



# Full wwPDB X-ray Structure Validation Report ⓘ

Dec 21, 2022 – 12:25 pm GMT

PDB ID : 7P18  
Title : Crystal structure of 3-ketosteroid delta1-dehydrogenase from *Sterolibacterium denitrificans* in complex with 1,4-androstadiene-3,17-dione  
Authors : Wojcik, P.; Mrugala, B.; Kurpiewska, K.; Szaleniec, M.  
Deposited on : 2021-07-01  
Resolution : 1.84 Å(reported)

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

We welcome your comments at [validation@mail.wwpdb.org](mailto:validation@mail.wwpdb.org)

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

---

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

MolProbity : 4.02b-467  
Mogul : 1.8.4, CSD as541be (2020)  
Xtriage (Phenix) : 1.13  
EDS : 2.31.3  
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) : 2.31.3

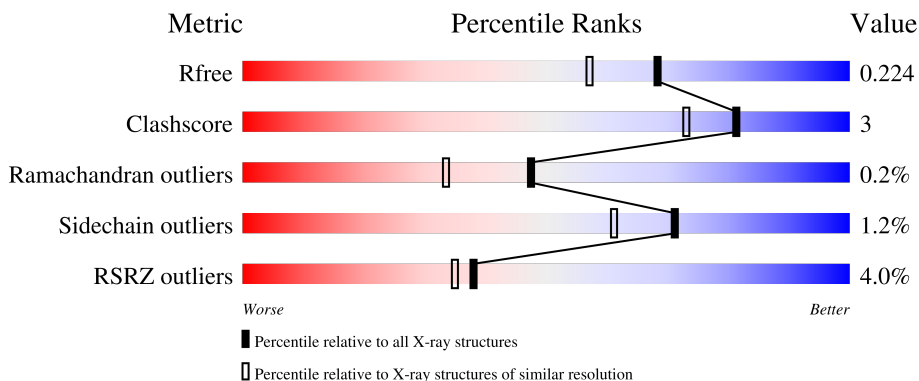
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 1.84 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	4003 (1.86-1.82)
Clashscore	141614	4233 (1.86-1.82)
Ramachandran outliers	138981	4185 (1.86-1.82)
Sidechain outliers	138945	4186 (1.86-1.82)
RSRZ outliers	127900	3957 (1.86-1.82)

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	561	
1	B	561	

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 criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
5	GOL	A	607	-	-	-	X

## 2 Entry composition i

There are 7 unique types of molecules in this entry. The entry contains 9665 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-oxosteroid 1-dehydrogenase.

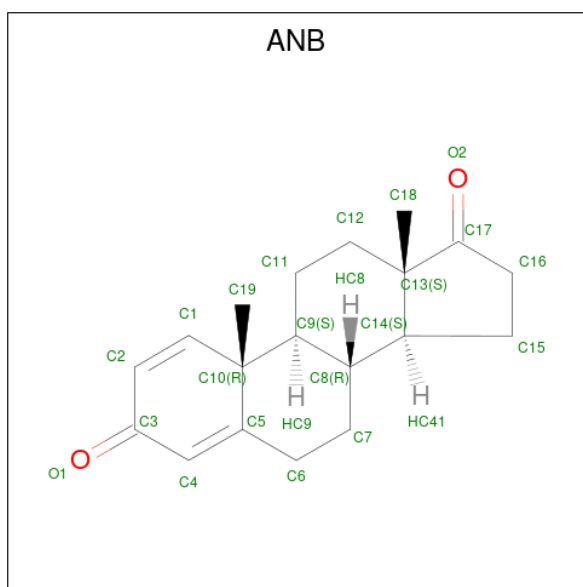
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	558	Total	C	N	O	S	0	4	0
			4322	2734	764	797	27			
1	B	557	Total	C	N	O	S	0	7	0
			4336	2739	770	801	26			

- Molecule 2 is FLAVIN-ADENINE DINUCLEOTIDE (three-letter code: FAD) (formula:  $C_{27}H_{33}N_9O_{15}P_2$ ) (labeled as "Ligand of Interest" by depositor).



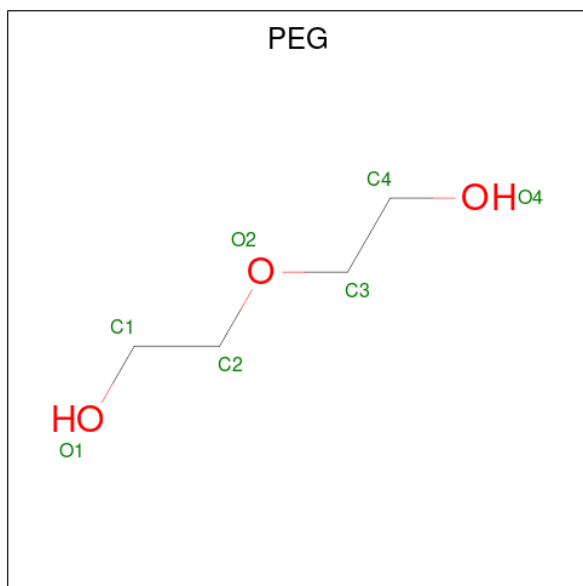
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	N	O	P		
2	A	1	Total	C	N	O	P	0	0
			53	27	9	15	2		
2	B	1	Total	C	N	O	P	0	0
			53	27	9	15	2		

- Molecule 3 is ANDROSTA-1,4-DIENE-3,17-DIONE (three-letter code: ANB) (formula:  $C_{19}H_{24}O_2$ ) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			21	19	2		

- Molecule 4 is DI(HYDROXYETHYL)ETHER (three-letter code: PEG) (formula:  $C_4H_{10}O_3$ ) (labeled as "Ligand of Interest" by depositor).



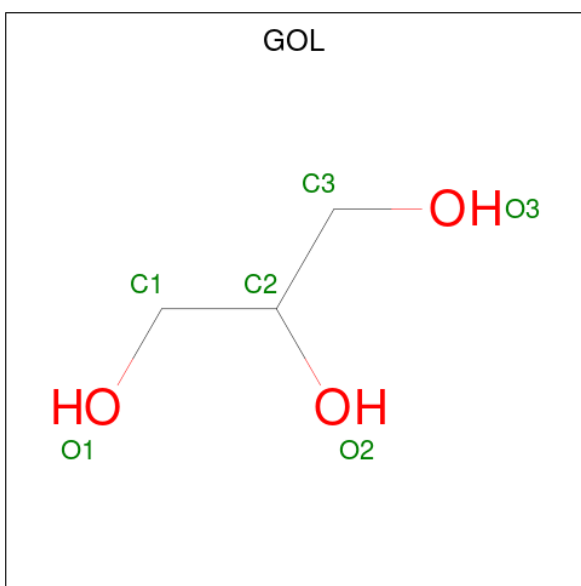
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			7	4	3		
4	A	1	Total	C	O	0	0
			7	4	3		

*Continued on next page...*

Continued from previous page...

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

- Molecule 5 is GLYCEROL (three-letter code: GOL) (formula: C<sub>3</sub>H<sub>8</sub>O<sub>3</sub>) (labeled as "Ligand of Interest" by depositor).



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

- Molecule 6 is SODIUM ION (three-letter code: NA) (formula: Na) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	1	Total	Na	0	0
			1	1		

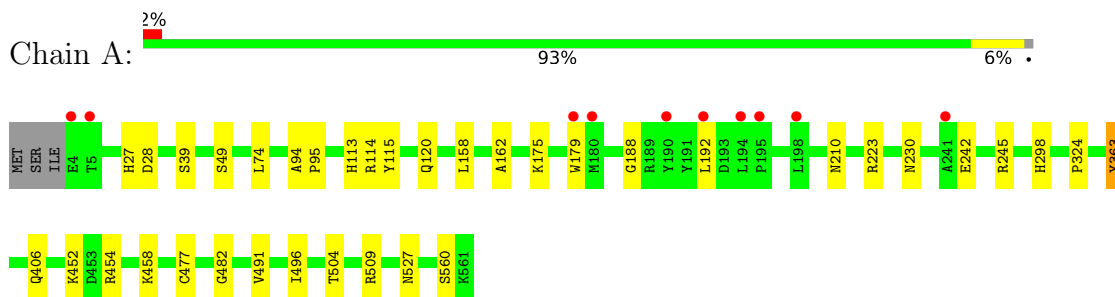
- Molecule 7 is water.

<b>Mol</b>	<b>Chain</b>	<b>Residues</b>	<b>Atoms</b>		<b>ZeroOcc</b>	<b>AltConf</b>
7	A	521	Total 521	O 521	0	0
7	B	305	Total 305	O 305	0	0

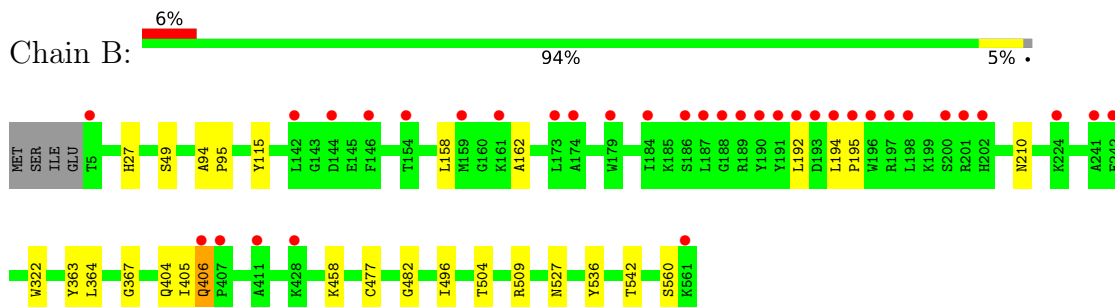
### 3 Residue-property plots [i](#)

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-oxosteroid 1-dehydrogenase



- Molecule 1: 3-oxosteroid 1-dehydrogenase





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	130.37Å 130.37Å 139.74Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	47.71 – 1.84 47.66 – 1.84	Depositor EDS
% Data completeness (in resolution range)	99.5 (47.71-1.84) 99.5 (47.66-1.84)	Depositor EDS
$R_{merge}$	0.11	Depositor
$R_{sym}$	0.13	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.60 (at 1.84Å)	Xtrriage
Refinement program	REFMAC 5.8.0258	Depositor
R, $R_{free}$	0.185 , 0.214 0.195 , 0.224	Depositor DCC
$R_{free}$ test set	2100 reflections (1.77%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	29.4	Xtrriage
Anisotropy	0.055	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.31 , 45.3	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.34$	Xtrriage
Estimated twinning fraction	0.016 for -h,-k,l	Xtrriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	9665	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	38.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.76% 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: PEG, NA, GOL, FAD, ANB

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	0.74	0/4426	0.83	0/5998
1	B	0.71	0/4440	0.81	0/6018
All	All	0.72	0/8866	0.82	0/12016

