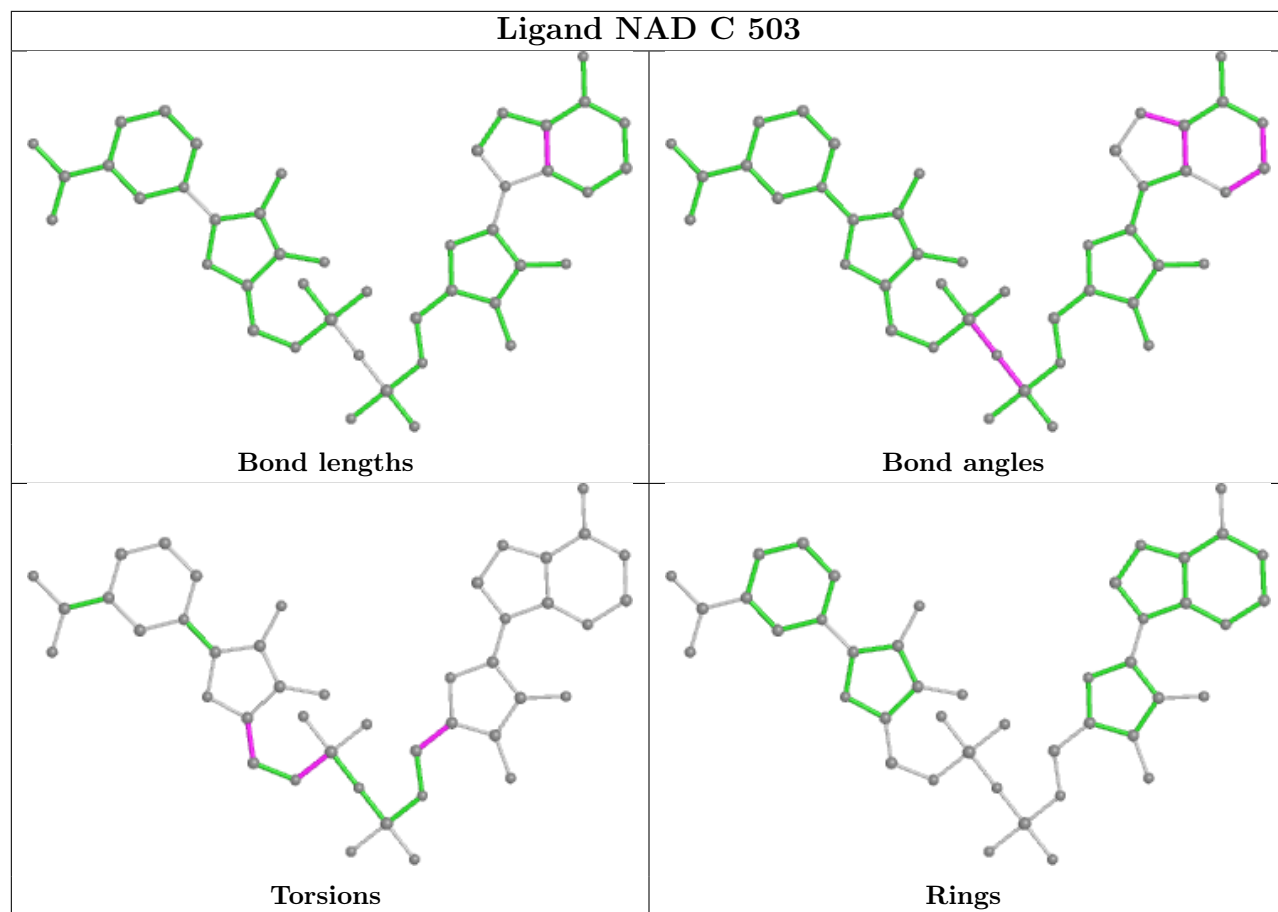






## Ligand G6P D 502





## 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	415/441 (94%)	0.13	6 (1%) 75 80	32, 43, 63, 79	0
1	B	423/441 (95%)	-0.00	2 (0%) 91 93	34, 44, 66, 99	0
1	C	421/441 (95%)	0.07	6 (1%) 75 80	32, 48, 75, 96	0
1	D	416/441 (94%)	0.01	10 (2%) 59 65	36, 51, 71, 90	0
All	All	1675/1764 (94%)	0.05	24 (1%) 75 80	32, 46, 69, 99	0

