



# Full wwPDB X-ray Structure Validation Report ⓘ

May 26, 2026 – 01:49 PM JST

PDB ID : 9V6F / pdb\_00009v6f  
Title : The crystal structure of a ThDP-dependent enzyme PpBFD  
Authors : Hou, X.L.; Zhou, J.H.  
Deposited on : 2025-05-27  
Resolution : 2.64 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 : 4-5-2 with Phenix2.0  
Mogul : 1.8.5 (274361), CSD as541be (2020)  
Xtriage (Phenix) : 2.0  
EDS : 3.0  
Buster-report : wwPDB partial adaption of 1.1.7 (2018)  
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)  
CCP4 : 9.0.010 (Gargrove)  
Density-Fitness : 1.0.12  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.49

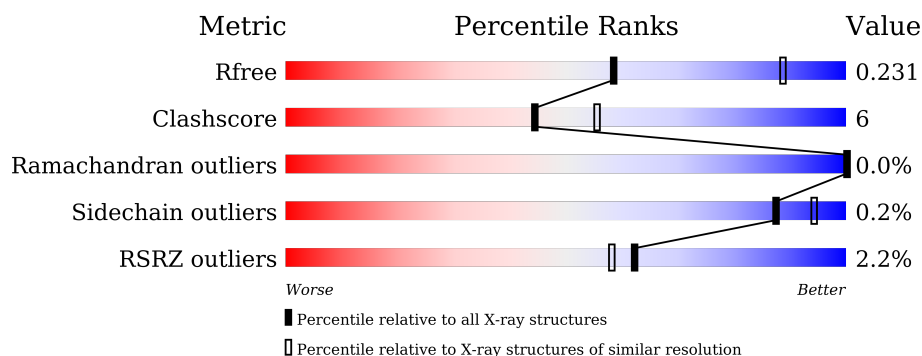
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.64 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))
$R_{free}$	180053	2053 (2.66-2.62)
Clashscore	190562	2097 (2.66-2.62)
Ramachandran outliers	187476	2066 (2.66-2.62)
Sidechain outliers	187428	2066 (2.66-2.62)
RSRZ outliers	180081	2052 (2.66-2.62)

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	530	<div> <div>3%</div> <div>88%</div> <div>11%</div> <div>.</div> </div>
1	B	530	<div> <div>2%</div> <div>79%</div> <div>15%</div> <div>5%</div> </div>
1	C	530	<div> <div>%</div> <div>84%</div> <div>11%</div> <div>5%</div> </div>
1	D	530	<div> <div>3%</div> <div>81%</div> <div>15%</div> <div>.</div> </div>

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

ria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	GOL	C	614	-	-	X	-

## 2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 15716 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 Benzoylformate decarboxylase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	524	Total	C	N	O	S	0	3	0
			3947	2497	683	748	19			
1	B	501	Total	C	N	O	S	0	3	0
			3774	2384	655	716	19			
1	C	503	Total	C	N	O	S	0	3	0
			3788	2394	657	718	19			
1	D	511	Total	C	N	O	S	0	3	0
			3844	2428	666	731	19			

There are 20 discrepancies between the modelled and reference sequences:

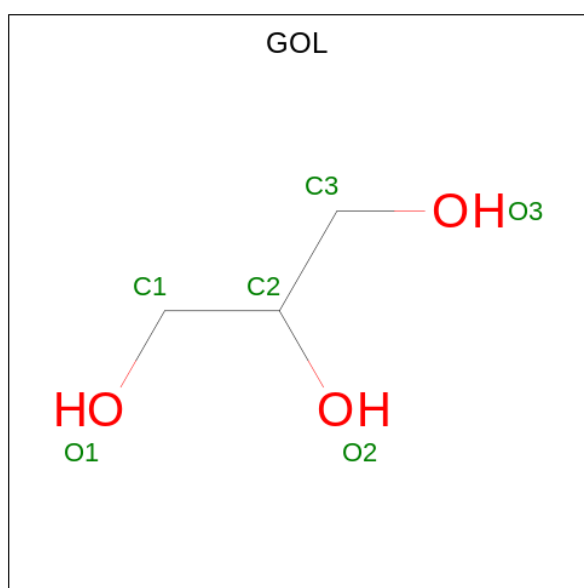
Chain	Residue	Modelled	Actual	Comment	Reference
A	109	PHE	LEU	conflict	UNP P20906
A	377	LEU	THR	conflict	UNP P20906
A	460	GLY	ALA	conflict	UNP P20906
A	529	LEU	-	expression tag	UNP P20906
A	530	GLU	-	expression tag	UNP P20906
B	109	PHE	LEU	conflict	UNP P20906
B	377	LEU	THR	conflict	UNP P20906
B	460	GLY	ALA	conflict	UNP P20906
B	529	LEU	-	expression tag	UNP P20906
B	530	GLU	-	expression tag	UNP P20906
C	109	PHE	LEU	conflict	UNP P20906
C	377	LEU	THR	conflict	UNP P20906
C	460	GLY	ALA	conflict	UNP P20906
C	529	LEU	-	expression tag	UNP P20906
C	530	GLU	-	expression tag	UNP P20906
D	109	PHE	LEU	conflict	UNP P20906
D	377	LEU	THR	conflict	UNP P20906
D	460	GLY	ALA	conflict	UNP P20906
D	529	LEU	-	expression tag	UNP P20906
D	530	GLU	-	expression tag	UNP P20906



- Molecule 2 is MAGNESIUM ION (CCD ID: MG) (formula: Mg) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	8	Total	Mg	0	0
			8	8		
2	B	11	Total	Mg	0	0
			11	11		
2	C	11	Total	Mg	0	0
			11	11		
2	D	2	Total	Mg	0	0
			2	2		

- Molecule 3 is GLYCEROL (CCD ID: 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
3	A	1	Total	C	O	0	0
			6	3	3		
3	A	1	Total	C	O	0	0
			6	3	3		
3	A	1	Total	C	O	0	0
			6	3	3		
3	A	1	Total	C	O	0	0
			6	3	3		
3	B	1	Total	C	O	0	0
			6	3	3		
3	B	1	Total	C	O	0	0
			6	3	3		

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

- Molecule 4 is CHLORIDE ION (CCD ID: CL) (formula: Cl) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	C	1	Total	Cl	0	0
			1	1		

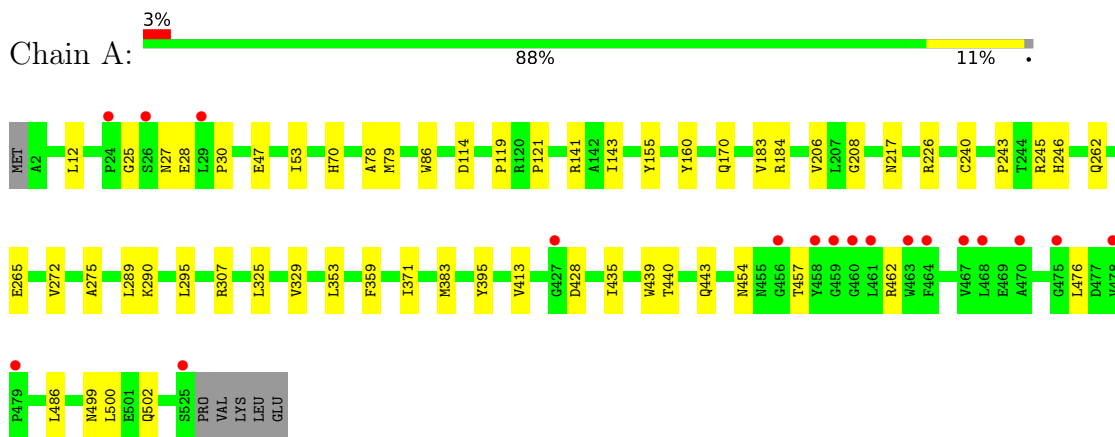
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	81	Total	O	0	0
			81	81		
5	B	39	Total	O	0	0
			39	39		
5	C	74	Total	O	0	0
			74	74		
5	D	52	Total	O	0	0
			52	52		

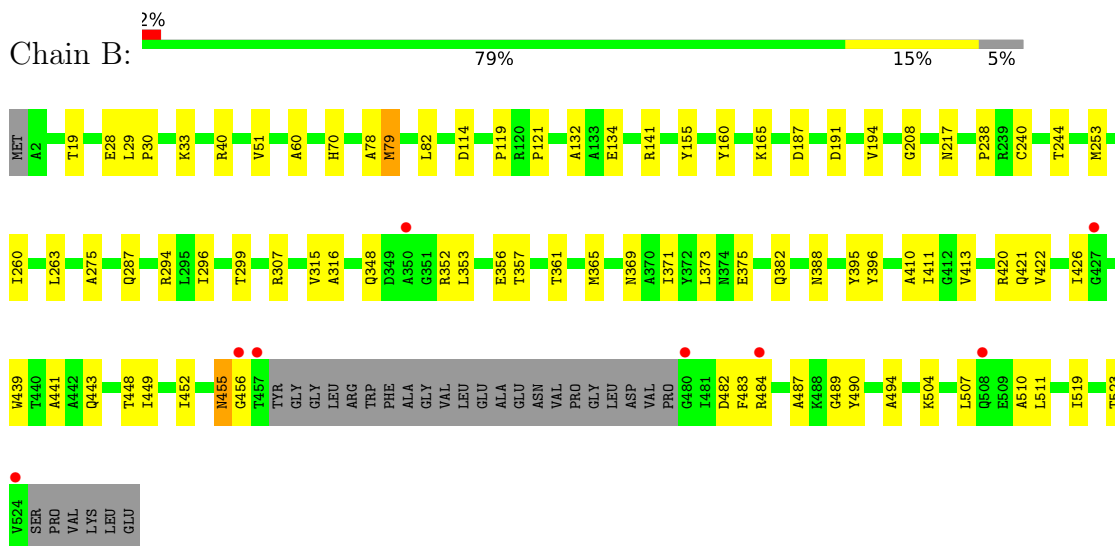
### 3 Residue-property plots

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.

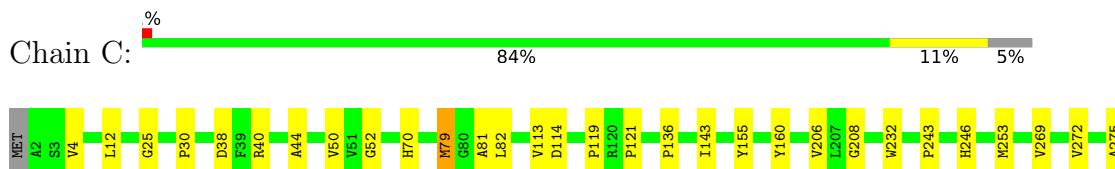
#### • Molecule 1: Benzoylformate decarboxylase

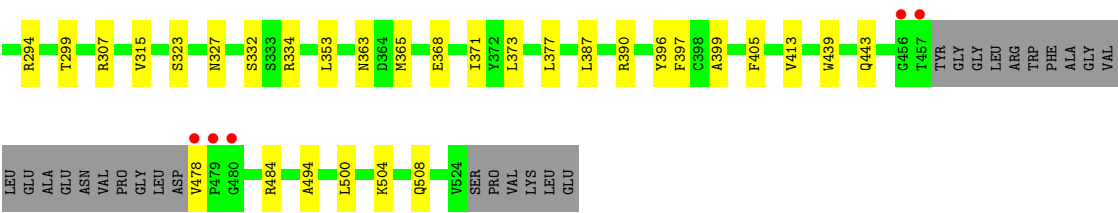


#### • Molecule 1: Benzoylformate decarboxylase

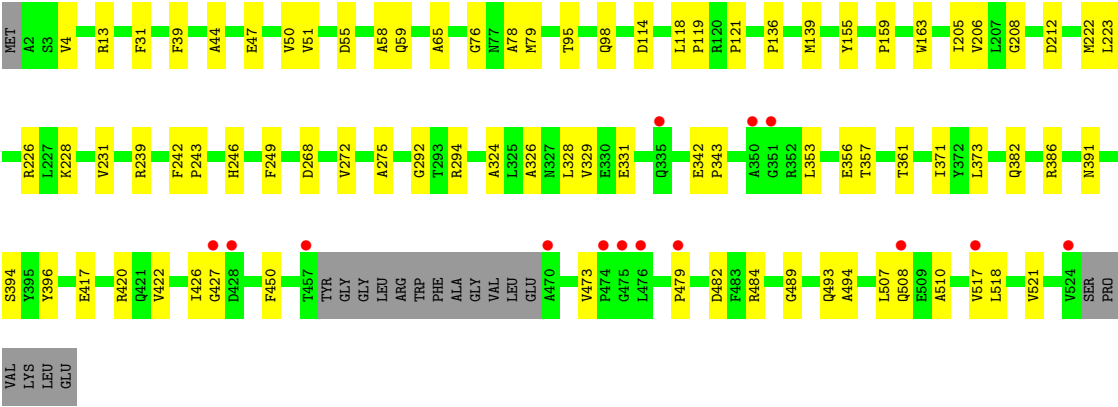
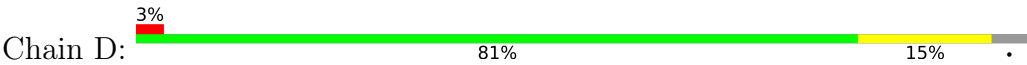