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

### 5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	4322	0	4250	28	0
1	B	4336	0	4247	15	0
2	A	53	0	31	2	0
2	B	53	0	31	3	0
3	A	21	0	24	1	0
4	A	21	0	30	5	0
4	B	14	0	20	0	0
5	A	18	0	24	4	0
6	A	1	0	0	0	0
7	A	521	0	0	6	0
7	B	305	0	0	4	0
All	All	9665	0	8657	45	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 (45) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:560[A]:SER:OG	7:B:701:HOH:O	1.90	0.88
1:B:406:GLN:O	1:B:406:GLN:HG3	1.92	0.70
1:A:242:GLU:HG3	7:A:1035:HOH:O	1.98	0.64
1:B:496:ILE:HB	2:B:601:FAD:H9	1.81	0.62
1:A:188:GLY:O	1:A:192:LEU:HG	2.00	0.61
1:A:406:GLN:NE2	7:A:703:HOH:O	2.36	0.59
1:A:28:ASP:HB3	5:A:607:GOL:H2	1.86	0.58
1:A:27:HIS:HD2	7:A:1050:HOH:O	1.89	0.55
1:A:158:LEU:HB3	1:A:162:ALA:HB3	1.89	0.54
1:A:496:ILE:HB	2:A:601:FAD:H9	1.89	0.54
1:A:120:GLN:HB3	4:A:603:PEG:H12	1.89	0.54
1:A:560:SER:HB2	7:A:724:HOH:O	2.07	0.54
1:A:179:TRP:CD2	5:A:606:GOL:O2	2.60	0.54
1:A:113:HIS:HB2	4:A:604:PEG:H31	1.91	0.53
1:A:452[B]:LYS:HG2	1:A:454:ARG:NH1	2.24	0.52
1:B:158:LEU:HB3	1:B:162:ALA:HB3	1.91	0.52
1:B:192:LEU:HD23	1:B:192:LEU:O	2.10	0.51
1:B:27:HIS:HD2	7:B:905:HOH:O	1.93	0.51
1:A:363:TYR:OH	3:A:602:ANB:C2	2.62	0.47
1:A:94:ALA:HB3	1:A:95:PRO:HD3	1.96	0.46
1:A:245:ARG:HD3	7:B:851:HOH:O	2.16	0.46
1:A:175:LYS:HB3	5:A:606:GOL:H31	1.97	0.45
1:B:49:SER:HA	2:B:601:FAD:C6	2.48	0.44
1:B:94:ALA:HB3	1:B:95:PRO:HD3	2.00	0.44
1:B:536:TYR:HB3	7:B:872:HOH:O	2.17	0.43
1:A:120:GLN:CB	4:A:603:PEG:H12	2.48	0.43
1:B:322:TRP:HB3	1:B:367:GLY:HA3	2.00	0.43
1:A:230:ASN:ND2	7:A:707:HOH:O	2.41	0.43
4:A:604:PEG:H32	7:A:860:HOH:O	2.19	0.43
1:B:194:LEU:CB	1:B:195:PRO:CD	2.97	0.43
1:A:39:SER:HB3	1:A:298:HIS:CD2	2.55	0.42
1:A:39:SER:HB2	1:A:298:HIS:NE2	2.34	0.42
2:B:601:FAD:H9	2:B:601:FAD:H1'1	1.83	0.42
1:B:458:LYS:HE2	1:B:477:CYS:HB3	2.01	0.42
1:A:27:HIS:CE1	1:A:223:ARG:HB3	2.55	0.41
1:A:504:THR:HA	1:A:509:ARG:O	2.21	0.41
1:A:179:TRP:CE3	5:A:606:GOL:O2	2.73	0.41

*Continued on next page...*

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:458:LYS:HE2	1:A:477:CYS:HB3	2.03	0.41
1:A:324:PRO:HB2	1:A:491:VAL:HG13	2.03	0.41
1:A:49:SER:HA	2:A:601:FAD:C6	2.51	0.41
1:B:364:LEU:HD22	1:B:536:TYR:HB2	2.03	0.40
1:A:74:LEU:HD23	1:A:74:LEU:HA	1.91	0.40
1:B:404[B]:GLN:HG3	1:B:405:ILE:HG23	2.03	0.40
1:A:114:ARG:CD	4:A:604:PEG:H22	2.51	0.40
1:B:504:THR:HA	1:B:509:ARG: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	560/561 (100%)	545 (97%)	14 (2%)	1 (0%)	47 33
1	B	562/561 (100%)	545 (97%)	16 (3%)	1 (0%)	47 33
All	All	1122/1122 (100%)	1090 (97%)	30 (3%)	2 (0%)	47 33

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	482	GLY
1	B	482	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	436/436 (100%)	432 (99%)	4 (1%)	78	71
1	B	436/436 (100%)	430 (99%)	6 (1%)	67	55
All	All	872/872 (100%)	862 (99%)	10 (1%)	71	64

All (10) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	115	TYR
1	A	210	ASN
1	A	363	TYR
1	A	527	ASN
1	B	115	TYR
1	B	210	ASN
1	B	363	TYR
1	B	406	GLN
1	B	527	ASN
1	B	542	THR

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

Mol	Chain	Res	Type
1	A	218	HIS
1	A	243	ASN
1	B	6	ASN
1	B	243	ASN

### 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

Of 12 ligands modelled in this entry, 1 is 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	A	608	-	5,5,5	0.13	0	5,5,5	0.14	0
2	FAD	B	601	-	53,58,58	0.86	1 (1%)	68,89,89	1.06	7 (10%)
4	PEG	A	605	-	6,6,6	0.42	0	5,5,5	0.29	0
5	GOL	A	607	-	5,5,5	0.11	0	5,5,5	0.30	0
4	PEG	B	603	-	6,6,6	0.16	0	5,5,5	0.11	0
3	ANB	A	602	-	24,24,24	0.41	0	39,39,39	1.06	2 (5%)
2	FAD	A	601	-	53,58,58	0.70	1 (1%)	68,89,89	0.91	2 (2%)
4	PEG	B	602	-	6,6,6	0.20	0	5,5,5	0.13	0
5	GOL	A	606	-	5,5,5	0.12	0	5,5,5	0.26	0
4	PEG	A	603	-	6,6,6	0.31	0	5,5,5	0.25	0
4	PEG	A	604	-	6,6,6	0.28	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
5	GOL	A	608	-	-	3/4/4/4	-
2	FAD	B	601	-	-	2/30/50/50	0/6/6/6
4	PEG	A	605	-	-	4/4/4/4	-
5	GOL	A	607	-	-	3/4/4/4	-
4	PEG	B	603	-	-	2/4/4/4	-
3	ANB	A	602	-	-	-	0/4/4/4
2	FAD	A	601	-	-	1/30/50/50	0/6/6/6
4	PEG	B	602	-	-	1/4/4/4	-
5	GOL	A	606	-	-	3/4/4/4	-
4	PEG	A	603	-	-	2/4/4/4	-
4	PEG	A	604	-	-	1/4/4/4	-

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	601	FAD	O2-C2	2.31	1.28	1.24
2	A	601	FAD	C2-N3	-2.10	1.34	1.39

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	601	FAD	O2P-P-O1P	2.98	126.97	112.24
2	B	601	FAD	C5A-C6A-N6A	2.92	124.80	120.35
3	A	602	ANB	C2-C3-C4	-2.74	114.33	117.13
3	A	602	ANB	C10-C9-C8	-2.71	109.79	112.08
2	A	601	FAD	C5A-C6A-N6A	2.60	124.30	120.35
2	B	601	FAD	O5'-P-O1P	-2.51	99.25	109.07
2	B	601	FAD	C1'-C2'-C3'	2.49	116.74	109.79
2	B	601	FAD	O2A-PA-O1A	2.37	123.96	112.24
2	B	601	FAD	C4X-C4-N3	2.11	118.56	113.19
2	A	601	FAD	O2P-P-O1P	2.07	122.47	112.24
2	B	601	FAD	O3B-C3B-C2B	2.02	118.36	111.82

There are no chirality outliers.

All (22) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	A	608	GOL	C1-C2-C3-O3
4	A	605	PEG	O1-C1-C2-O2
5	A	607	GOL	O1-C1-C2-C3
5	A	607	GOL	O1-C1-C2-O2
2	B	601	FAD	O4B-C4B-C5B-O5B
4	A	604	PEG	O1-C1-C2-O2
5	A	608	GOL	O2-C2-C3-O3
4	A	603	PEG	O2-C3-C4-O4
4	B	603	PEG	O1-C1-C2-O2
5	A	606	GOL	O2-C2-C3-O3
4	A	603	PEG	C4-C3-O2-C2
4	A	605	PEG	C1-C2-O2-C3
5	A	606	GOL	O1-C1-C2-O2
2	A	601	FAD	O4B-C4B-C5B-O5B
4	A	605	PEG	C4-C3-O2-C2
4	A	605	PEG	O2-C3-C4-O4
4	B	602	PEG	O2-C3-C4-O4
5	A	608	GOL	O1-C1-C2-C3
4	B	603	PEG	C4-C3-O2-C2

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type	Atoms
2	B	601	FAD	C3B-C4B-C5B-O5B
5	A	607	GOL	O2-C2-C3-O3
5	A	606	GOL	C1-C2-C3-O3