The worst 5 of 24 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	311	ARG	4.0
1	B	315	ILE	4.0
1	C	312	PHE	3.4
1	D	296	PHE	3.1
1	C	25	ARG	2.6

### 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.

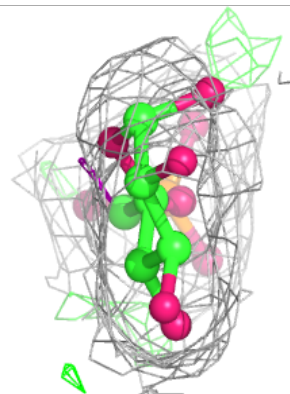
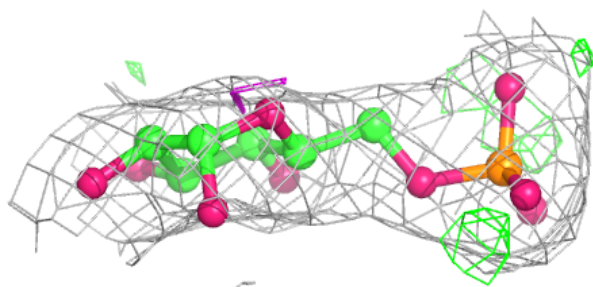
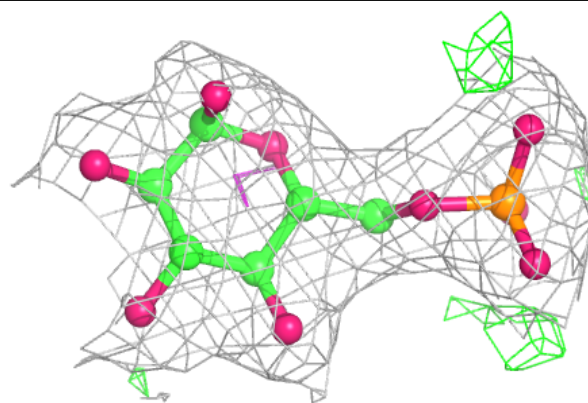


Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
5	GOL	C	505	6/6	0.81	0.23	83,86,94,95	0
5	GOL	C	504	6/6	0.92	0.15	56,62,65,70	0
5	GOL	B	504	6/6	0.93	0.14	52,60,63,64	0
3	G6P	D	502	16/16	0.94	0.12	46,49,55,65	0
3	G6P	C	502	16/16	0.94	0.10	38,42,46,56	0
4	NAD	C	503	44/44	0.94	0.11	41,49,53,55	0
4	NAD	B	503	44/44	0.95	0.10	38,42,46,48	0
4	NAD	D	503	44/44	0.95	0.11	47,51,56,57	0
4	NAD	A	503	44/44	0.96	0.10	35,40,43,43	0
3	G6P	B	502	16/16	0.96	0.11	31,41,48,57	0
3	G6P	A	502	16/16	0.96	0.14	36,40,46,48	0
2	MN	D	501	1/1	0.99	0.11	43,43,43,43	0
2	MN	C	501	1/1	0.99	0.08	36,36,36,36	0
2	MN	A	501	1/1	0.99	0.14	39,39,39,39	0
2	MN	B	501	1/1	1.00	0.08	33,33,33,33	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.