#### • Molecule 1: Benzoylformate decarboxylase





● Molecule 1: Benzoylformate decarboxylase



VAL  
LYS  
LEU  
GLU

## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	95.73Å 114.37Å 202.16Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.77 – 2.64 49.77 – 2.64	Depositor EDS
% Data completeness (in resolution range)	97.3 (49.77-2.64) 97.3 (49.77-2.64)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.18 (at 2.65Å)	Xtriage
Refinement program	PHENIX 1.12_2829, PHENIX 1.12_2829	Depositor
R, $R_{free}$	0.191 , 0.230 0.191 , 0.231	Depositor DCC
$R_{free}$ test set	3193 reflections (4.84%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	44.0	Xtriage
Anisotropy	0.117	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.32 , 45.3	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	15716	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	51.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.87% 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: GOL, MG, CL

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	0.18	0/4055	0.37	0/5542
1	B	0.17	0/3875	0.36	0/5293
1	C	0.18	0/3890	0.37	0/5315
1	D	0.16	0/3947	0.35	0/5394
All	All	0.17	0/15767	0.36	0/21544

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	3947	0	3871	41	0
1	B	3774	0	3706	58	0
1	C	3788	0	3722	43	0
1	D	3844	0	3773	54	0
2	A	8	0	0	0	0
2	B	11	0	0	0	0
2	C	11	0	0	0	0
2	D	2	0	0	0	0
3	A	24	0	32	3	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	18	0	24	3	0
3	C	42	0	56	7	0
4	C	1	0	0	1	0
5	A	81	0	0	0	0
5	B	39	0	0	0	0
5	C	74	0	0	1	0
5	D	52	0	0	1	0
All	All	15716	0	15184	184	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 6.

All (184) 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:79:MET:HE1	1:A:119:PRO:HA	1.56	0.86
1:C:81:ALA:HA	3:C:614:GOL:H2	1.59	0.84
1:C:405:PHE:HA	3:C:614:GOL:H12	1.57	0.84
1:B:60:ALA:HB2	1:B:413:VAL:HG23	1.65	0.77
1:D:76:GLY:HA2	1:D:79:MET:HE3	1.65	0.76
1:C:52:GLY:HA2	3:C:614:GOL:H31	1.67	0.73
1:C:206:VAL:HG12	1:C:232:TRP:HB2	1.70	0.73
1:B:352:ARG:NH1	1:B:523:THR:O	2.23	0.72
1:C:478:VAL:N	3:C:618:GOL:HO2	1.88	0.71
1:B:361:THR:HG23	1:B:365:MET:HE3	1.73	0.71
3:A:610:GOL:H11	1:C:113:VAL:HB	1.72	0.70
1:B:253:MET:HE3	1:B:263:LEU:HD11	1.75	0.69
1:B:411:ILE:HG12	1:B:441:ALA:HB2	1.75	0.68
1:C:307:ARG:HH12	3:C:617:GOL:H32	1.59	0.67
1:D:222:MET:SD	1:D:226:ARG:NH2	2.66	0.67
1:B:121:PRO:HG3	1:D:114:ASP:HB3	1.78	0.66
1:B:29:LEU:HD22	1:D:473:VAL:HG11	1.78	0.66
1:B:482:ASP:OD1	1:B:484:ARG:HG2	1.96	0.66
1:C:323:SER:O	1:C:327:ASN:ND2	2.28	0.65
1:D:508:GLN:NE2	1:D:508:GLN:O	2.30	0.65
1:A:27:ASN:H	1:A:70:HIS:CE1	2.17	0.62
1:C:368:GLU:HG2	1:C:390:ARG:HE	1.64	0.62
1:B:30:PRO:HA	1:B:33:LYS:HD3	1.80	0.62
1:A:114:ASP:HB3	1:C:121:PRO:HG3	1.81	0.61
1:B:208:GLY:HA3	1:B:275:ALA:HB2	1.82	0.61
1:D:373:LEU:HG	1:D:396:TYR:HB2	1.82	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:228:LYS:NZ	1:D:331:GLU:OE2	2.31	0.61
1:C:208:GLY:HA3	1:C:275:ALA:HB2	1.83	0.60
1:D:226:ARG:HG2	1:D:329:VAL:HG23	1.83	0.60
1:B:114:ASP:HB3	1:D:121:PRO:HG3	1.84	0.60
1:D:59:GLN:HB3	1:D:239:ARG:HH12	1.66	0.60
1:C:206:VAL:HG23	1:C:272:VAL:HG13	1.84	0.60
1:C:399:ALA:N	4:C:612:CL:CL	2.70	0.58
1:D:47:GLU:HB3	1:D:78:ALA:HB2	1.85	0.58
1:B:134:GLU:HG3	3:C:613:GOL:H32	1.85	0.57
1:D:493:GLN:HB3	1:D:517:VAL:HG23	1.85	0.57
1:B:375:GLU:HB3	1:B:426:ILE:HG23	1.87	0.56
1:C:377:LEU:HD22	1:C:397:PHE:HE1	1.71	0.56
1:A:208:GLY:HA3	1:A:275:ALA:HB2	1.86	0.56
1:B:132:ALA:HB2	3:B:612:GOL:H12	1.86	0.56
1:C:371:ILE:HG21	1:C:413:VAL:HG11	1.87	0.56
1:C:79:MET:HE1	1:C:119:PRO:HA	1.86	0.56
1:A:226:ARG:NH1	1:A:329:VAL:O	2.40	0.55
1:D:342:GLU:HG3	1:D:343:PRO:HD2	1.89	0.55
1:A:86:TRP:CD1	3:A:610:GOL:HO2	2.24	0.54
1:B:253:MET:HE2	1:B:260:ILE:HG12	1.89	0.54
1:D:231:VAL:HG21	1:D:242:PHE:CE2	2.42	0.53
1:B:455:ASN:OD1	1:B:456:GLY:N	2.42	0.53
1:D:208:GLY:HA3	1:D:275:ALA:HB2	1.91	0.53
1:B:353:LEU:H	1:B:353:LEU:HD12	1.75	0.52
1:B:373:LEU:HD11	1:B:413:VAL:HG11	1.92	0.52
1:C:500:LEU:HG	1:C:504:LYS:HE2	1.92	0.52
1:C:206:VAL:HA	1:C:232:TRP:O	2.10	0.52
1:B:507:LEU:HD12	1:B:519:ILE:HD13	1.92	0.52
1:C:368:GLU:HG2	1:C:390:ARG:NE	2.25	0.51
1:B:487:ALA:HB3	1:B:494:ALA:HB2	1.91	0.51
1:D:356:GLU:HG3	1:D:382:GLN:HE21	1.74	0.51
1:D:357:THR:O	1:D:361:THR:OG1	2.23	0.51
1:B:371:ILE:HB	1:B:422:VAL:HG22	1.92	0.51
1:D:231:VAL:HG23	1:D:249:PHE:HA	1.92	0.51
1:B:413:VAL:HG13	1:B:422:VAL:HG21	1.92	0.51
1:C:269:VAL:HG22	1:C:294:ARG:HD3	1.91	0.51
1:B:79:MET:HE3	1:B:82:LEU:HD12	1.92	0.50
1:B:132:ALA:H	3:B:612:GOL:H31	1.76	0.50
1:B:165:LYS:HG3	3:B:612:GOL:H11	1.91	0.50
1:D:95:THR:HG21	1:D:139:MET:HE1	1.93	0.50
1:A:79:MET:CE	1:A:119:PRO:HA	2.36	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:119:PRO:HG3	1:A:155:TYR:CD2	2.47	0.50
1:A:243:PRO:HG2	1:A:246:HIS:HB2	1.94	0.49
1:D:212:ASP:HB3	1:D:239:ARG:HE	1.77	0.49
1:C:484:ARG:HG2	1:C:494:ALA:HB1	1.93	0.49
1:D:119:PRO:HG3	1:D:155:TYR:CG	2.47	0.49
1:D:371:ILE:HB	1:D:422:VAL:HG22	1.93	0.49
1:B:373:LEU:HD23	1:B:396:TYR:HB2	1.94	0.49
1:B:410:ALA:HB1	1:B:448:THR:HG21	1.94	0.49
1:D:205:ILE:HB	1:D:231:VAL:HG12	1.95	0.49
1:D:391:ASN:O	1:D:394:SER:OG	2.27	0.49
1:D:450:PHE:HB2	1:D:518:LEU:HD12	1.94	0.49
1:D:507:LEU:O	1:D:510:ALA:N	2.46	0.48
1:A:121:PRO:HG3	1:C:114:ASP:HB2	1.94	0.48
1:C:439:TRP:CZ2	1:C:443:GLN:HG3	2.49	0.48
1:D:4:VAL:HG13	1:D:136:PRO:HD3	1.95	0.48
1:A:184:ARG:NH2	1:B:187:ASP:OD1	2.45	0.48
1:A:265:GLU:HB2	1:A:290:LYS:HD2	1.95	0.48
1:A:457:THR:OG1	1:A:476:LEU:O	2.22	0.48
1:D:382:GLN:HE22	1:D:386:ARG:HH21	1.61	0.48
1:A:428:ASP:HB3	1:A:454:ASN:HA	1.96	0.47
1:B:452:ILE:HG21	1:B:483:PHE:CD1	2.49	0.47
1:A:462:ARG:HG2	1:A:476:LEU:HD13	1.95	0.47
1:B:449:ILE:HG12	1:B:510:ALA:HB1	1.96	0.47
1:A:353:LEU:HD21	1:A:500:LEU:HA	1.96	0.47
1:B:28:GLU:OE1	1:B:70:HIS:HA	2.15	0.47
1:D:51:VAL:HG21	1:D:78:ALA:HB1	1.96	0.47
1:B:294:ARG:CZ	1:B:296:ILE:HD11	2.44	0.47
1:B:482:ASP:HB3	1:D:489:GLY:HA2	1.96	0.47
1:D:417:GLU:HB3	1:D:420:ARG:HG3	1.96	0.47
1:A:119:PRO:HG3	1:A:155:TYR:CG	2.49	0.47
1:C:439:TRP:CE2	1:C:443:GLN:HG3	2.50	0.47
1:C:12:LEU:HD23	1:C:143:ILE:HD11	1.96	0.47
1:C:332:SER:OG	1:C:334:ARG:NE	2.47	0.46
1:D:356:GLU:HG3	1:D:382:GLN:NE2	2.30	0.46
1:C:119:PRO:HG3	1:C:155:TYR:CD2	2.50	0.46
1:D:58:ALA:HB2	1:D:65:ALA:HB2	1.98	0.46
1:A:499:ASN:OD1	1:A:502:GLN:HG3	2.16	0.46
1:C:377:LEU:HD22	1:C:397:PHE:CE1	2.51	0.46
1:A:325:LEU:O	1:A:329:VAL:HG22	2.14	0.46
1:A:25:GLY:HA3	1:A:70:HIS:CD2	2.51	0.46
1:B:369:ASN:O	1:B:420:ARG:NE	2.45	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:299:THR:O	1:C:315:VAL:HA	2.16	0.46
1:A:27:ASN:H	1:A:70:HIS:HE1	1.61	0.45
1:D:79:MET:HE1	1:D:118:LEU:HG	1.98	0.45
1:D:294:ARG:NH2	5:D:701:HOH:O	2.38	0.45
1:A:53:ILE:HD13	1:A:440:THR:HG21	1.97	0.45
1:B:191:ASP:HA	1:B:194:VAL:HG12	1.98	0.45
1:B:439:TRP:HB2	1:B:490:TYR:CD1	2.51	0.45
1:C:79:MET:HE3	1:C:82:LEU:HD12	1.99	0.45
1:D:353:LEU:HD13	1:D:521:VAL:HG13	1.99	0.45
1:A:371:ILE:HG21	1:A:413:VAL:HG11	1.99	0.44
1:B:365:MET:HE1	1:B:504:LYS:HA	1.98	0.44
1:D:223:LEU:HD12	1:D:329:VAL:HG21	1.99	0.44
1:A:439:TRP:CZ2	1:A:443:GLN:HG3	2.53	0.44
1:A:245:ARG:HA	1:A:245:ARG:HD3	1.68	0.44
1:C:405:PHE:HD1	3:C:614:GOL:H32	1.81	0.44
1:B:299:THR:O	1:B:315:VAL:HA	2.18	0.44
1:B:238:PRO:O	1:B:396:TYR:HA	2.17	0.44
1:D:13:ARG:HG3	1:D:39:PHE:HE1	1.83	0.44
1:B:30:PRO:HG2	1:B:160:TYR:HB2	1.99	0.43
1:D:31:PHE:HB2	1:D:163:TRP:CE2	2.53	0.43
1:B:348:GLN:HG2	1:B:357:THR:HG21	2.00	0.43
1:D:324:ALA:O	1:D:328:LEU:HG	2.18	0.43
1:B:439:TRP:CZ2	1:B:443:GLN:HG3	2.53	0.43
1:C:119:PRO:HG3	1:C:155:TYR:CG	2.53	0.43
1:C:232:TRP:HB3	1:C:253:MET:HG3	2.00	0.43
1:A:12:LEU:HD23	1:A:143:ILE:HD11	2.01	0.43
1:C:44:ALA:HB3	1:C:50:VAL:HG22	2.01	0.43
1:D:222:MET:HG2	1:D:326:ALA:HB1	2.01	0.43
1:A:25:GLY:HA3	1:A:70:HIS:CG	2.53	0.43
1:C:365:MET:HE3	1:C:508:GLN:OE1	2.19	0.43
1:A:30:PRO:HG2	1:A:160:TYR:HD1	1.83	0.43
1:C:25:GLY:HA3	1:C:70:HIS:CG	2.54	0.43
1:B:244:THR:OG1	1:B:388:ASN:HA	2.19	0.42
1:C:243:PRO:HG2	1:C:246:HIS:HB2	2.01	0.42
1:D:484:ARG:HG2	1:D:494:ALA:HB1	2.01	0.42
1:B:217:ASN:OD1	1:B:217:ASN:N	2.52	0.42
1:C:30:PRO:HG2	1:C:160:TYR:HB2	2.02	0.42
1:B:119:PRO:HG3	1:B:155:TYR:CG	2.54	0.42
1:B:443:GLN:NE2	1:D:479:PRO:O	2.52	0.42
1:A:183:VAL:HG11	1:B:316:ALA:HB2	2.02	0.42
1:A:262:GLN:OE1	1:A:290:LYS:HE3	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:435:ILE:HG12	1:A:486:LEU:HD13	2.01	0.42
1:D:226:ARG:HD3	1:D:331:GLU:OE1	2.20	0.42
1:C:363:ASN:HB2	1:C:387:LEU:HD23	2.01	0.42
1:C:353:LEU:HD21	1:C:500:LEU:HA	2.01	0.42
1:A:217:ASN:OD1	1:A:217:ASN:N	2.53	0.42
1:A:307:ARG:HA	1:B:141:ARG:HG2	2.02	0.41
1:B:240:CYS:HB2	1:B:395:TYR:HB3	2.02	0.41
1:D:55:ASP:O	1:D:59:GLN:HG3	2.21	0.41
1:B:489:GLY:HA2	1:D:482:ASP:HB3	2.01	0.41
1:D:243:PRO:HG2	1:D:246:HIS:HB2	2.02	0.41
1:A:141:ARG:HG2	1:B:307:ARG:HA	2.02	0.41
1:A:28:GLU:OE2	1:A:70:HIS:ND1	2.52	0.41
1:B:421:GLN:HG2	1:B:511:LEU:HD23	2.02	0.41
1:D:268:ASP:HA	1:D:292:GLY:O	2.21	0.41
1:D:426:ILE:HG22	1:D:427:GLY:O	2.21	0.41
1:D:98:GLN:HB3	1:D:159:PRO:HA	2.03	0.41
1:A:170:GLN:CD	1:B:287:GLN:HA	2.45	0.41
1:A:359:PHE:CE1	1:A:383:MET:HG3	2.56	0.41
1:D:59:GLN:HB3	1:D:239:ARG:NH1	2.32	0.41
1:A:47:GLU:HB3	1:A:78:ALA:HB2	2.03	0.41
1:B:28:GLU:CD	1:B:70:HIS:HA	2.46	0.41
1:B:79:MET:HE1	1:B:119:PRO:HA	2.03	0.40
1:B:356:GLU:HG3	1:B:382:GLN:HG2	2.03	0.40
1:D:44:ALA:HB3	1:D:50:VAL:HG22	2.03	0.40
3:A:611:GOL:H31	5:C:754:HOH:O	2.22	0.40
1:B:51:VAL:HG21	1:B:78:ALA:HB1	2.03	0.40
1:C:4:VAL:HG13	1:C:136:PRO:HD3	2.03	0.40
1:C:38:ASP:OD1	1:C:38:ASP:N	2.45	0.40
1:A:206:VAL:HB	1:A:272:VAL:HG22	2.04	0.40
1:A:240:CYS:HB2	1:A:395:TYR:HB3	2.03	0.40
1:B:19:THR:HG23	1:B:40:ARG:HG3	2.03	0.40
1:C:373:LEU:HG	1:C:396:TYR:HB2	2.04	0.40
1:A:289:LEU:HD11	1:A:295:LEU:HB2	2.03	0.40
1:D:206:VAL:O	1:D:272:VAL:HA	2.22	0.40
1:D:356:GLU:OE1	1:D:356:GLU:N	2.40	0.40