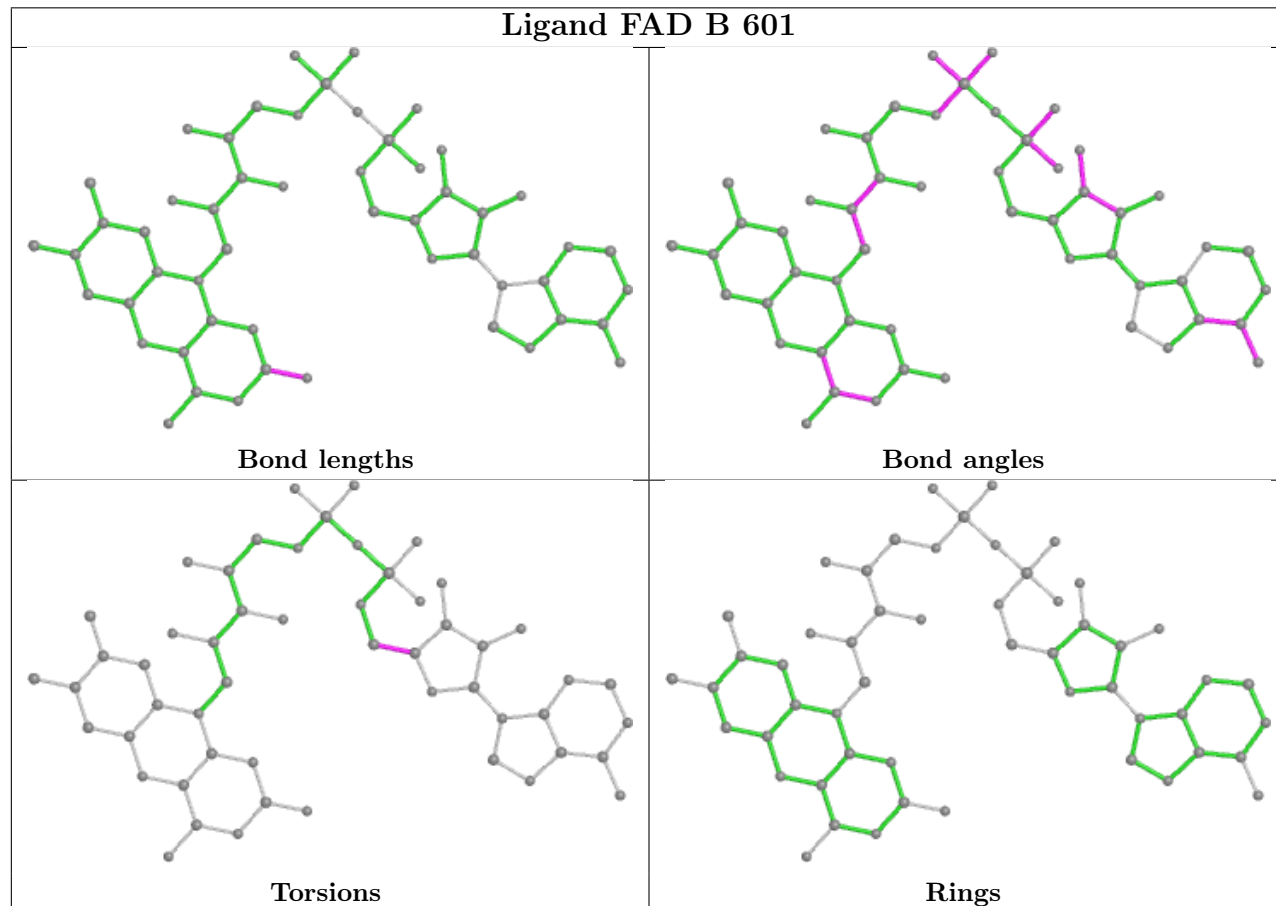
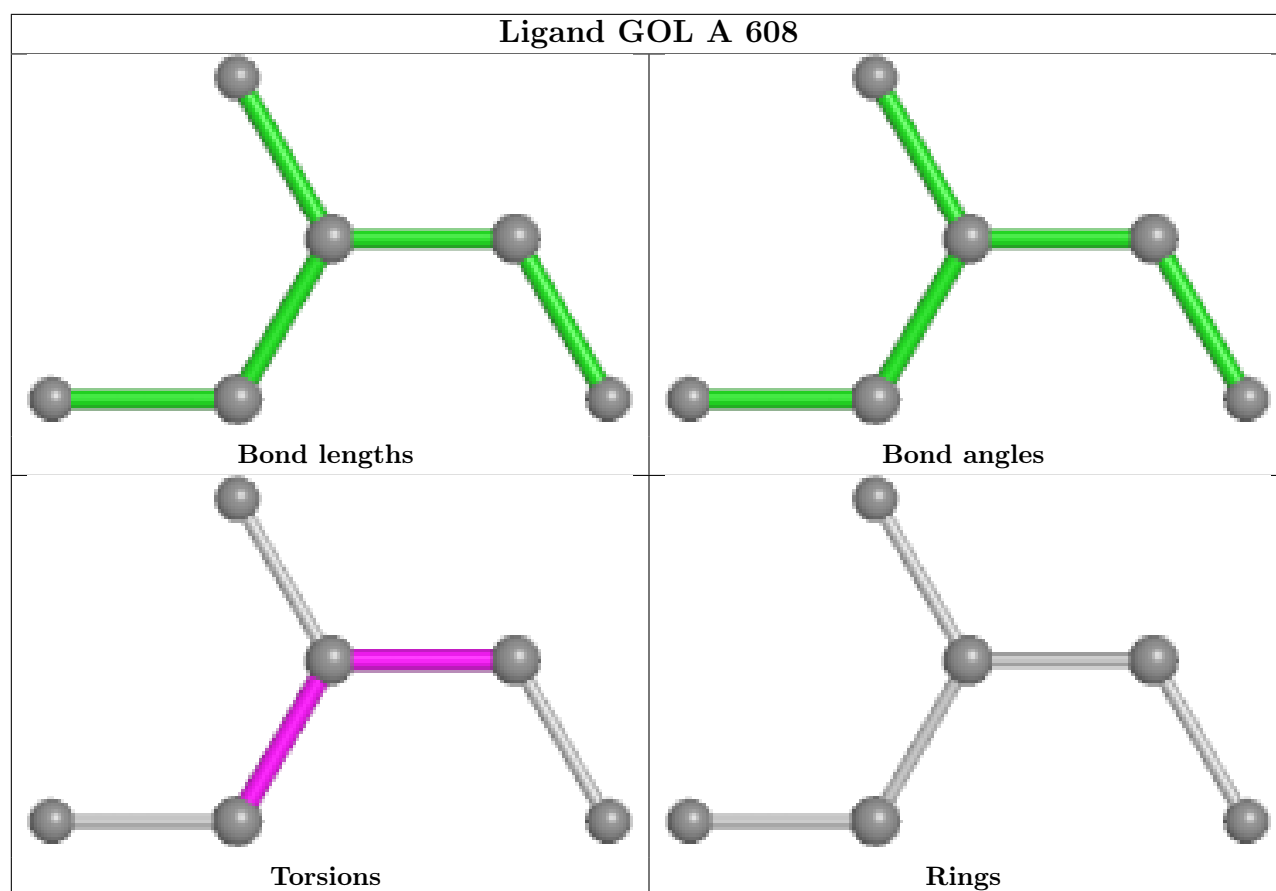
There are no ring outliers.

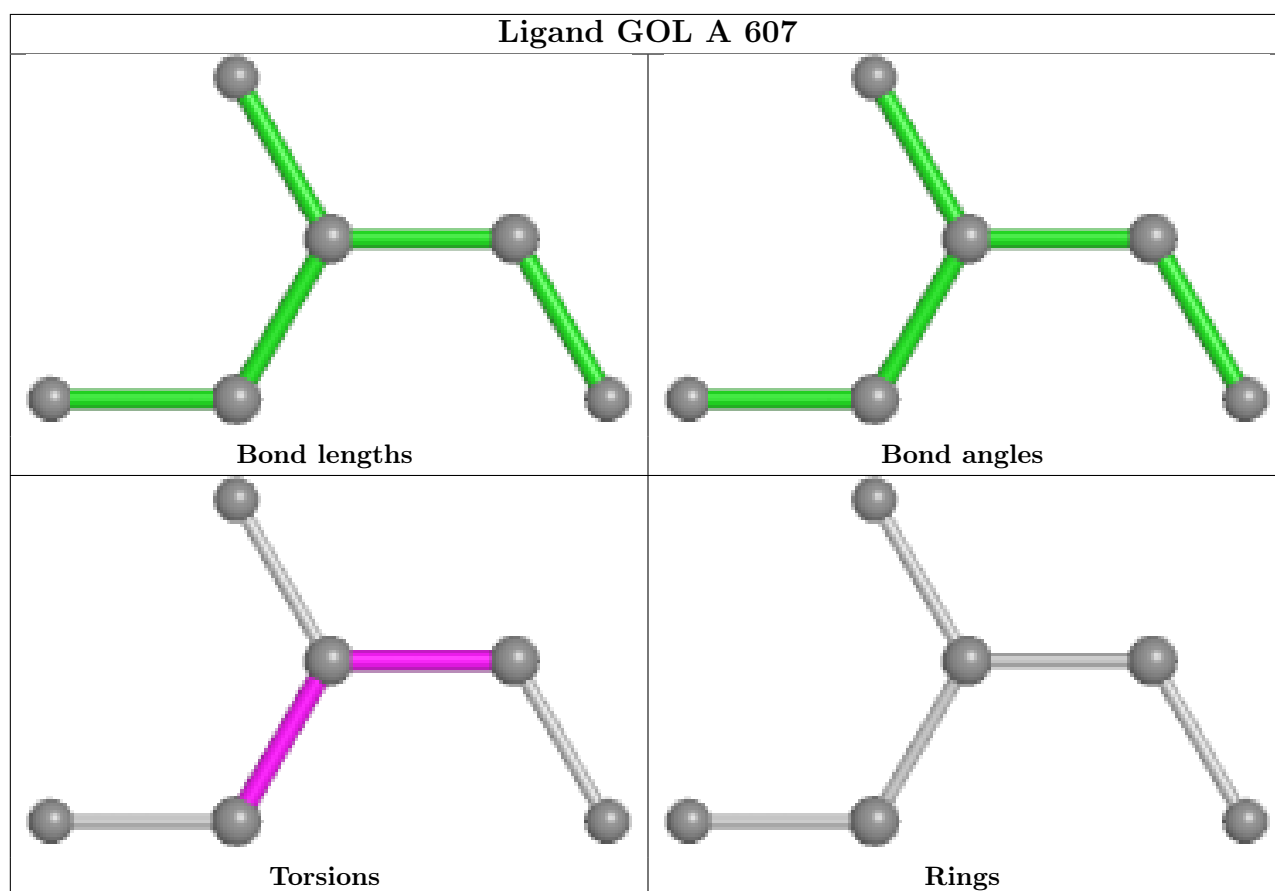
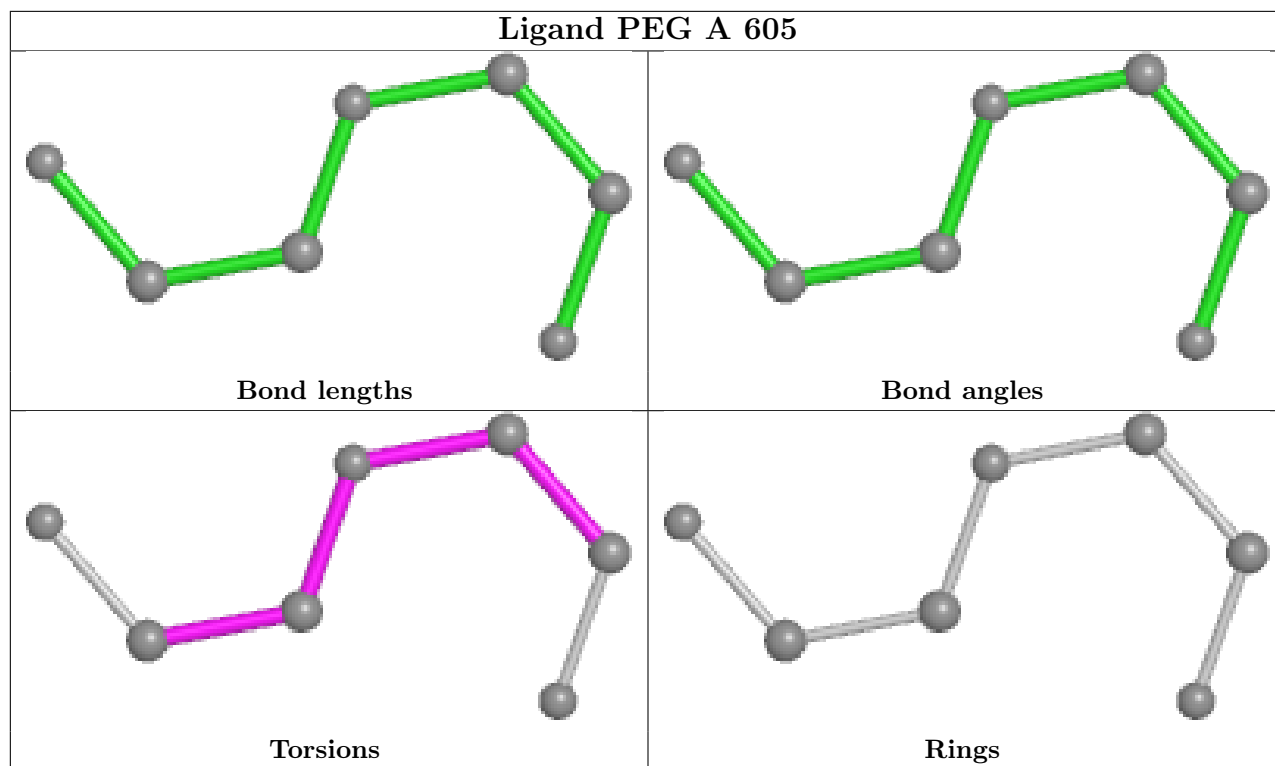
7 monomers are involved in 15 short contacts:

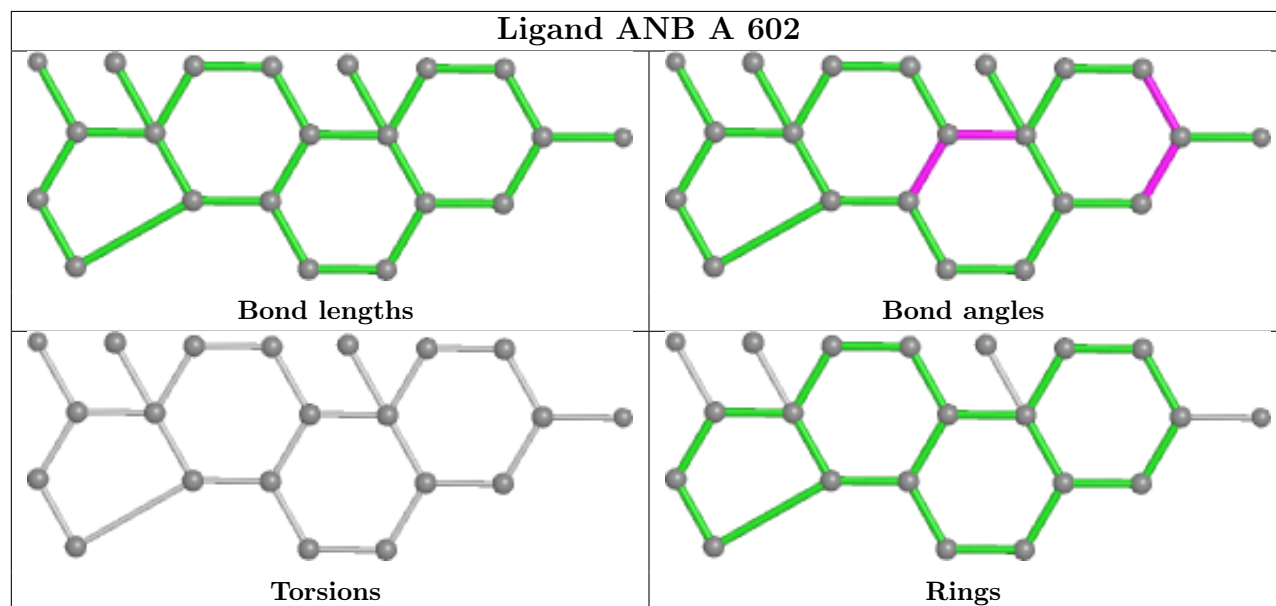
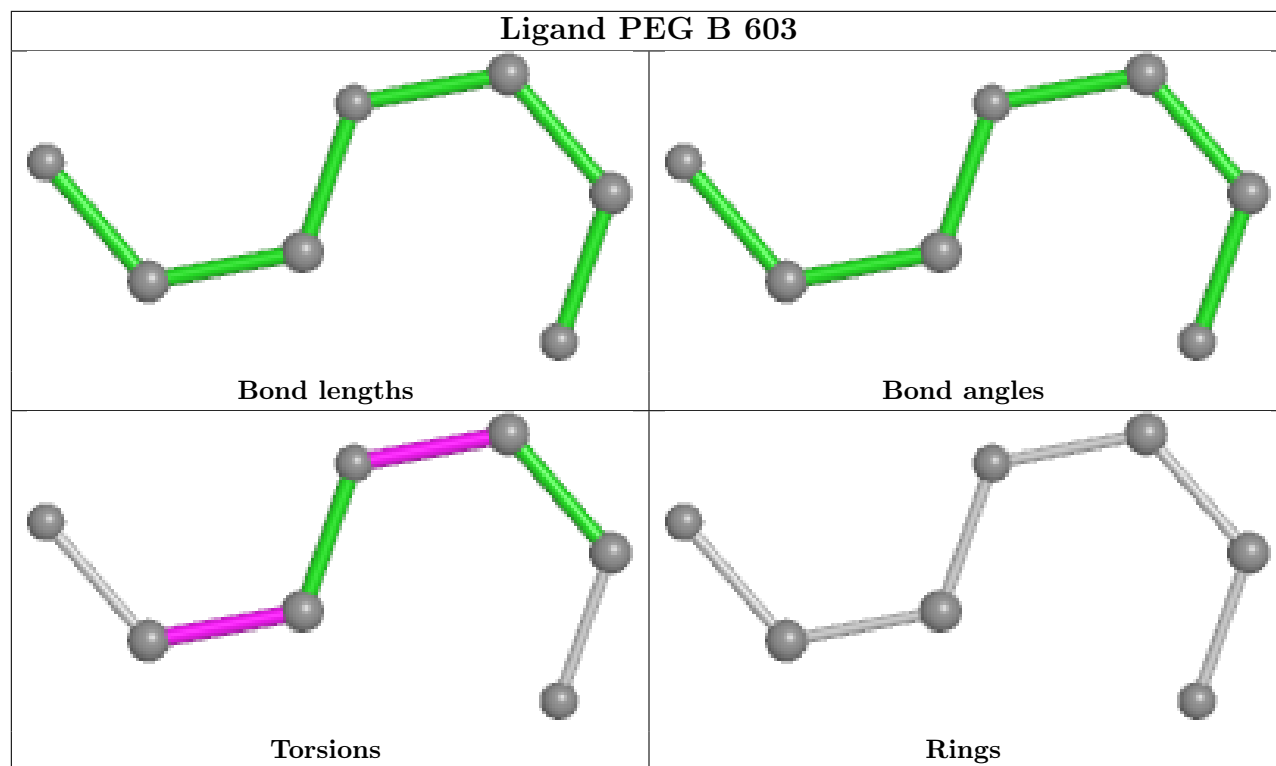
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	601	FAD	3	0
5	A	607	GOL	1	0
3	A	602	ANB	1	0
2	A	601	FAD	2	0
5	A	606	GOL	3	0
4	A	603	PEG	2	0
4	A	604	PEG	3	0

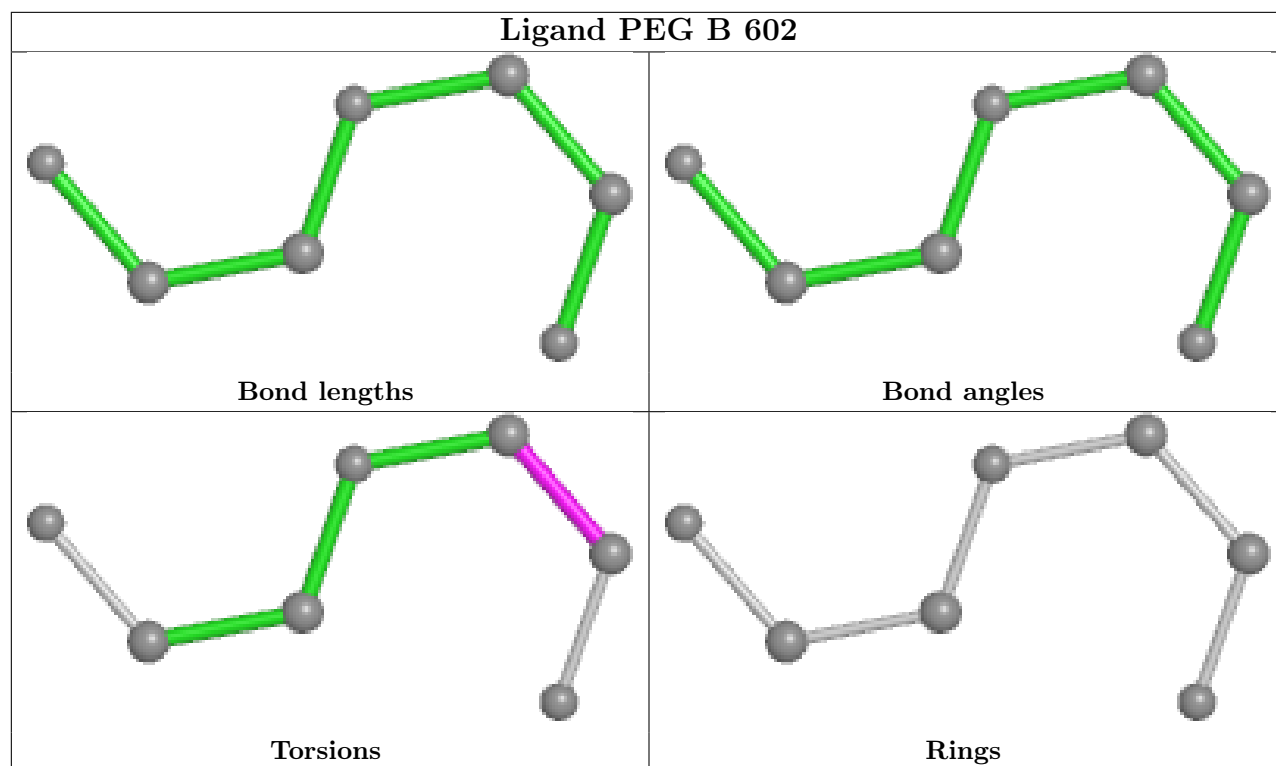
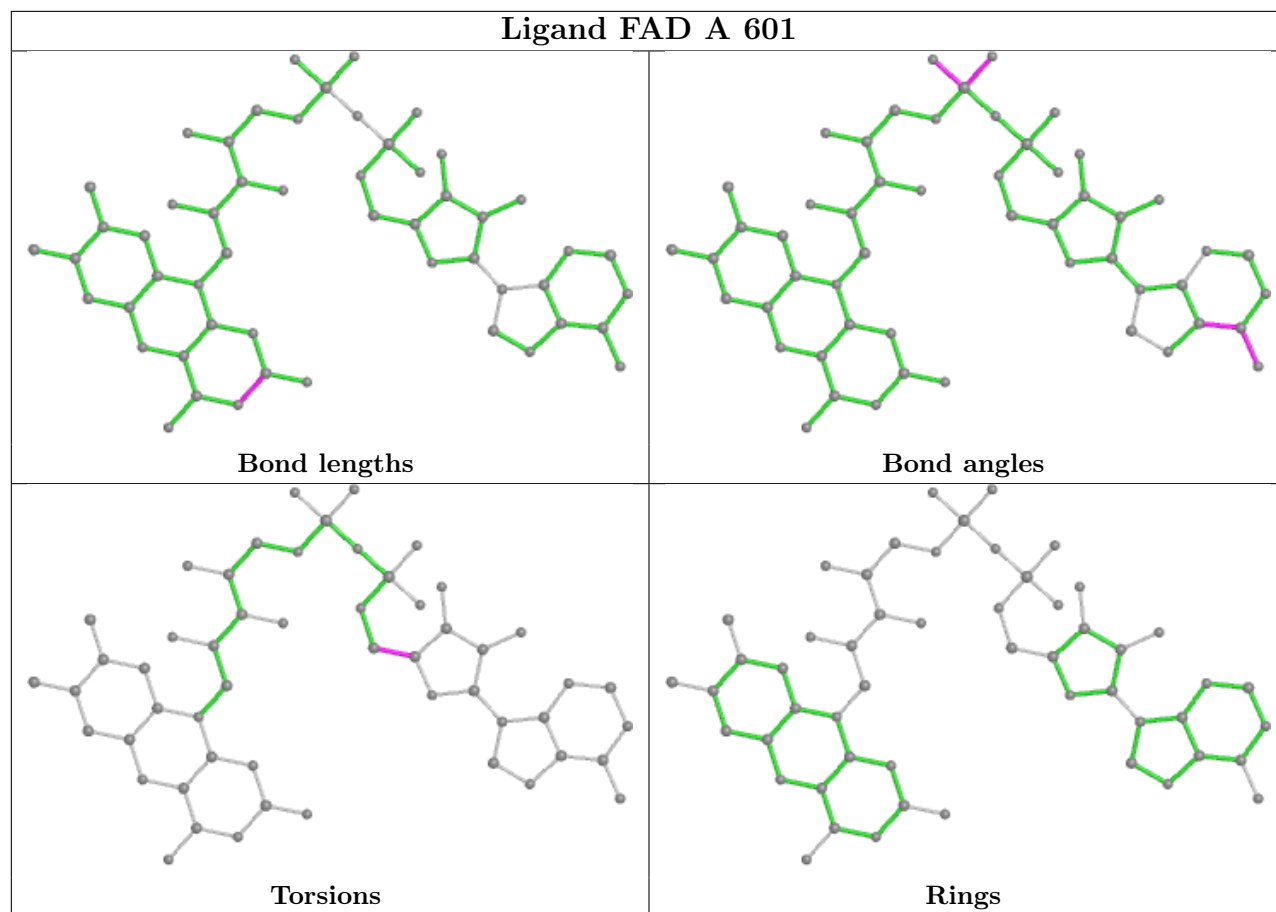
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