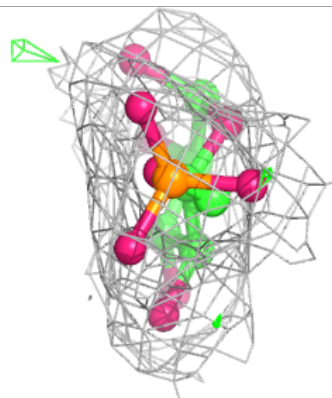
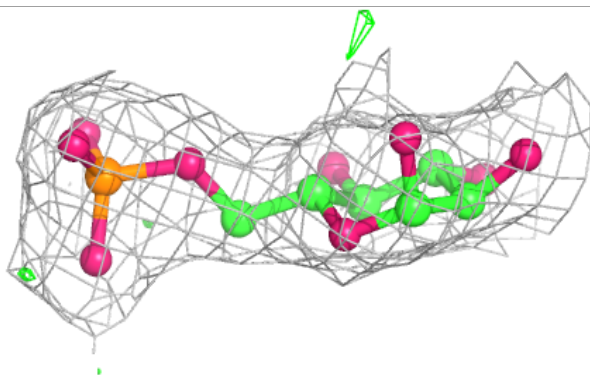
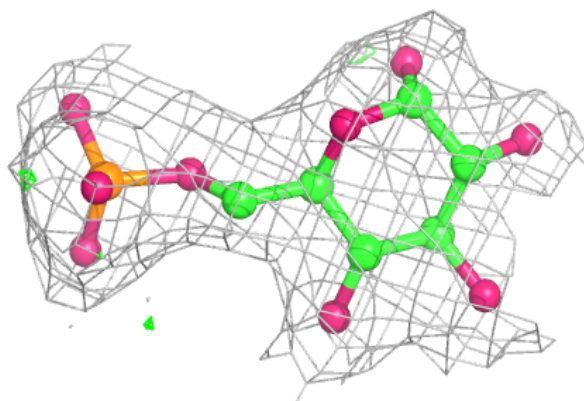
**Electron density around G6P D 502:**

2mF<sub>o</sub>-DF<sub>c</sub> (at 0.7 rmsd) in gray  
mF<sub>o</sub>-DF<sub>c</sub> (at 3 rmsd) in purple (negative)  
and green (positive)

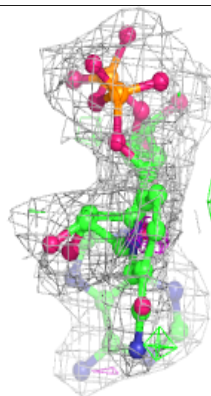
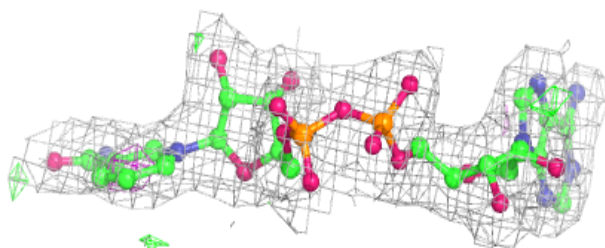
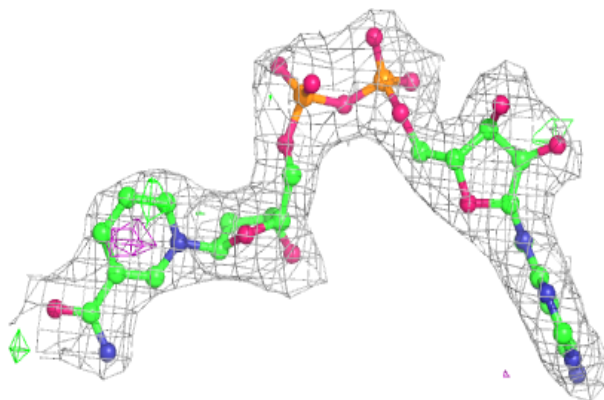


**Electron density around G6P C 502:**

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

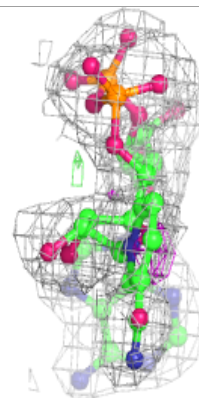
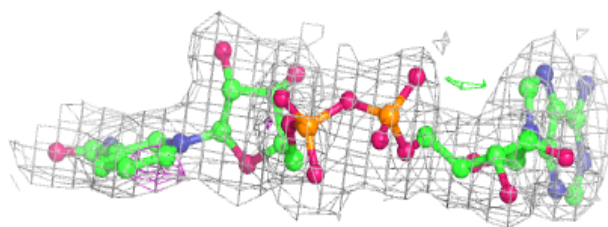
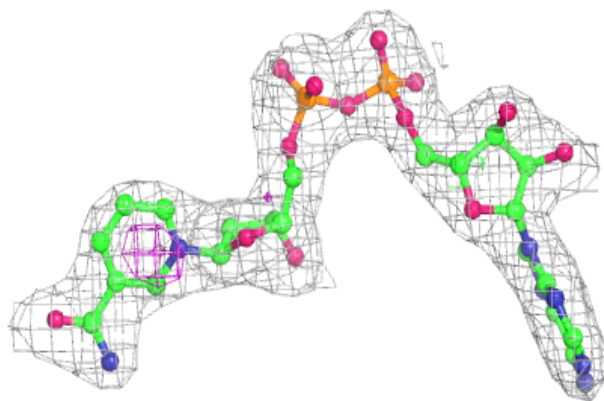
**Electron density around NAD C 503:**