There are no symmetry-related clashes.

## 5.3 Torsion angles

### 5.3.1 Protein backbone

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	525/530 (99%)	515 (98%)	10 (2%)	0	100	100
1	B	500/530 (94%)	483 (97%)	16 (3%)	1 (0%)	43	58
1	C	502/530 (95%)	492 (98%)	10 (2%)	0	100	100
1	D	510/530 (96%)	501 (98%)	9 (2%)	0	100	100
All	All	2037/2120 (96%)	1991 (98%)	45 (2%)	1 (0%)	100	100

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	455	ASN

### 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	413/416 (99%)	413 (100%)	0	100	100
1	B	396/416 (95%)	395 (100%)	1 (0%)	86	93
1	C	398/416 (96%)	396 (100%)	2 (0%)	81	89
1	D	404/416 (97%)	404 (100%)	0	100	100
All	All	1611/1664 (97%)	1608 (100%)	3 (0%)	87	94

All (3) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	79	MET
1	C	40	ARG
1	C	79	MET

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (19) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	23	ASN
1	A	59	GLN
1	A	432	ASN
1	A	508	GLN
1	B	23	ASN
1	B	46	GLN
1	B	87	ASN
1	B	178	HIS
1	B	363	ASN
1	B	382	GLN
1	B	443	GLN
1	C	363	ASN
1	C	385	GLN
1	C	388	ASN
1	D	186	ASN
1	D	281	HIS
1	D	382	GLN
1	D	385	GLN
1	D	432	ASN

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

## 5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

## 5.6 Ligand geometry

Of 47 ligands modelled in this entry, 33 are monoatomic - leaving 14 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	$\# Z  > 2$
3	GOL	A	609	-	5,5,5	0.91	0	5,5,5	0.99	0
3	GOL	B	612	-	5,5,5	0.98	0	5,5,5	1.02	0
3	GOL	B	614	-	5,5,5	0.79	0	5,5,5	1.07	0
3	GOL	C	616	-	5,5,5	1.00	0	5,5,5	0.89	0
3	GOL	C	615	-	5,5,5	0.89	0	5,5,5	0.99	0
3	GOL	C	618	-	5,5,5	0.88	0	5,5,5	1.01	0
3	GOL	C	619	-	5,5,5	0.87	0	5,5,5	1.03	0
3	GOL	C	613	-	5,5,5	1.03	0	5,5,5	0.85	0
3	GOL	A	610	-	5,5,5	1.00	0	5,5,5	0.99	0
3	GOL	B	613	-	5,5,5	0.95	0	5,5,5	0.98	0
3	GOL	A	612	-	5,5,5	1.14	0	5,5,5	0.81	0
3	GOL	C	617	-	5,5,5	0.78	0	5,5,5	0.99	0
3	GOL	C	614	-	5,5,5	1.11	0	5,5,5	0.92	0
3	GOL	A	611	-	5,5,5	0.90	0	5,5,5	1.07	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	GOL	A	609	-	-	2/4/4/4	-
3	GOL	B	612	-	-	4/4/4/4	-
3	GOL	B	614	-	-	4/4/4/4	-
3	GOL	C	616	-	-	1/4/4/4	-
3	GOL	C	615	-	-	2/4/4/4	-
3	GOL	C	618	-	-	3/4/4/4	-
3	GOL	C	619	-	-	2/4/4/4	-
3	GOL	C	613	-	-	0/4/4/4	-
3	GOL	A	610	-	-	2/4/4/4	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	GOL	B	613	-	-	4/4/4/4	-
3	GOL	A	612	-	-	2/4/4/4	-
3	GOL	C	617	-	-	2/4/4/4	-
3	GOL	C	614	-	-	3/4/4/4	-
3	GOL	A	611	-	-	0/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (31) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	609	GOL	C1-C2-C3-O3
3	A	610	GOL	O1-C1-C2-O2
3	A	610	GOL	O1-C1-C2-C3
3	B	612	GOL	O1-C1-C2-O2
3	B	612	GOL	O1-C1-C2-C3
3	B	612	GOL	C1-C2-C3-O3
3	B	613	GOL	O1-C1-C2-C3
3	B	613	GOL	C1-C2-C3-O3
3	C	614	GOL	O1-C1-C2-C3
3	C	614	GOL	C1-C2-C3-O3
3	C	615	GOL	C1-C2-C3-O3
3	C	618	GOL	C1-C2-C3-O3
3	C	619	GOL	O1-C1-C2-C3
3	B	614	GOL	O2-C2-C3-O3
3	C	615	GOL	O2-C2-C3-O3
3	A	612	GOL	C1-C2-C3-O3
3	B	614	GOL	O1-C1-C2-C3
3	B	614	GOL	C1-C2-C3-O3
3	C	617	GOL	O1-C1-C2-C3
3	C	618	GOL	O1-C1-C2-C3
3	A	609	GOL	O2-C2-C3-O3
3	C	618	GOL	O2-C2-C3-O3
3	C	619	GOL	O1-C1-C2-O2
3	B	612	GOL	O2-C2-C3-O3
3	B	613	GOL	O2-C2-C3-O3
3	C	614	GOL	O1-C1-C2-O2
3	C	616	GOL	O1-C1-C2-C3
3	A	612	GOL	O2-C2-C3-O3

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Mol	Chain	Res	Type	Atoms
3	B	614	GOL	O1-C1-C2-O2
3	C	617	GOL	O1-C1-C2-O2
3	B	613	GOL	O1-C1-C2-O2