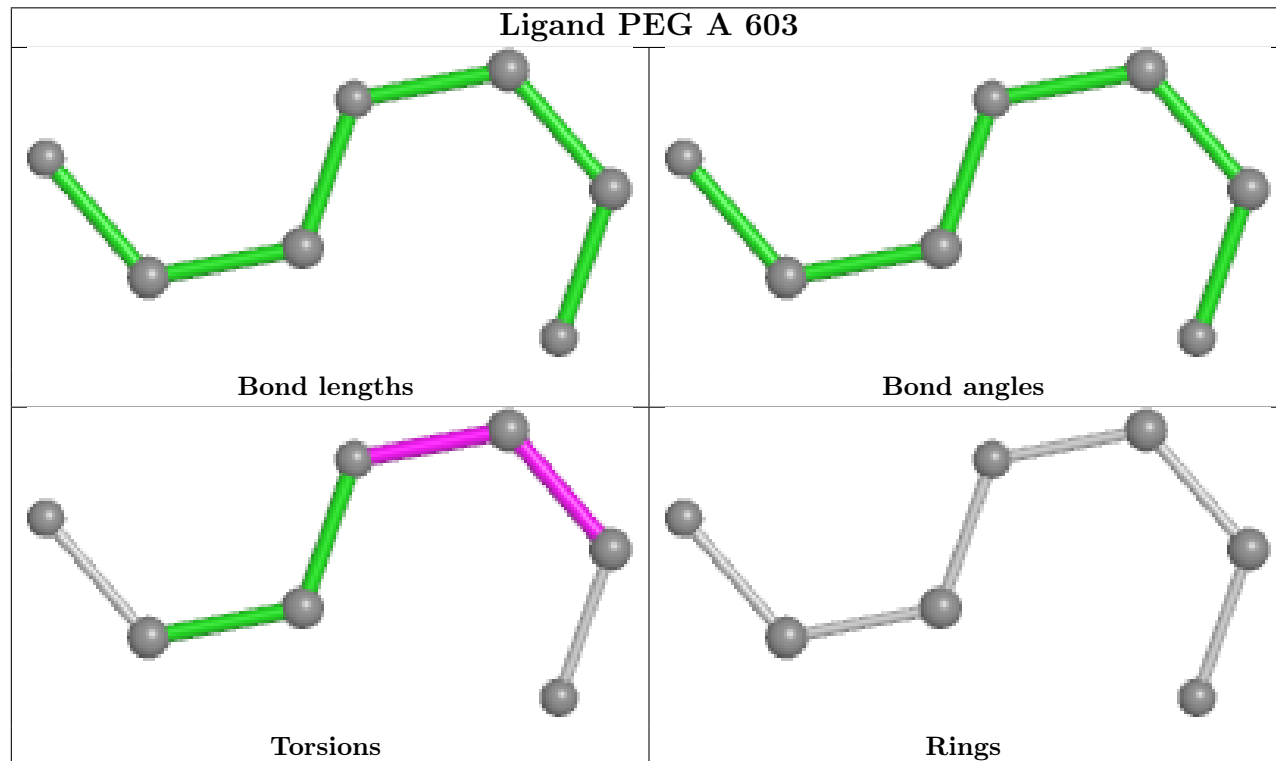
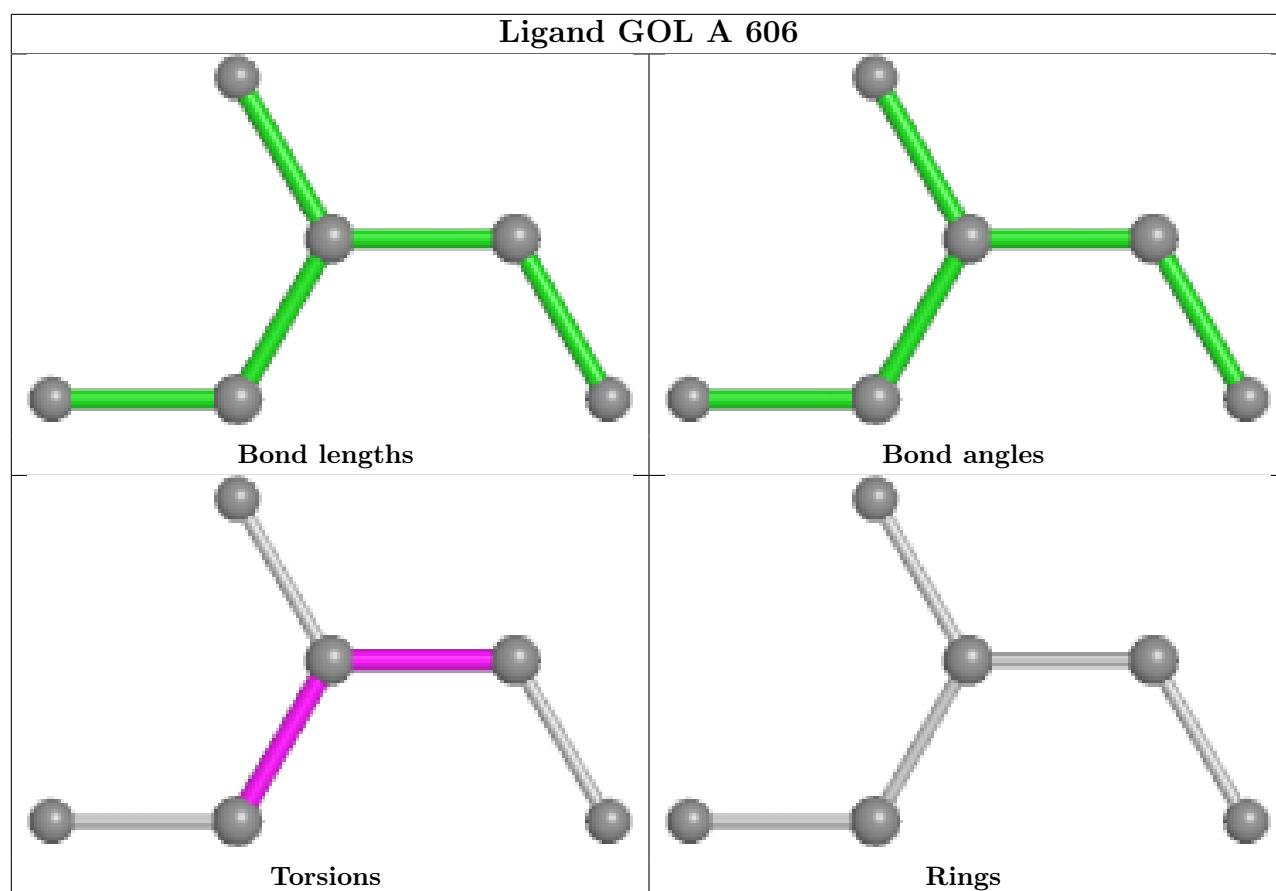