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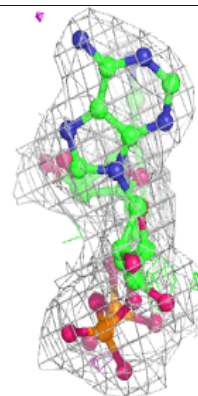
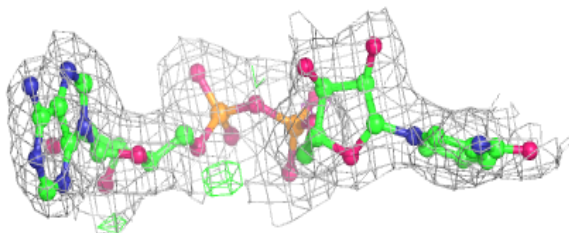
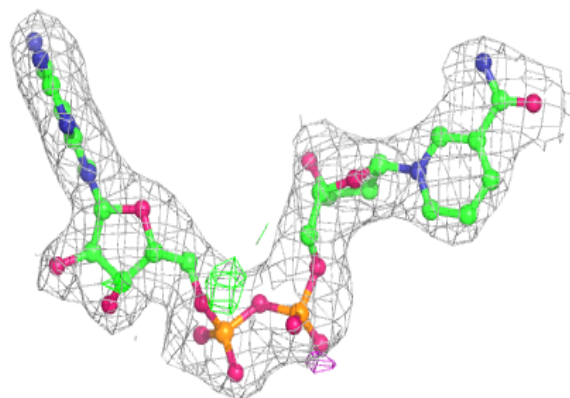


**Electron density around NAD B 503:**

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and green (positive)

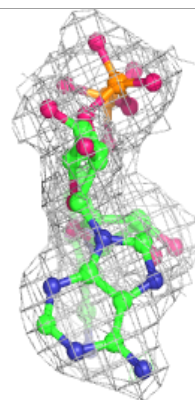
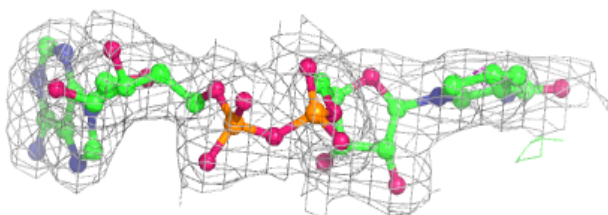
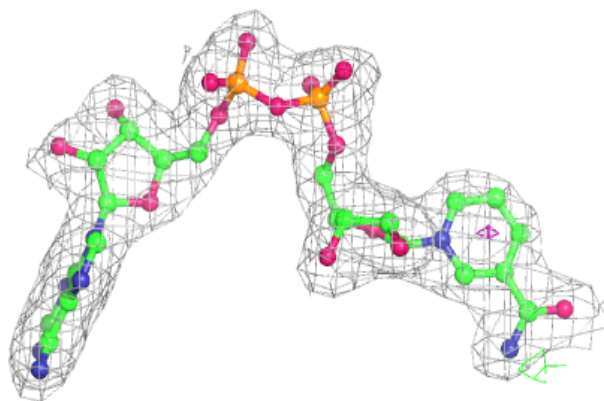
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and green (positive)