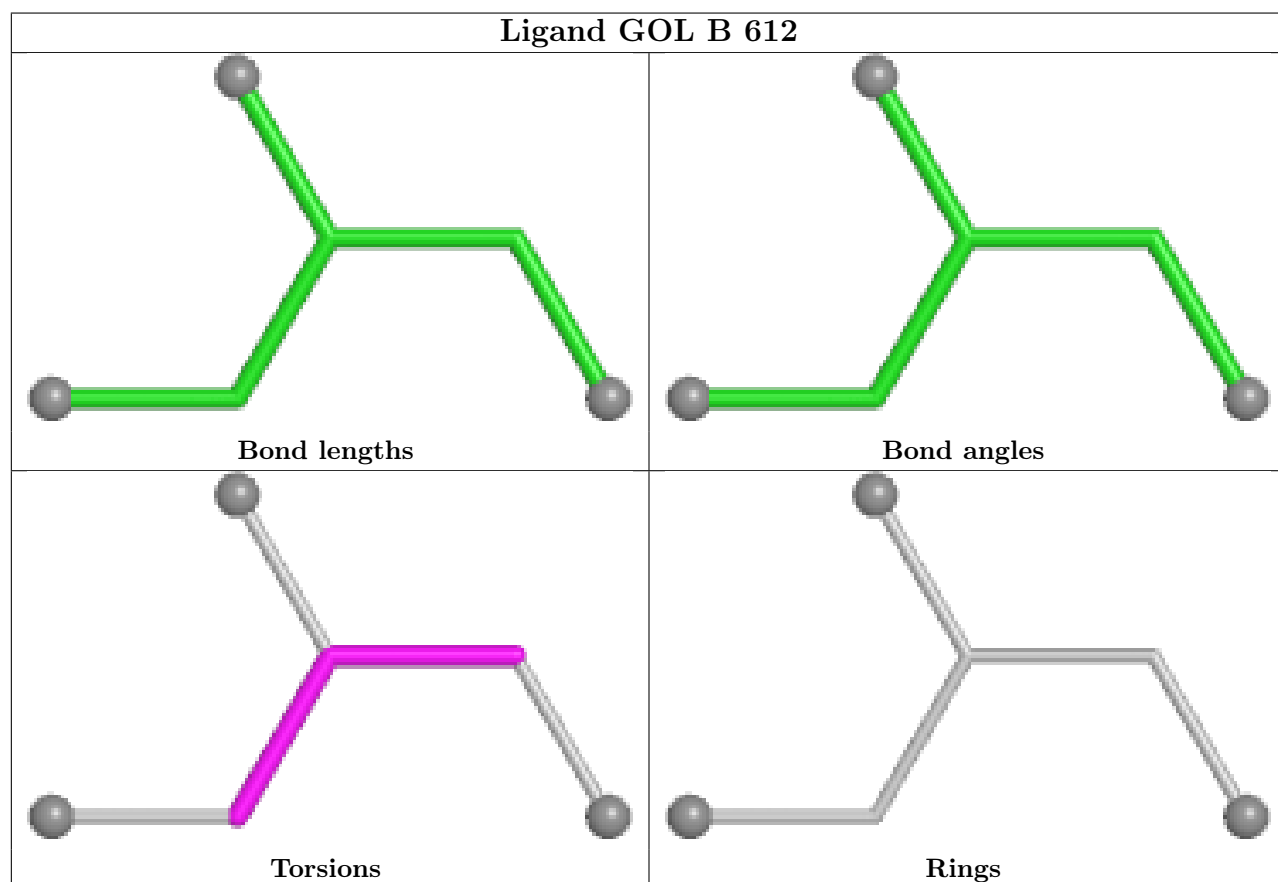
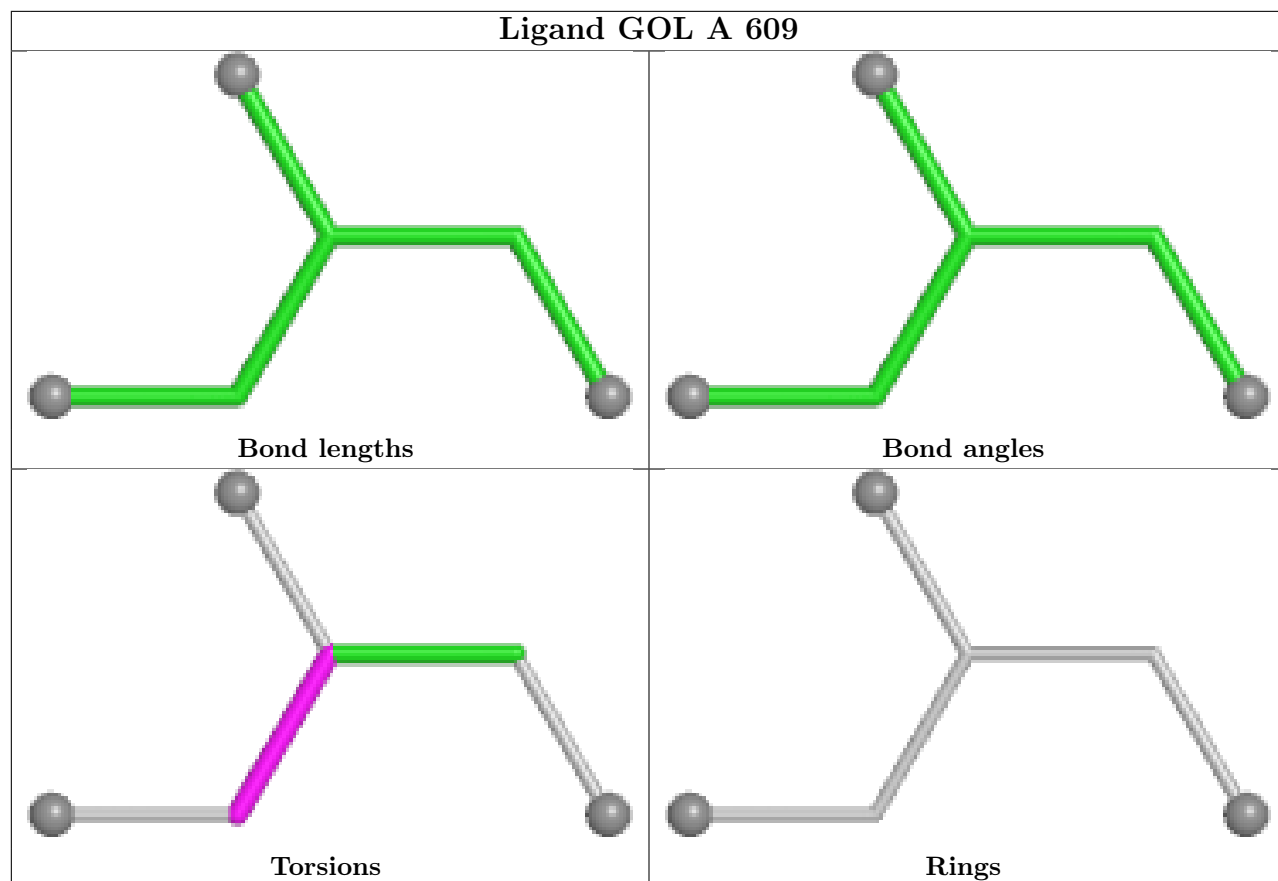
There are no ring outliers.

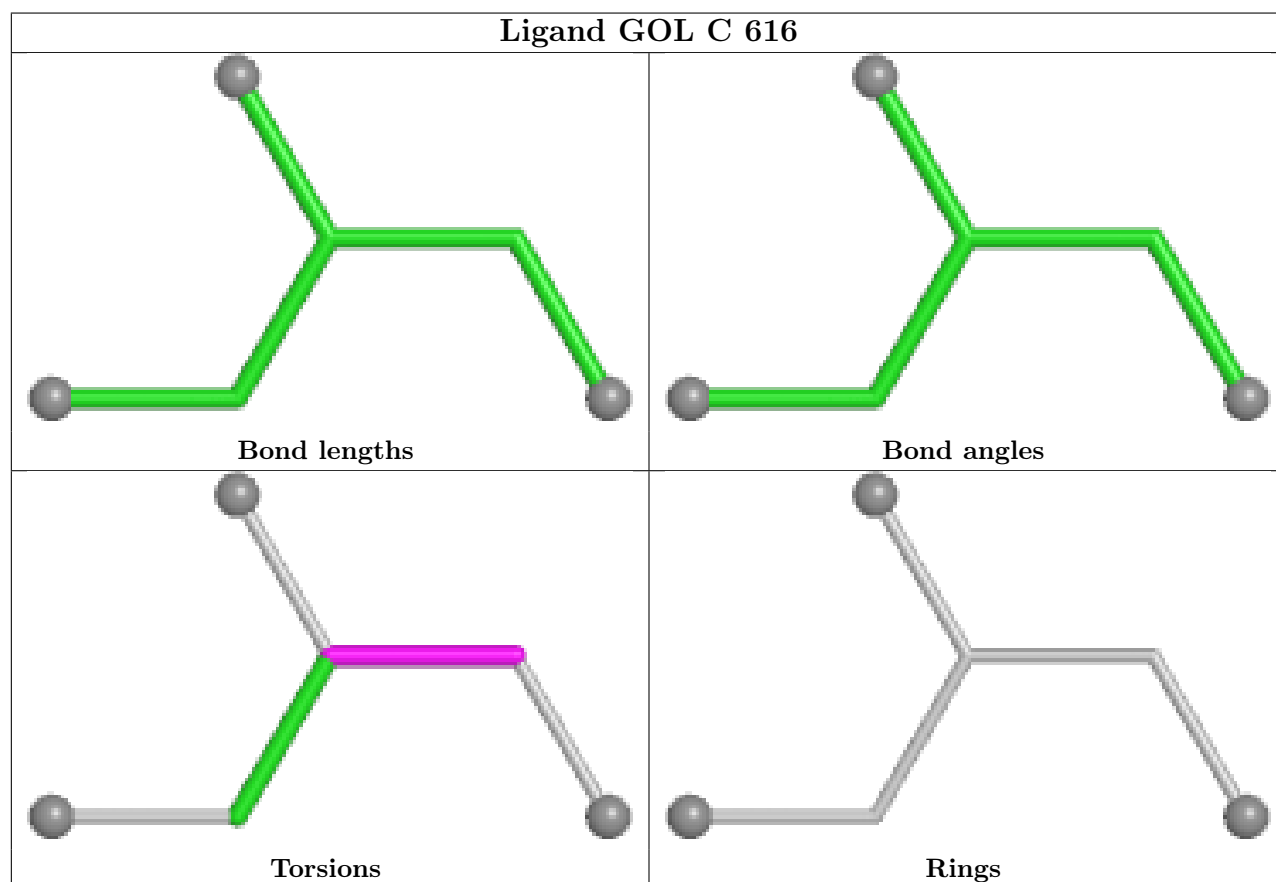
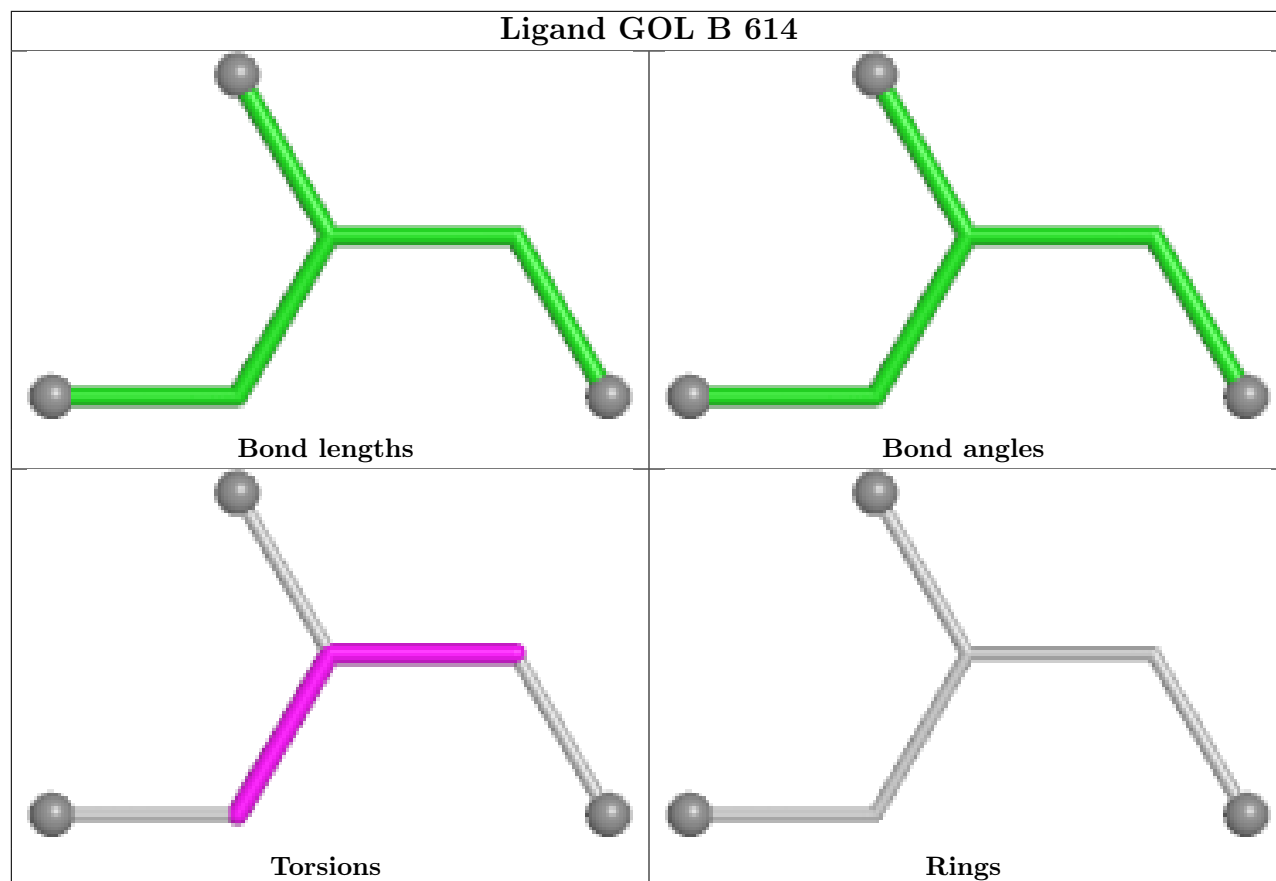
7 monomers are involved in 13 short contacts:

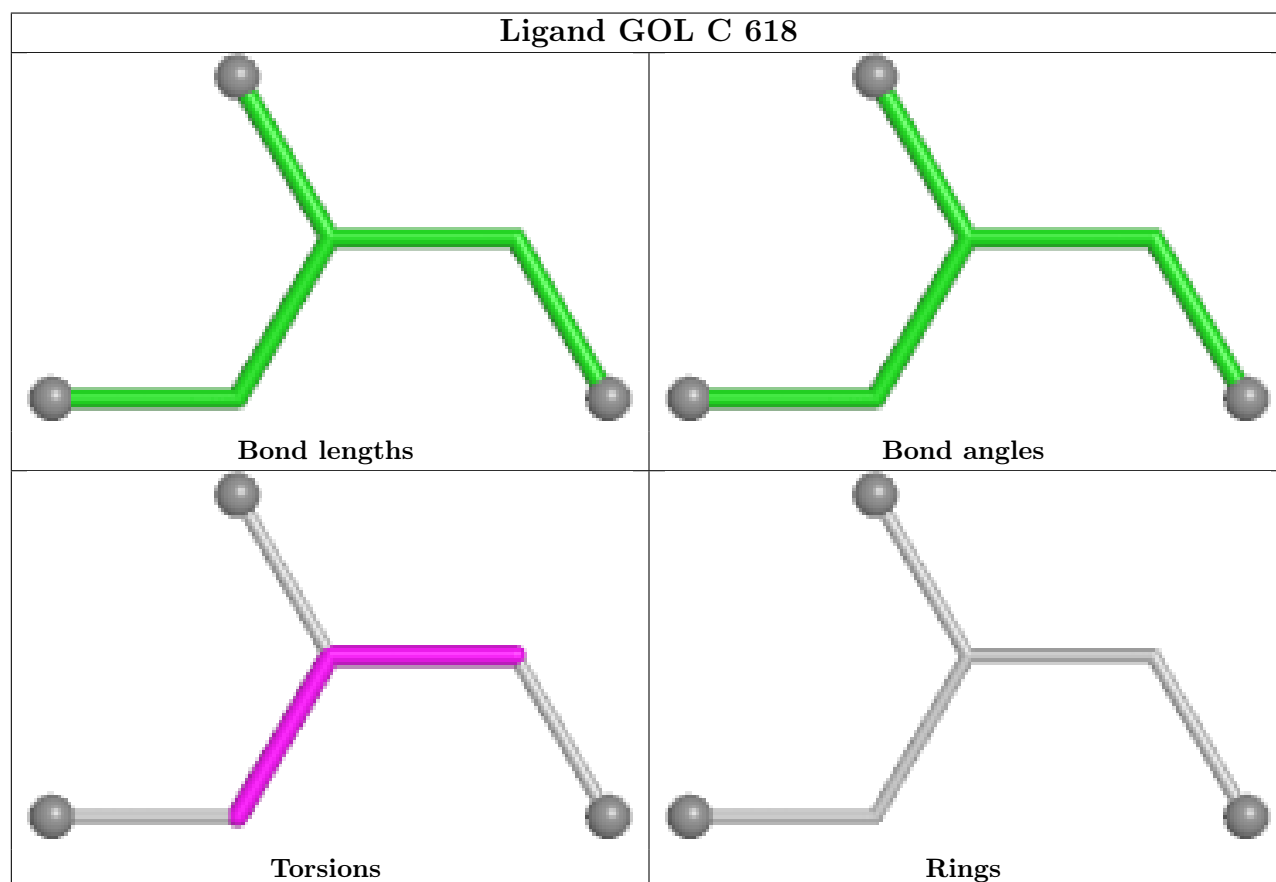
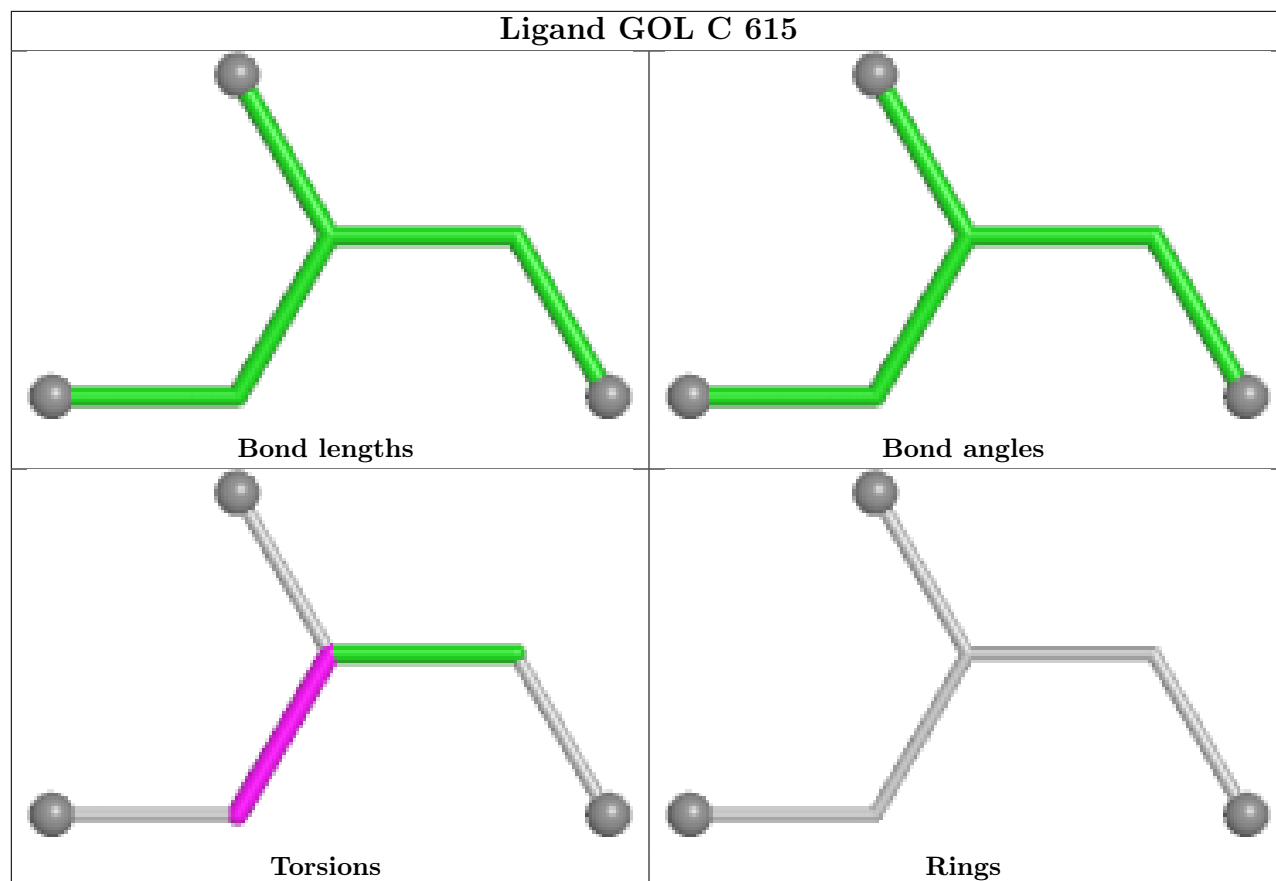
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	612	GOL	3	0
3	C	618	GOL	1	0
3	C	613	GOL	1	0
3	A	610	GOL	2	0
3	C	617	GOL	1	0
3	C	614	GOL	4	0
3	A	611	GOL	1	0

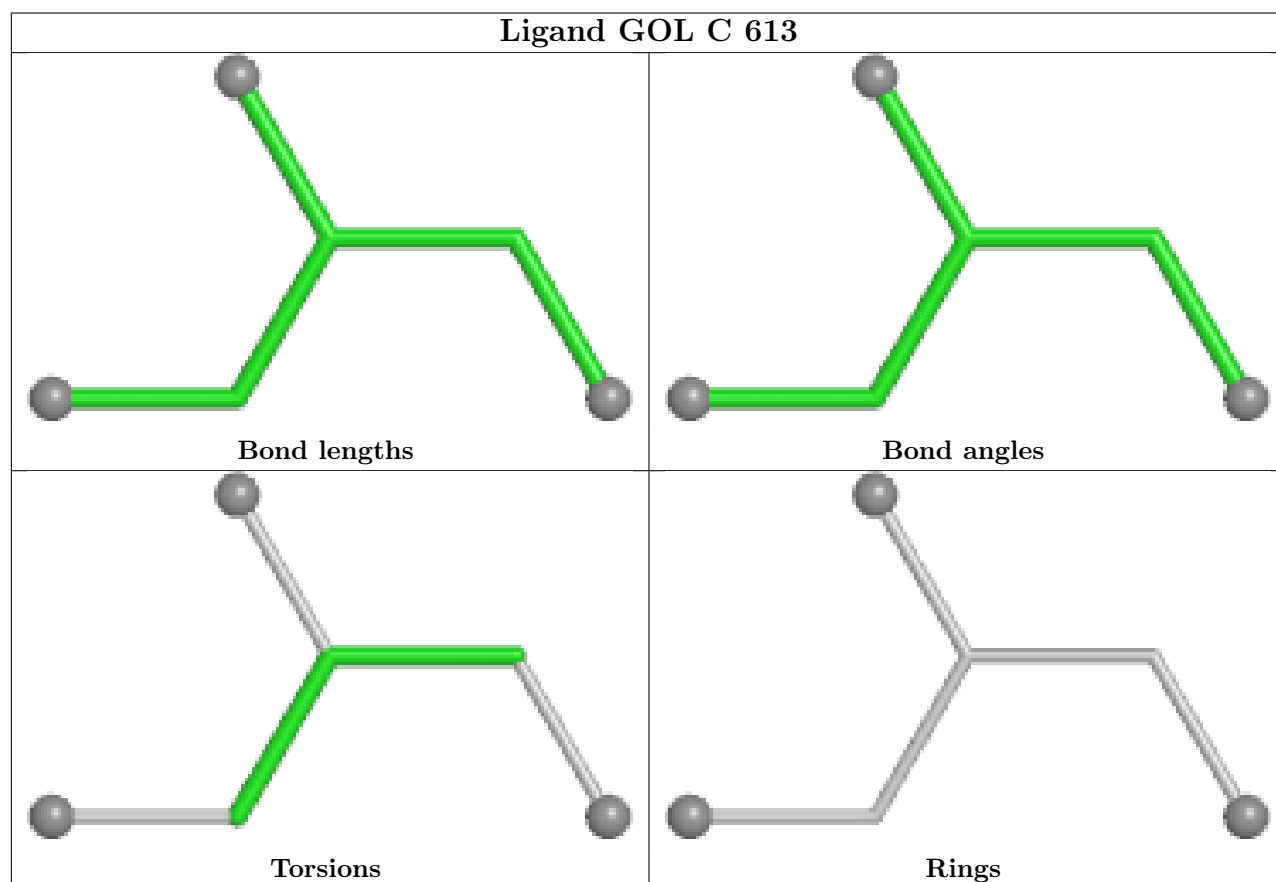
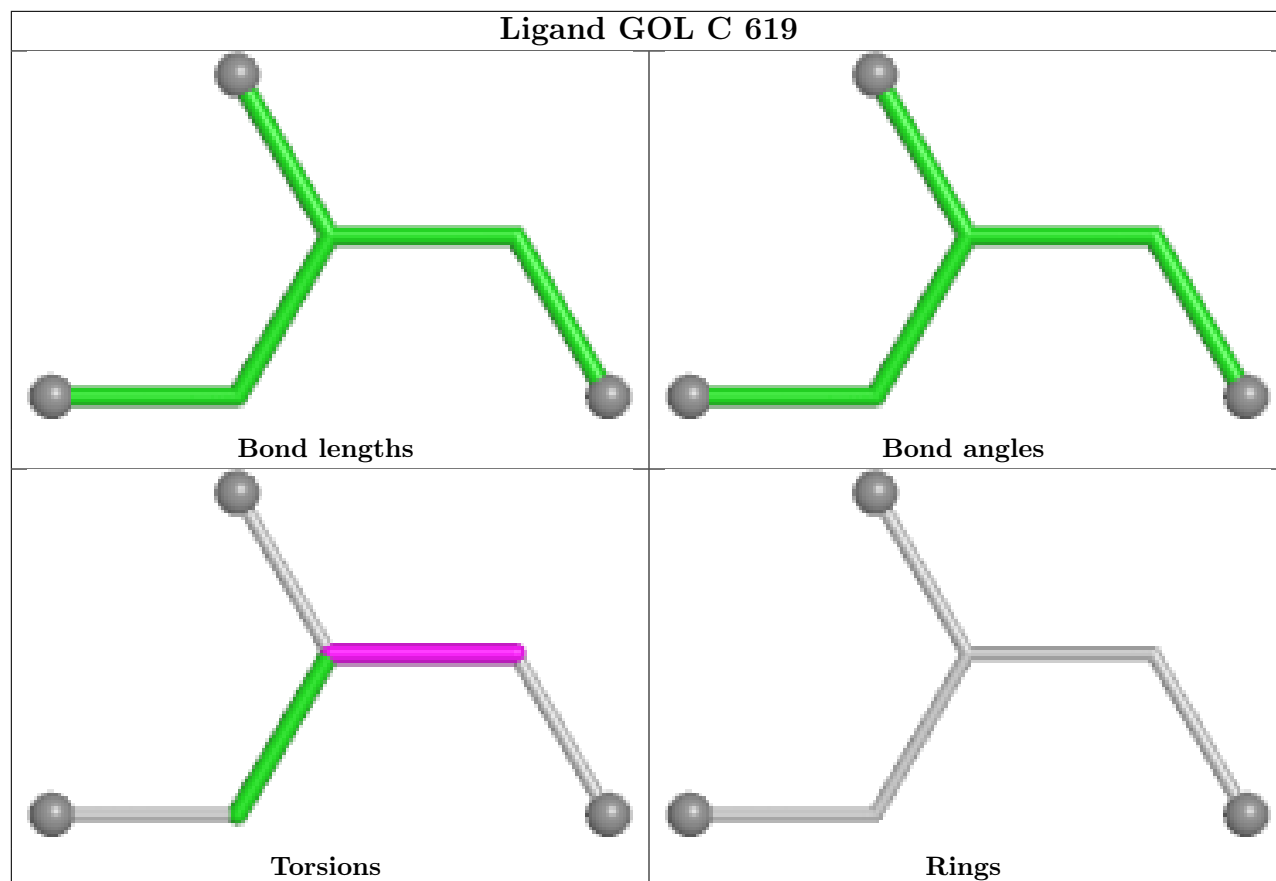
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