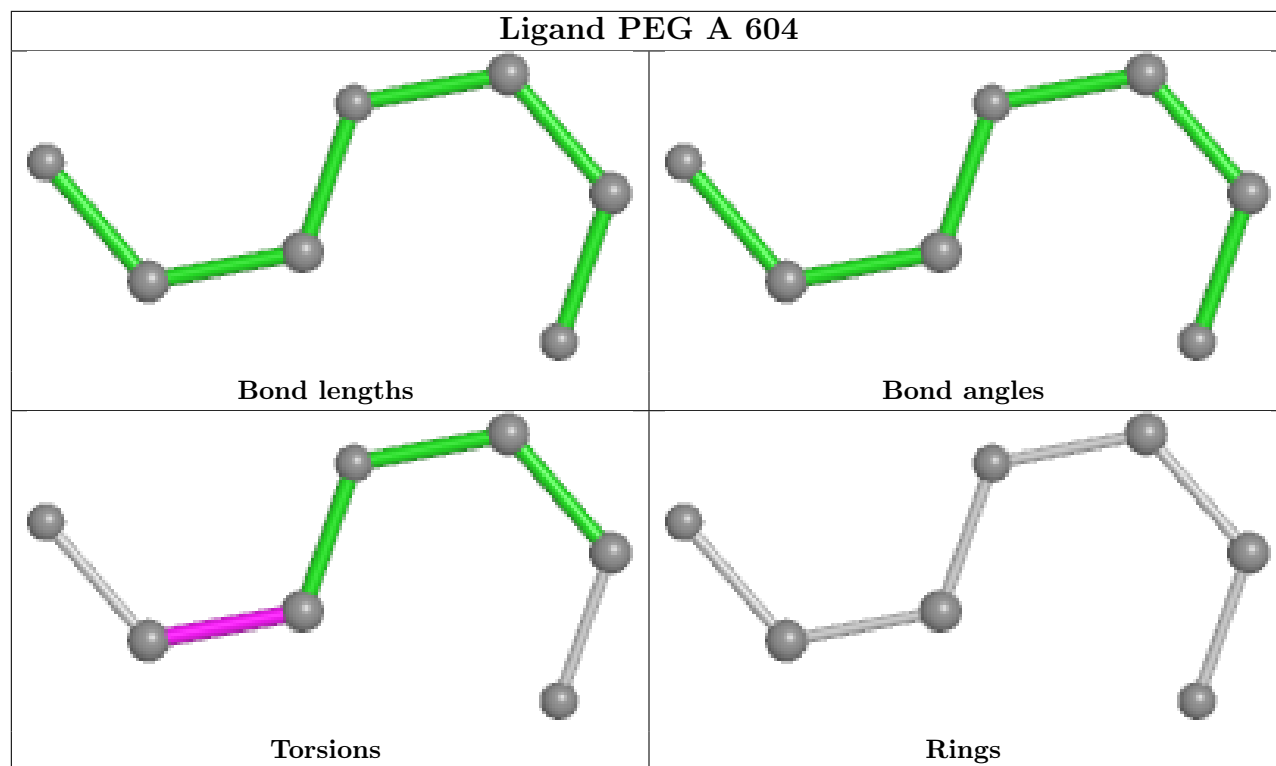












## 5.7 Other polymers [i](#)

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	558/561 (99%)	-0.27	10 (1%) 68 67	19, 27, 56, 109	0
1	B	557/561 (99%)	0.08	35 (6%) 20 17	22, 40, 74, 104	0
All	All	1115/1122 (99%)	-0.10	45 (4%) 38 35	19, 34, 68, 109	0

All (45) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	198	LEU	6.5
1	B	241	ALA	6.3
1	B	194	LEU	6.1
1	B	190	TYR	4.9
1	B	159	MET	4.7
1	A	194	LEU	4.4
1	A	198	LEU	4.3
1	B	201	ARG	3.8
1	B	174	ALA	3.7
1	B	179	TRP	3.5
1	A	4	GLU	3.4
1	B	411	ALA	3.3
1	B	195	PRO	3.2
1	B	196	TRP	3.1
1	A	192	LEU	3.1
1	B	192	LEU	3.1
1	B	561	LYS	3.1
1	B	191	TYR	3.0
1	A	5	THR	2.9
1	A	180	MET	2.9
1	B	184	ILE	2.9
1	A	179	TRP	2.9
1	B	189	ARG	2.9
1	B	428	LYS	2.7

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type	RSRZ
1	B	186	SER	2.7
1	B	200	SER	2.6
1	B	188	GLY	2.6
1	B	242	GLU	2.6
1	B	224	LYS	2.6
1	B	187	LEU	2.6
1	B	193	ASP	2.5
1	B	154	THR	2.5
1	B	197	ARG	2.5
1	B	5	THR	2.5
1	A	241	ALA	2.4
1	B	202	HIS	2.4
1	A	195	PRO	2.4
1	B	406	GLN	2.3
1	B	161	LYS	2.3
1	B	144	ASP	2.3
1	B	142	LEU	2.2
1	B	146	PHE	2.2
1	B	407	PRO	2.2
1	A	190	TYR	2.0
1	B	173	LEU	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	PEG	A	605	7/7	0.67	0.29	38,46,50,53	7
5	GOL	A	608	6/6	0.68	0.15	45,46,47,49	6

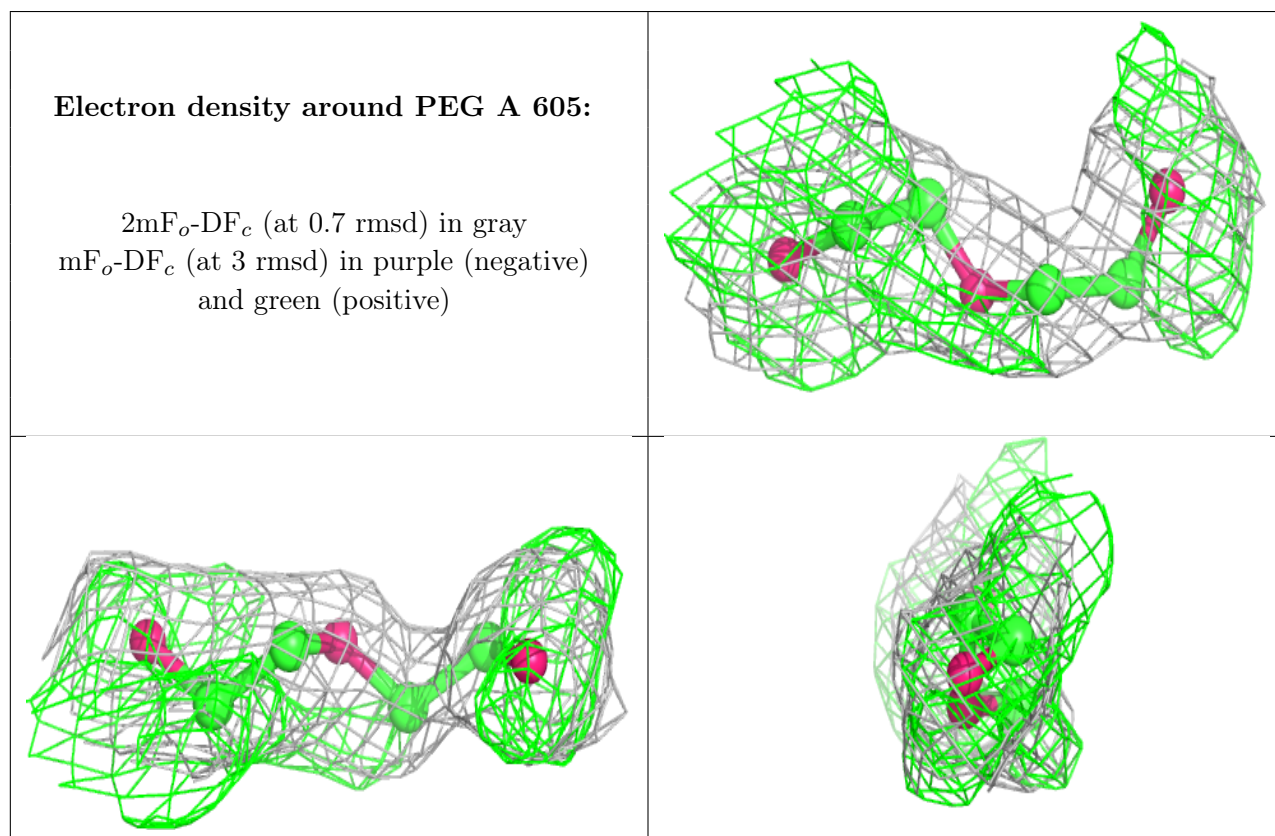
*Continued on next page...*



Continued from previous page...

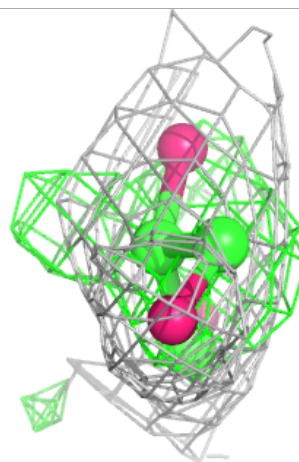
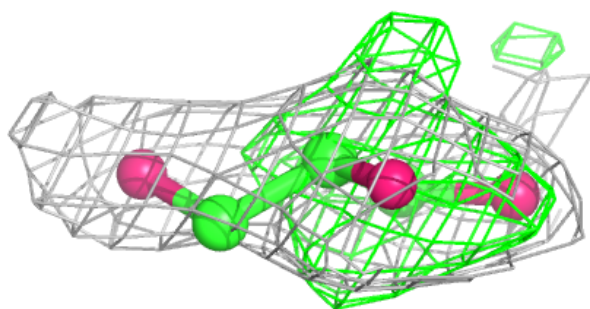
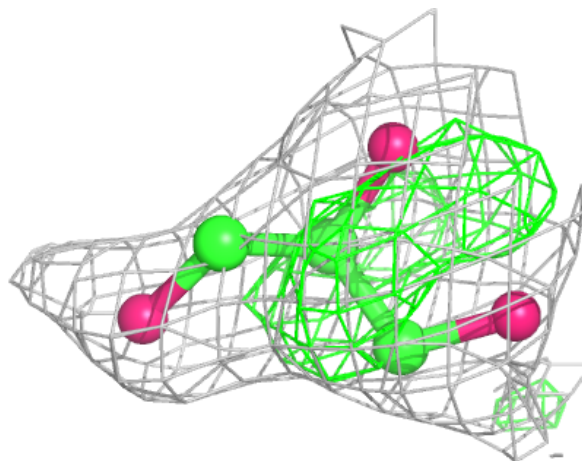
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
5	GOL	A	607	6/6	0.77	0.58	41,42,44,49	6
3	ANB	A	602	21/21	0.79	0.30	29,60,71,73	21
4	PEG	B	603	7/7	0.81	0.18	47,49,52,53	7
5	GOL	A	606	6/6	0.82	0.22	59,60,62,62	6
4	PEG	A	604	7/7	0.83	0.20	41,43,47,50	5
4	PEG	B	602	7/7	0.90	0.10	41,43,53,57	7
4	PEG	A	603	7/7	0.91	0.17	28,41,49,50	3
2	FAD	B	601	53/53	0.95	0.09	27,32,40,41	0
2	FAD	A	601	53/53	0.97	0.10	19,22,24,27	0
6	NA	A	609	1/1	0.98	0.13	28,28,28,28	0