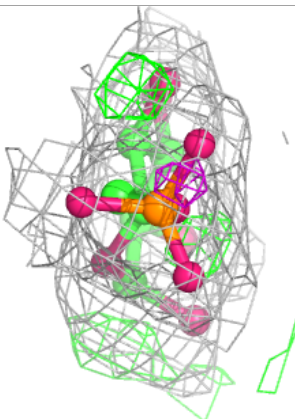
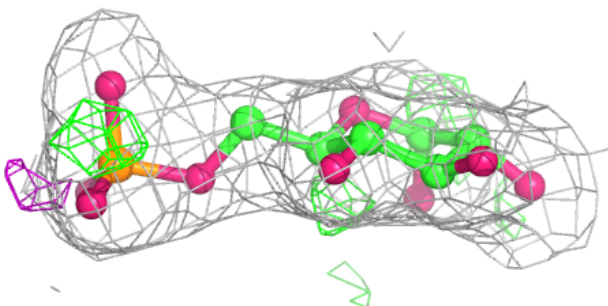
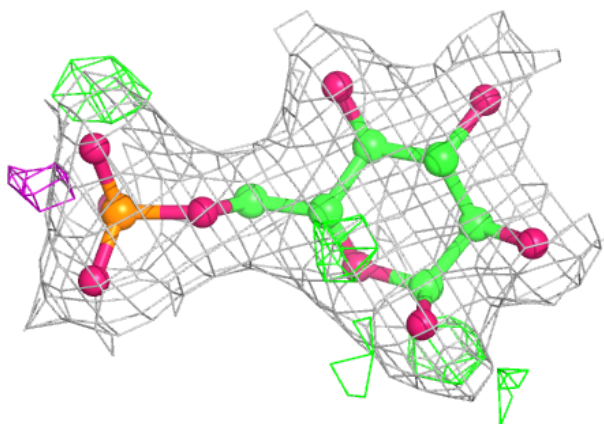


**Electron density around NAD A 503:**

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

**Electron density around G6P B 502:**

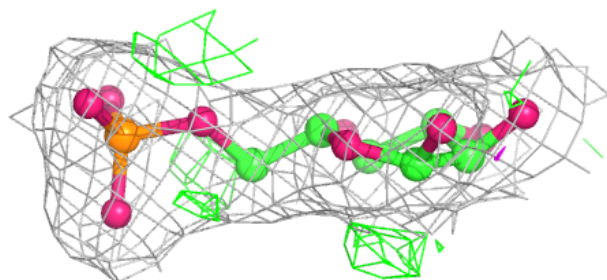
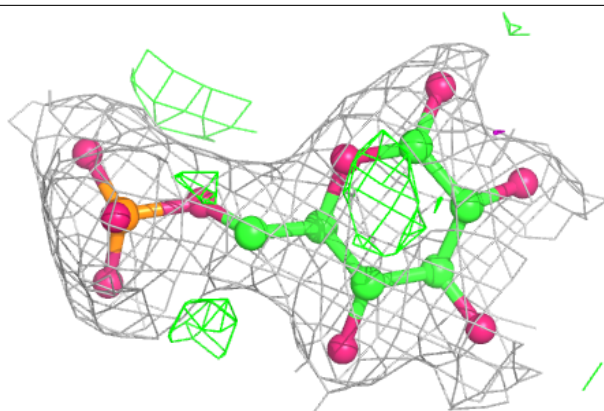
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 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)