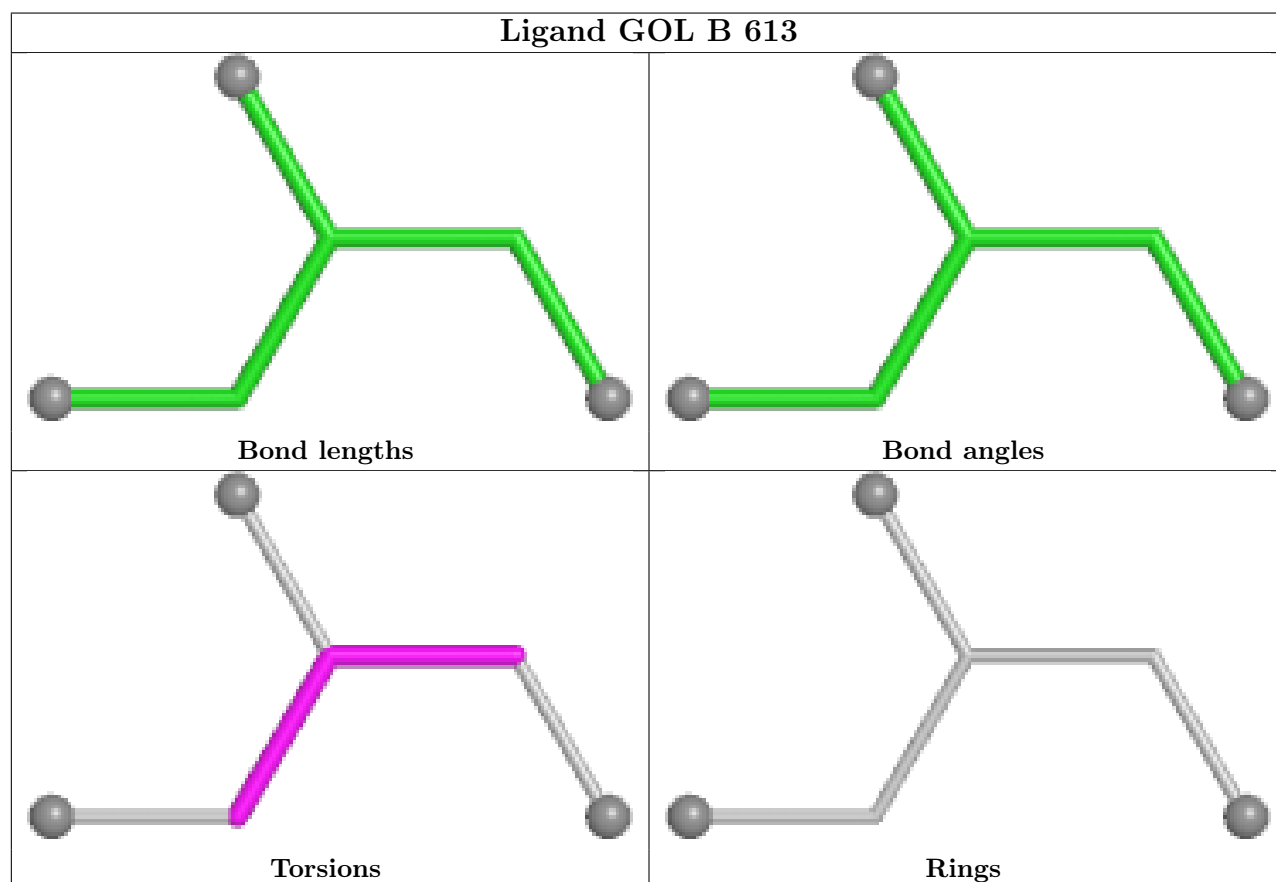
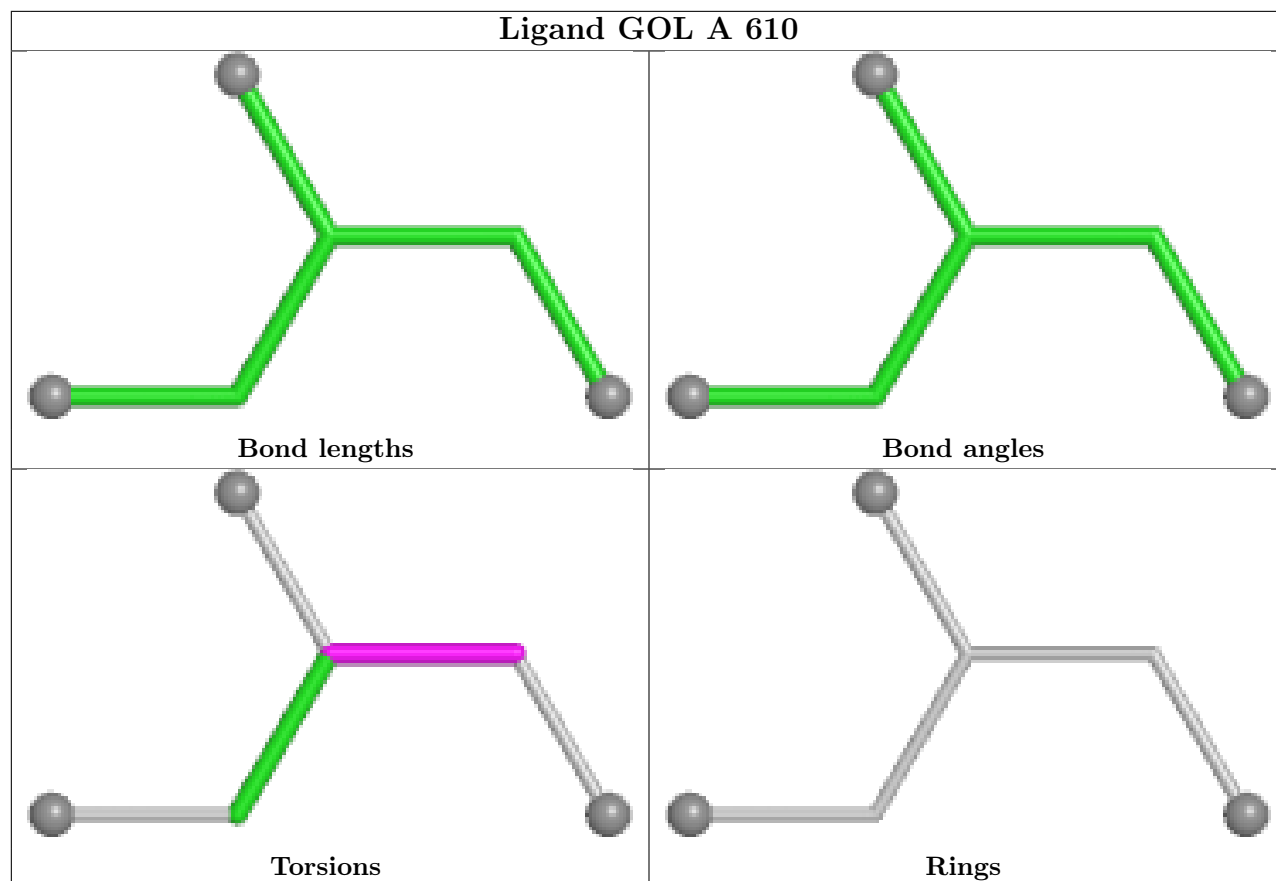


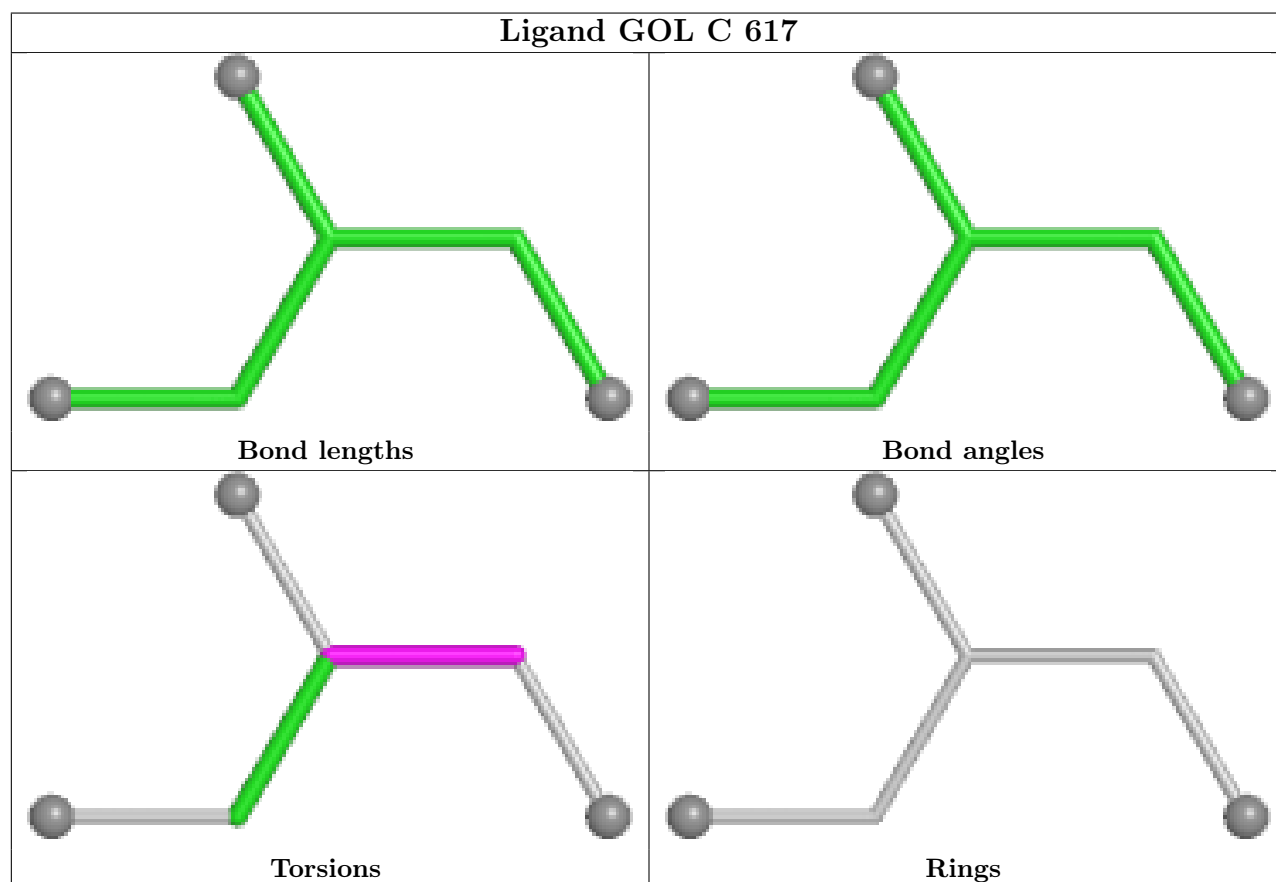
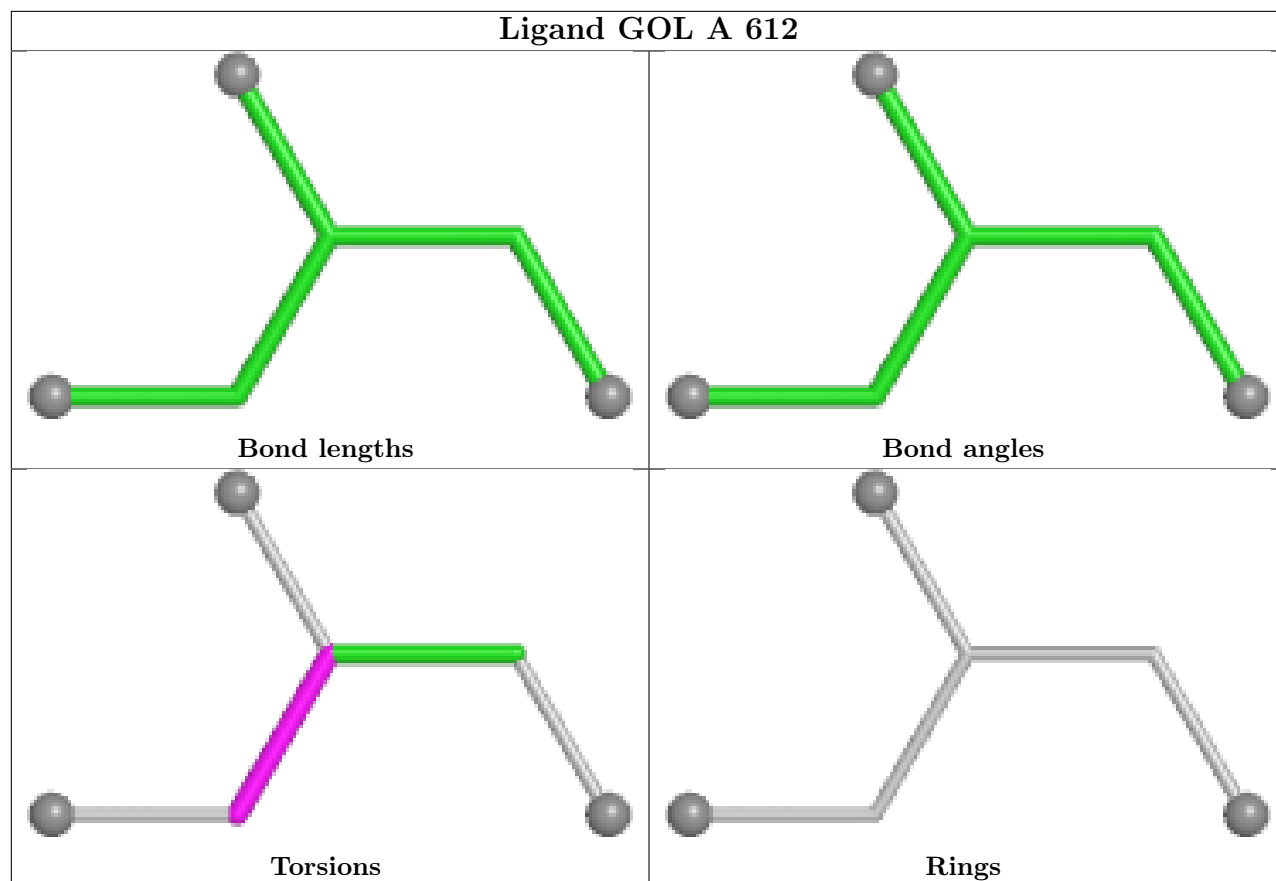