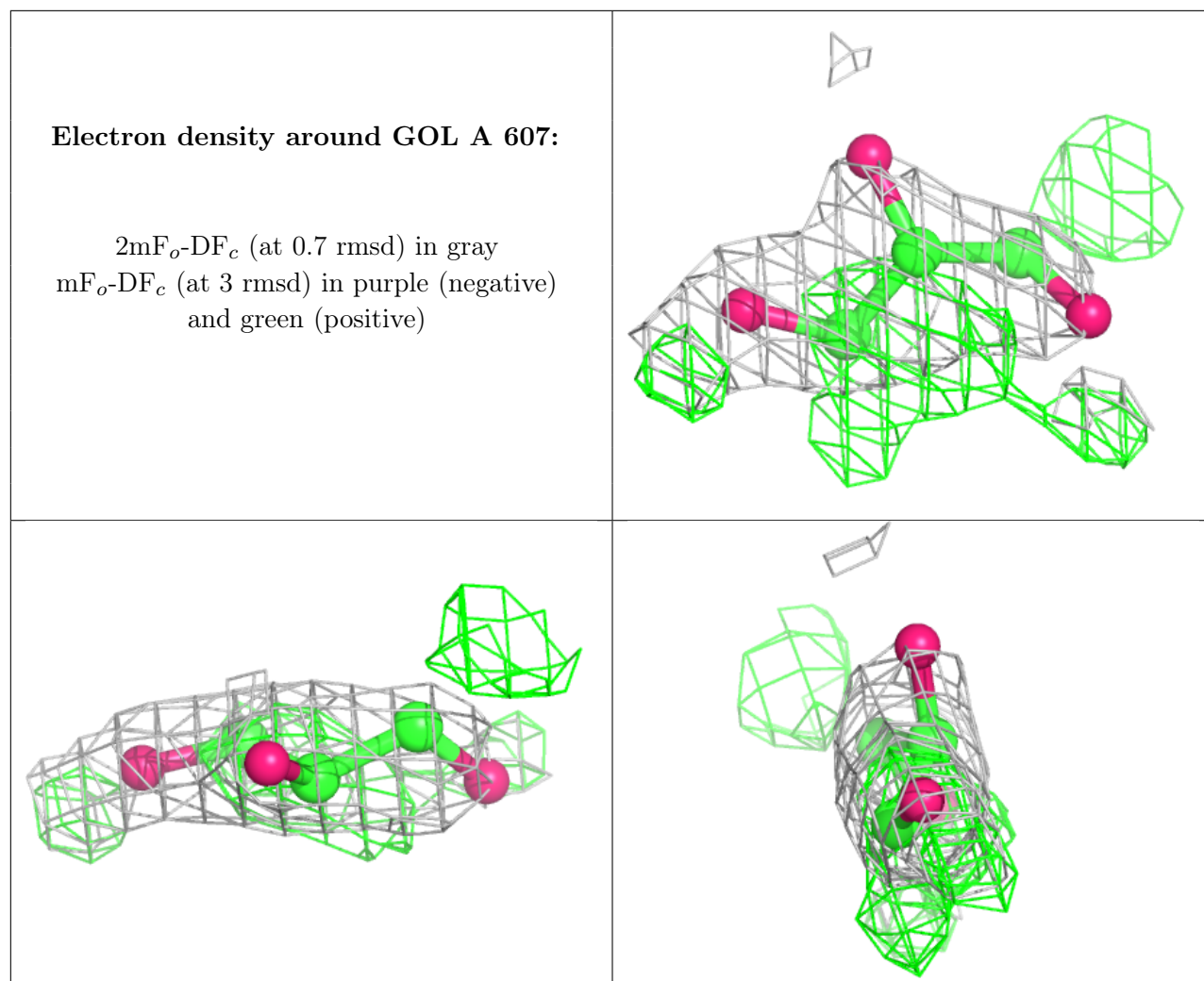
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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



**Electron density around GOL A 608:**

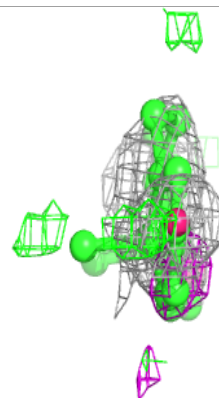
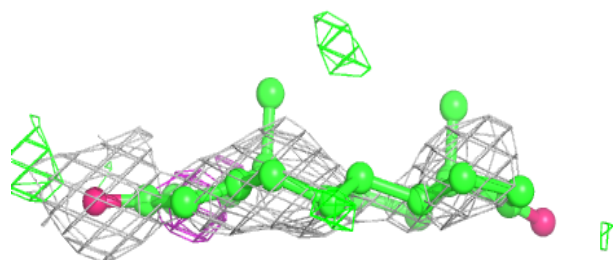
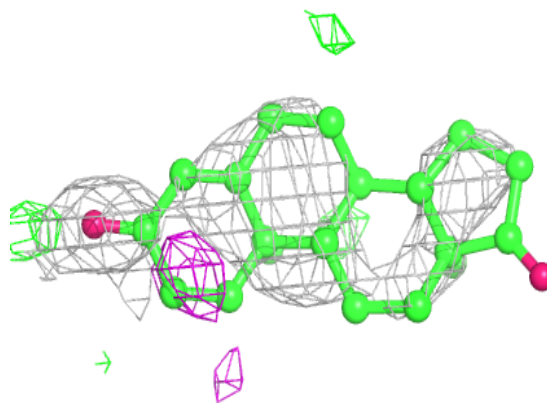
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)



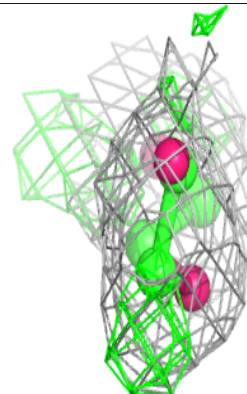
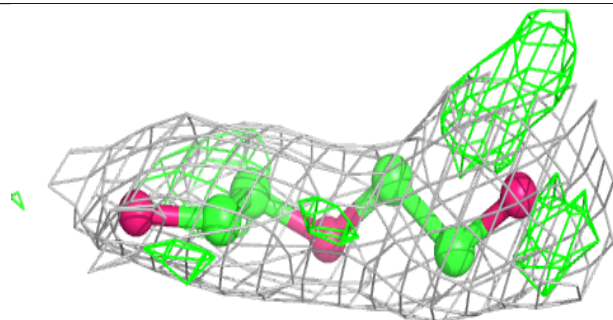
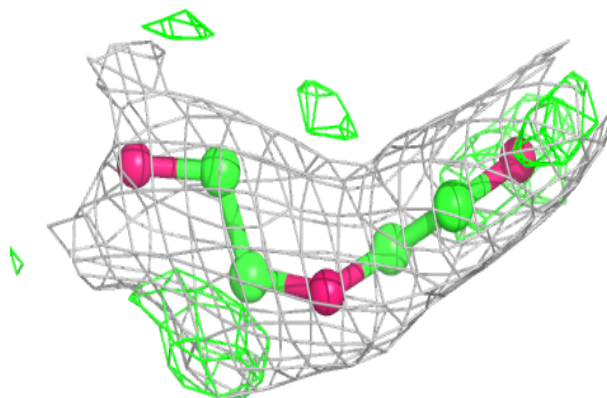