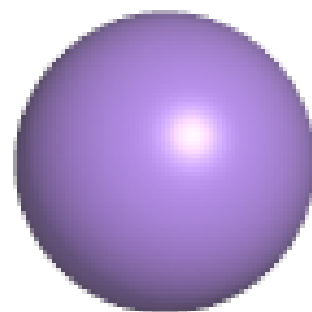
**Electron density around G6P A 502:**

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



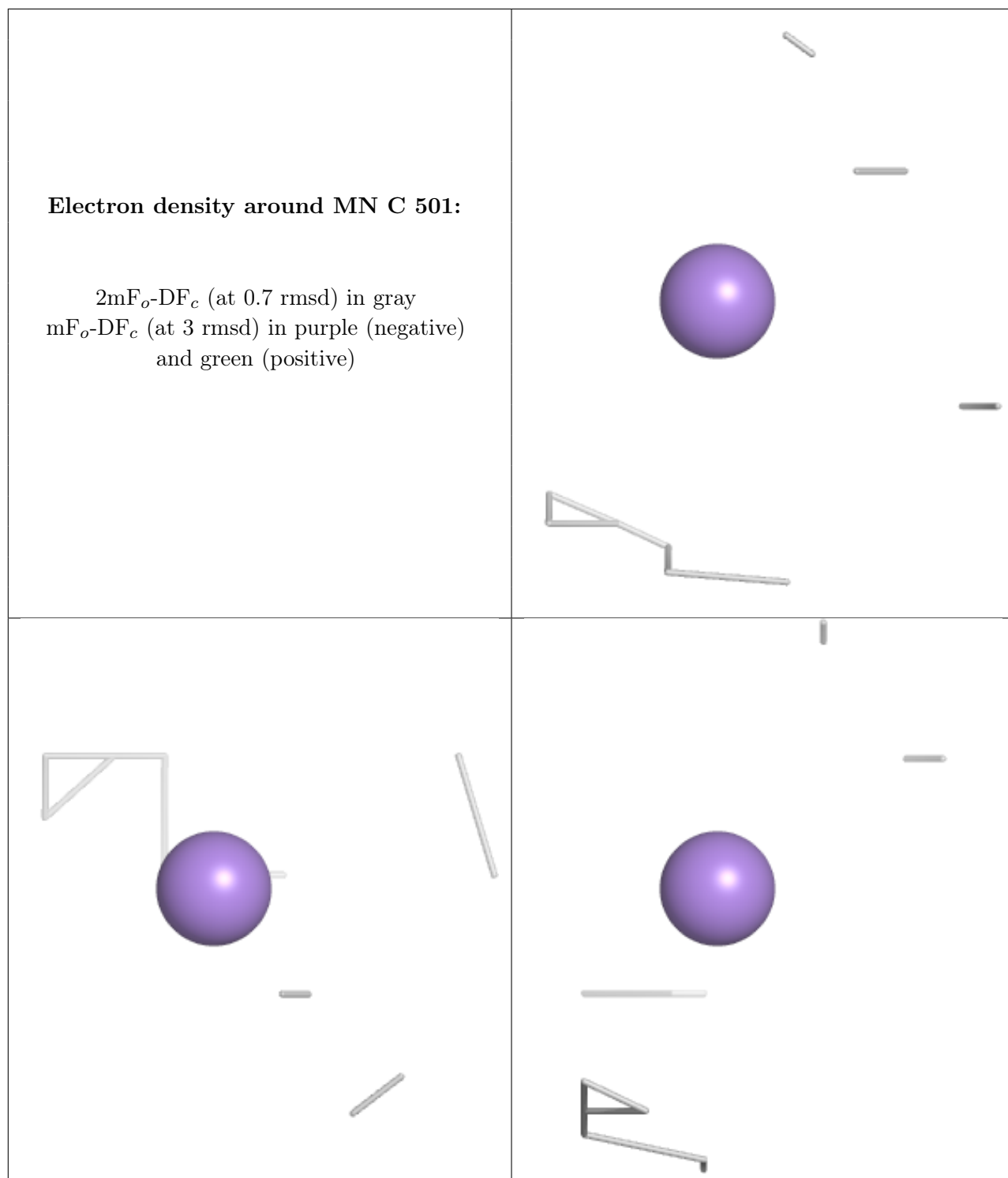
**Electron density around MN D 501:**

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mF<sub>o</sub>-DF<sub>c</sub> (at 3 rmsd) in purple (negative)  
and green (positive)



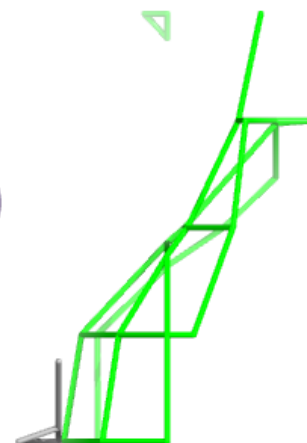
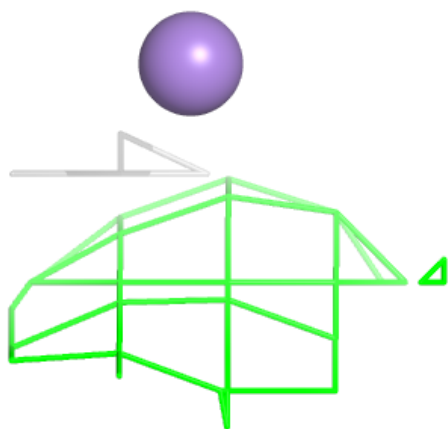
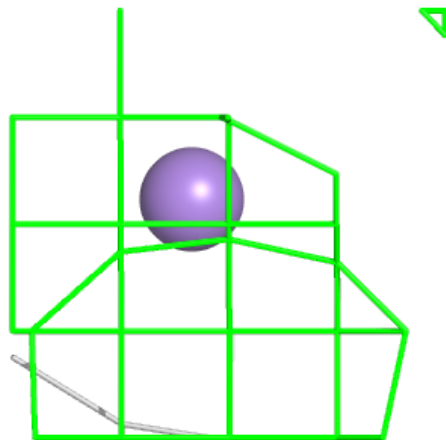
**Electron density around MN C 501:**