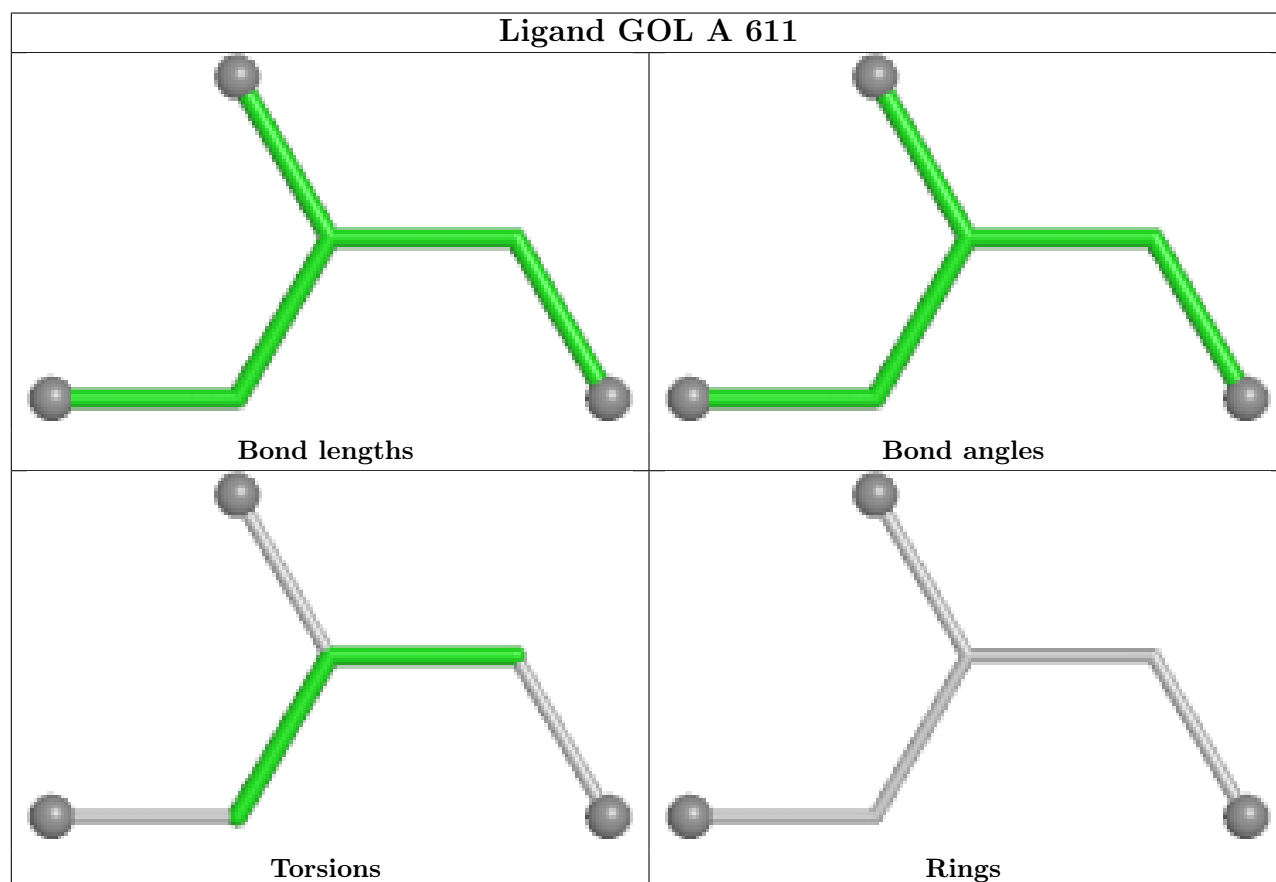
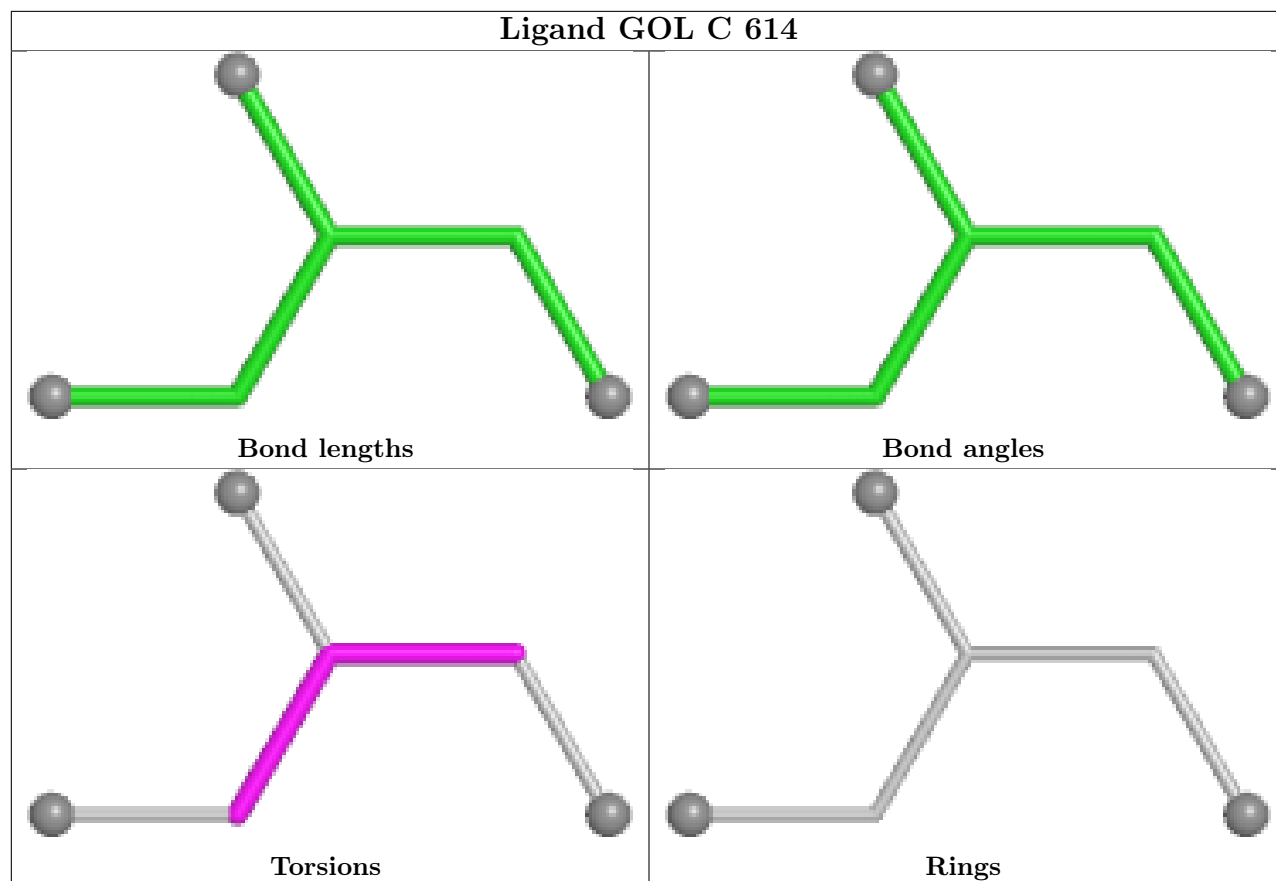












## 5.7 Other polymers [i](#)

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	524/530 (98%)	-0.24	18 (3%)	48 42	21, 41, 83, 145	3 (0%)
1	B	501/530 (94%)	0.06	8 (1%)	70 67	26, 55, 93, 113	3 (0%)
1	C	503/530 (94%)	-0.35	5 (0%)	79 76	23, 38, 68, 109	3 (0%)
1	D	511/530 (96%)	0.01	14 (2%)	56 51	26, 52, 95, 117	3 (0%)
All	All	2039/2120 (96%)	-0.13	45 (2%)	62 58	21, 46, 90, 145	12 (0%)

All (45) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	467	VAL	5.2
1	A	464	PHE	4.3
1	B	457	THR	4.2
1	A	461	LEU	4.2
1	D	427	GLY	4.1
1	B	524	VAL	3.7
1	D	524	VAL	3.6
1	A	463	TRP	3.5
1	B	480	GLY	3.5
1	A	470	ALA	3.4
1	C	478	VAL	3.4
1	C	479	PRO	3.4
1	A	458	TYR	3.3
1	C	457	THR	3.2
1	A	460	GLY	3.1
1	A	468	LEU	2.9
1	B	456	GLY	2.9
1	D	350	ALA	2.8
1	D	476	LEU	2.8
1	D	475	GLY	2.7
1	C	456	GLY	2.7

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Mol	Chain	Res	Type	RSRZ
1	A	24	PRO	2.6
1	A	26	SER	2.6
1	B	508	GLN	2.6
1	A	459	GLY	2.6
1	D	351	GLY	2.6
1	B	350	ALA	2.5
1	D	470	ALA	2.5
1	D	474	PRO	2.5
1	D	457	THR	2.5
1	A	29	LEU	2.5
1	C	480	GLY	2.4
1	B	427	GLY	2.4
1	A	427	GLY	2.4
1	A	525	SER	2.3
1	D	428	ASP	2.3
1	A	456	GLY	2.2
1	D	517	VAL	2.2
1	B	484	ARG	2.2
1	D	508	GLN	2.2
1	A	475	GLY	2.2
1	A	479	PRO	2.1
1	D	479	PRO	2.1
1	D	335[A]	GLN	2.1
1	A	478	VAL	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 oligosaccharides 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.