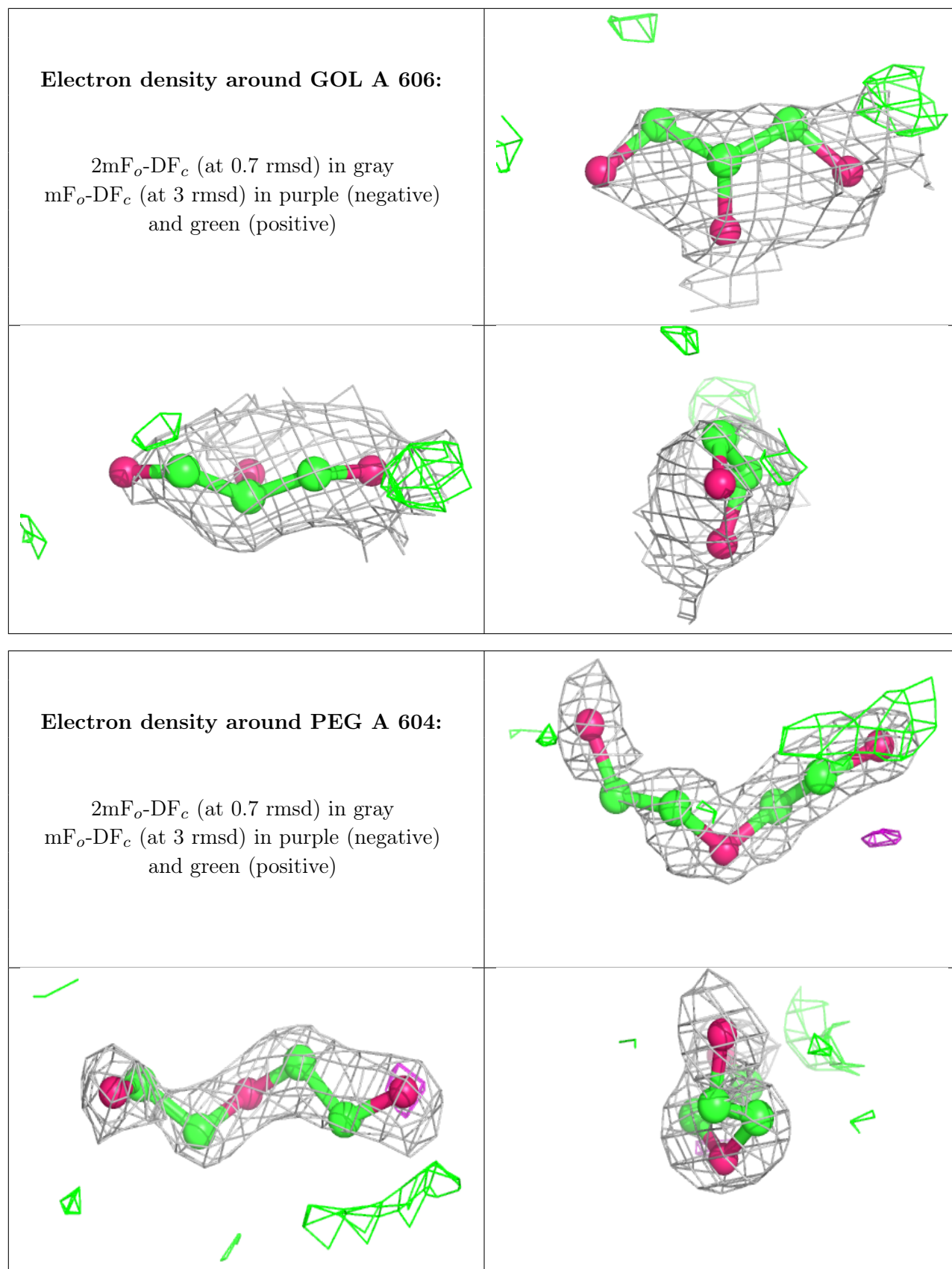
**Electron density around ANB A 602:**

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

**Electron density around PEG B 603:**

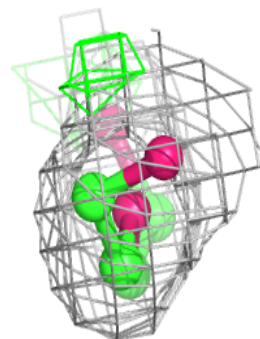
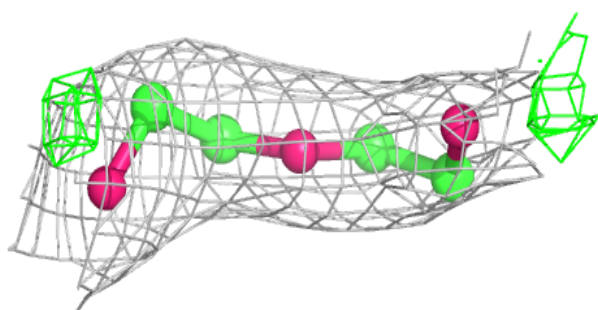
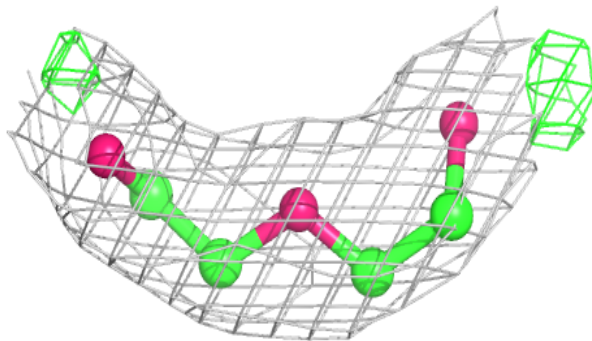
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)



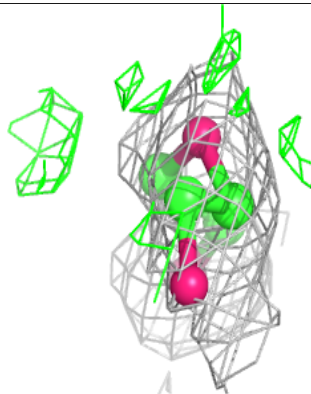
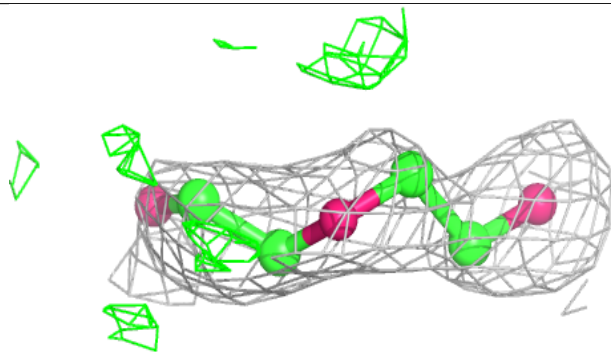
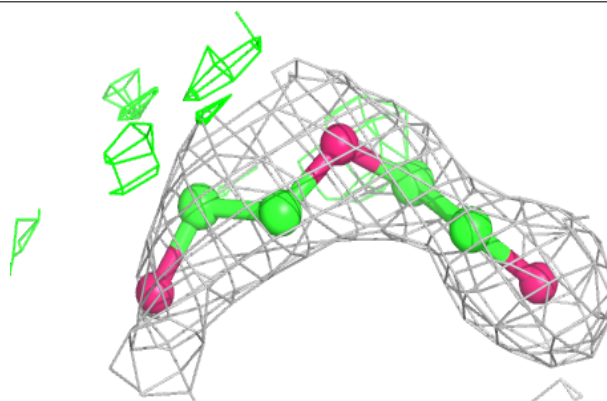


**Electron density around PEG B 602:**

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

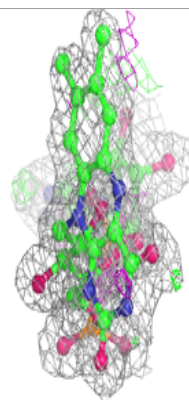
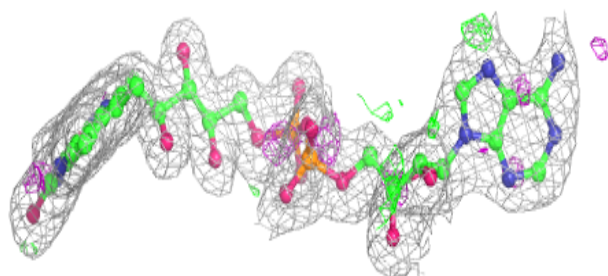
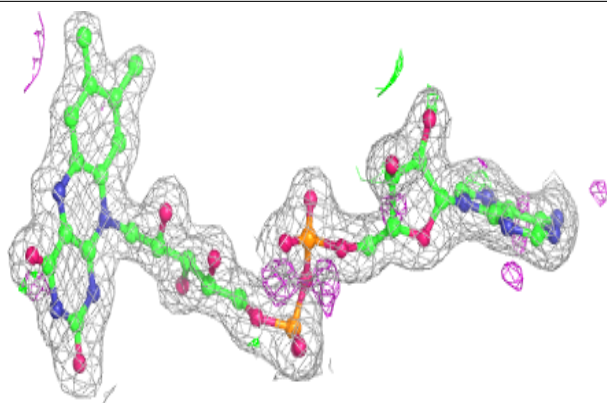
**Electron density around PEG A 603:**

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

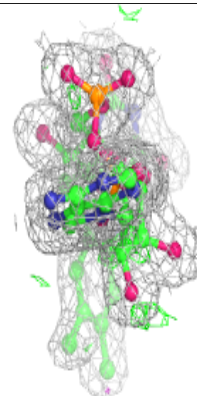
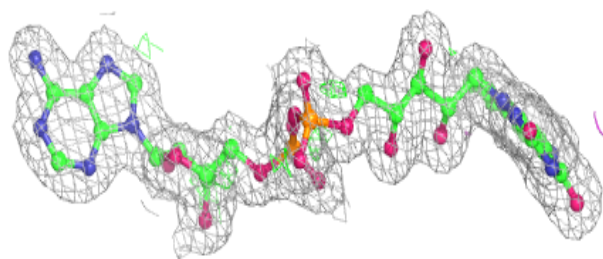
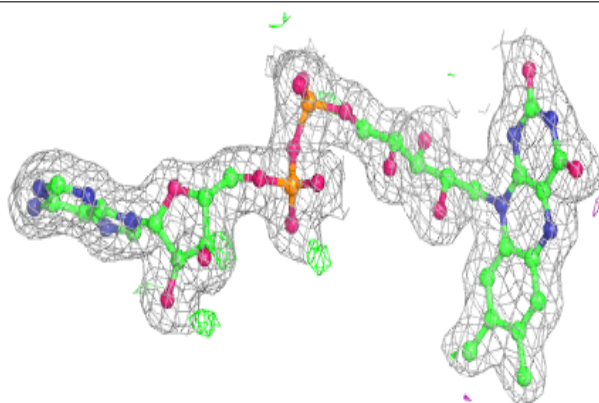


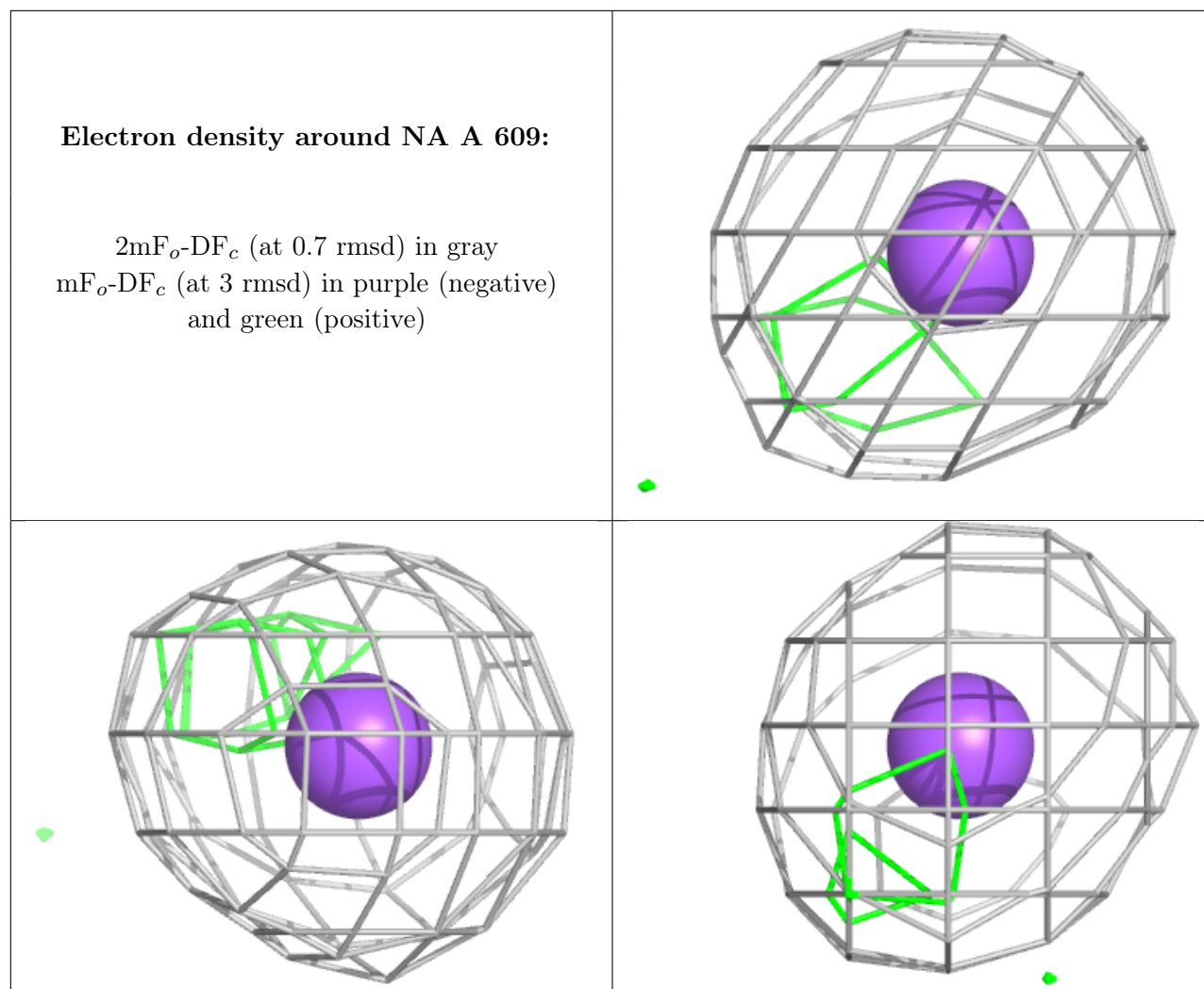
**Electron density around FAD B 601:**

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

**Electron density around FAD A 601:**

$2mF_o-DF_c$  (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.