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and green (positive)

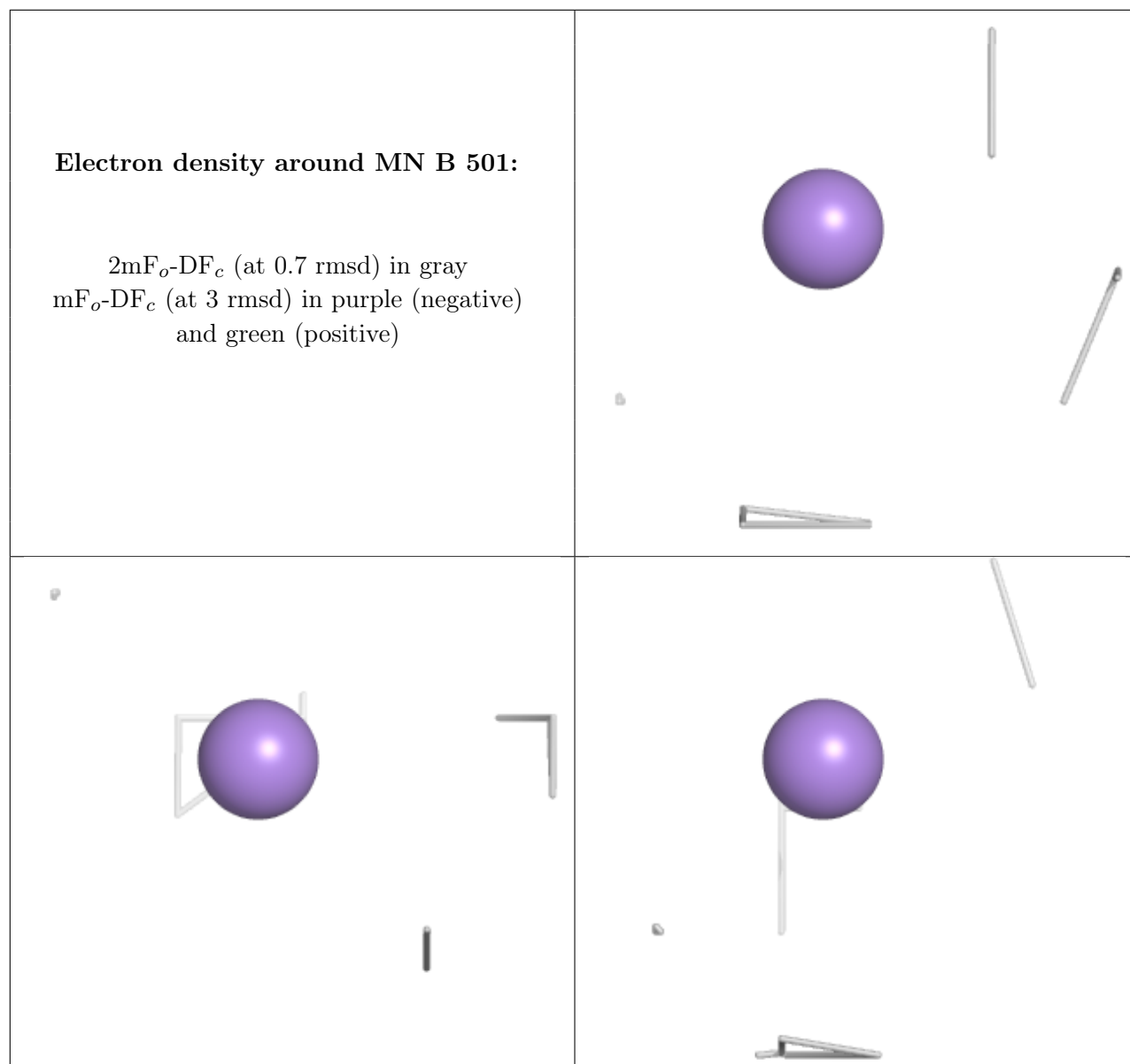


**Electron density around MN A 501:**

$2mF_o-DF_c$  (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.