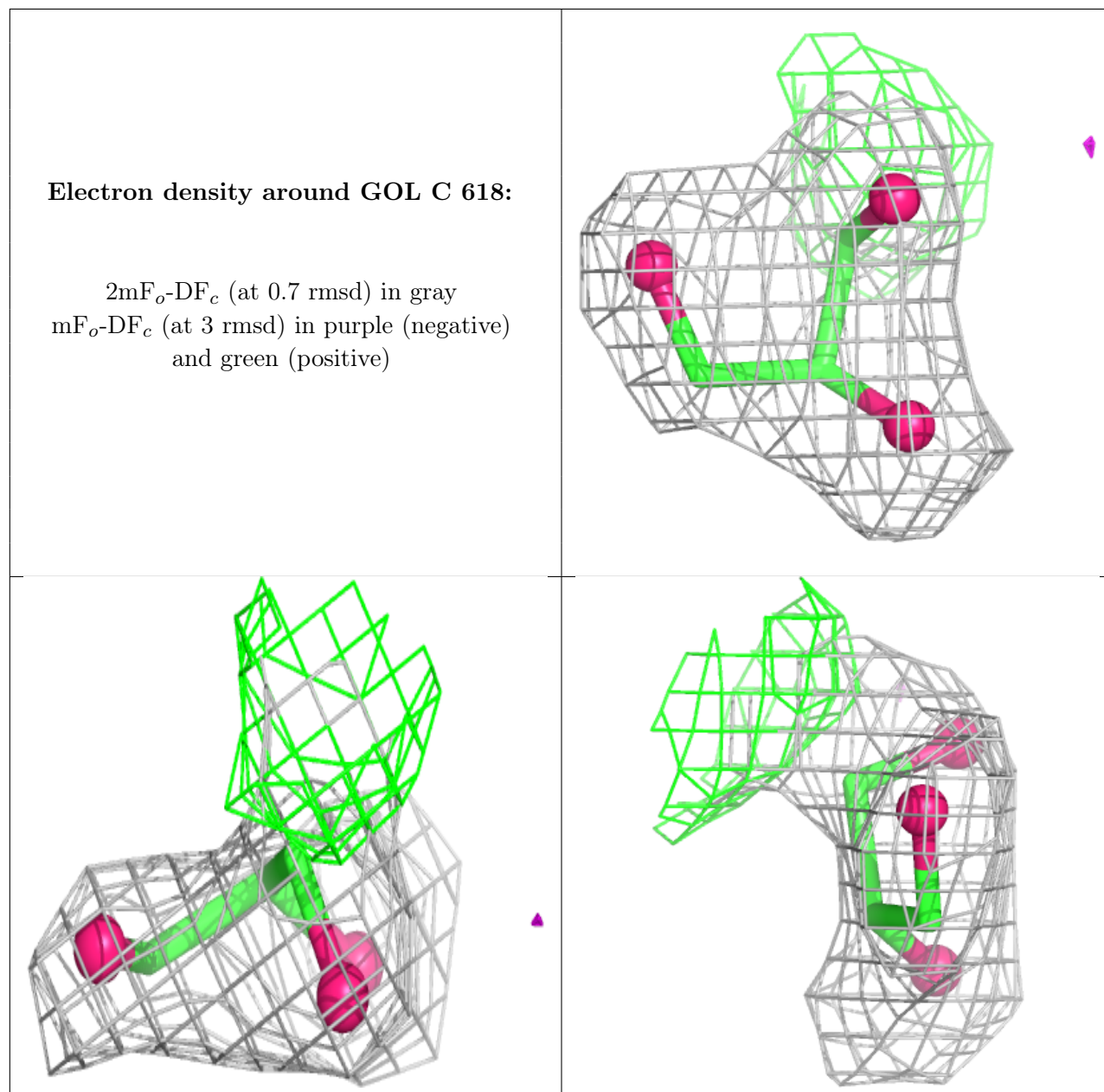
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
3	GOL	C	618	6/6	0.45	0.20	77,79,79,79	0
3	GOL	B	614	6/6	0.65	0.16	67,68,69,69	0
2	MG	B	609	1/1	0.72	0.26	61,61,61,61	0
3	GOL	A	612	6/6	0.76	0.20	59,62,62,63	0
3	GOL	C	616	6/6	0.77	0.17	62,65,66,66	0
3	GOL	A	610	6/6	0.81	0.21	24,28,29,30	6
3	GOL	B	612	6/6	0.82	0.24	75,76,77,78	0
3	GOL	C	615	6/6	0.83	0.14	66,67,69,69	0
3	GOL	C	613	6/6	0.84	0.17	63,65,66,66	0
2	MG	C	611	1/1	0.85	0.13	57,57,57,57	0
3	GOL	B	613	6/6	0.86	0.14	65,66,67,68	0
2	MG	A	607	1/1	0.86	0.22	63,63,63,63	0
2	MG	B	610	1/1	0.86	0.22	62,62,62,62	0
3	GOL	A	611	6/6	0.87	0.16	48,50,51,51	0
3	GOL	C	614	6/6	0.88	0.26	53,55,56,57	0
2	MG	C	604	1/1	0.88	0.14	60,60,60,60	0
3	GOL	C	619	6/6	0.88	0.16	37,39,40,41	6
2	MG	B	611	1/1	0.89	0.12	55,55,55,55	0
3	GOL	C	617	6/6	0.89	0.16	52,54,55,56	0
4	CL	C	612	1/1	0.89	0.18	91,91,91,91	0
2	MG	A	608	1/1	0.91	0.09	52,52,52,52	0
2	MG	A	602	1/1	0.91	0.12	48,48,48,48	0
2	MG	B	607	1/1	0.92	0.13	64,64,64,64	0
2	MG	A	601	1/1	0.92	0.08	44,44,44,44	0
3	GOL	A	609	6/6	0.93	0.10	55,56,56,56	0
2	MG	B	601	1/1	0.93	0.12	64,64,64,64	0
2	MG	B	604	1/1	0.93	0.17	68,68,68,68	0
2	MG	D	602	1/1	0.94	0.10	44,44,44,44	0
2	MG	C	605	1/1	0.94	0.14	62,62,62,62	0
2	MG	A	603	1/1	0.95	0.10	40,40,40,40	0
2	MG	B	605	1/1	0.95	0.16	48,48,48,48	0
2	MG	C	607	1/1	0.95	0.12	39,39,39,39	0
2	MG	B	602	1/1	0.95	0.10	51,51,51,51	0
2	MG	B	608	1/1	0.96	0.09	31,31,31,31	0
2	MG	D	601	1/1	0.96	0.13	44,44,44,44	0
2	MG	B	603	1/1	0.96	0.08	58,58,58,58	0
2	MG	C	601	1/1	0.96	0.09	41,41,41,41	0
2	MG	C	609	1/1	0.96	0.07	49,49,49,49	0
2	MG	C	610	1/1	0.96	0.09	51,51,51,51	0
2	MG	A	605	1/1	0.97	0.07	42,42,42,42	0
2	MG	B	606	1/1	0.97	0.08	40,40,40,40	0
2	MG	C	602	1/1	0.97	0.15	36,36,36,36	0
2	MG	C	608	1/1	0.97	0.08	41,41,41,41	0

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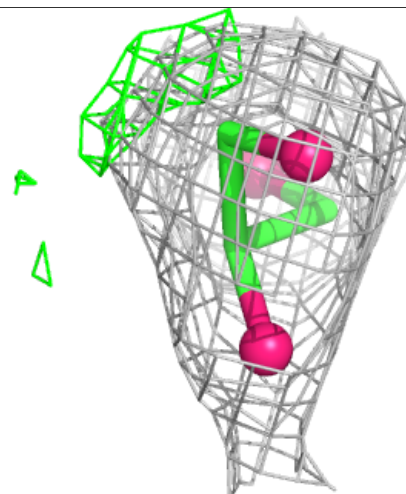
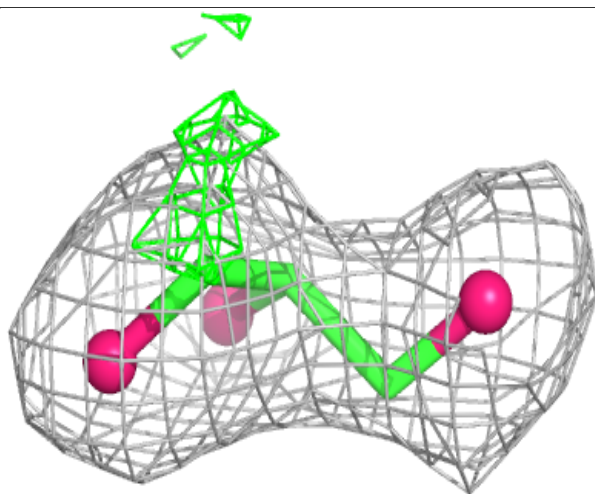
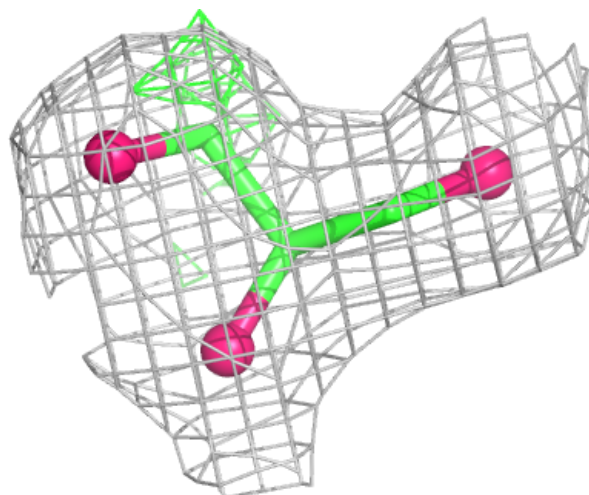
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
2	MG	C	603	1/1	0.98	0.11	47,47,47,47	0
2	MG	C	606	1/1	0.98	0.12	42,42,42,42	0
2	MG	A	606	1/1	0.98	0.05	40,40,40,40	0
2	MG	A	604	1/1	0.99	0.06	25,25,25,25	0

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.



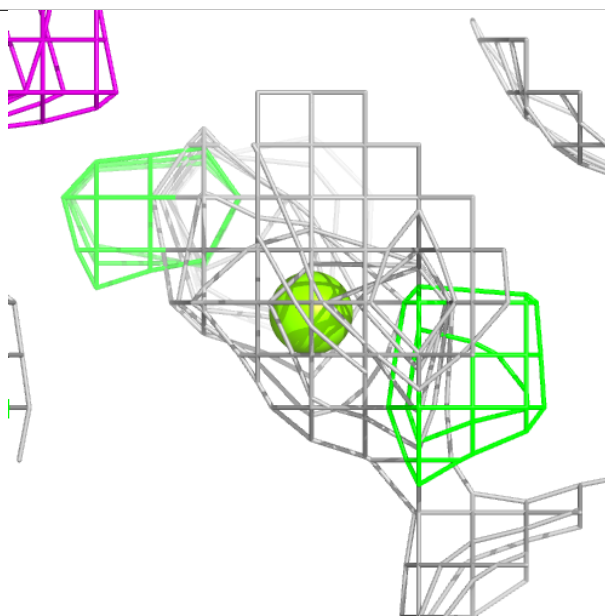
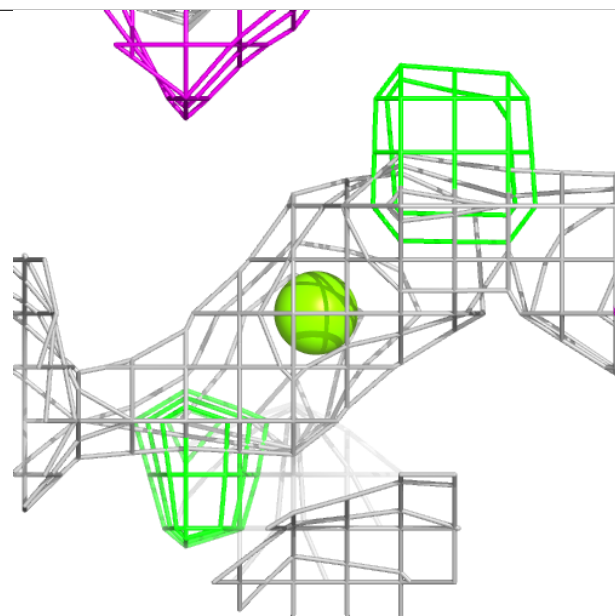
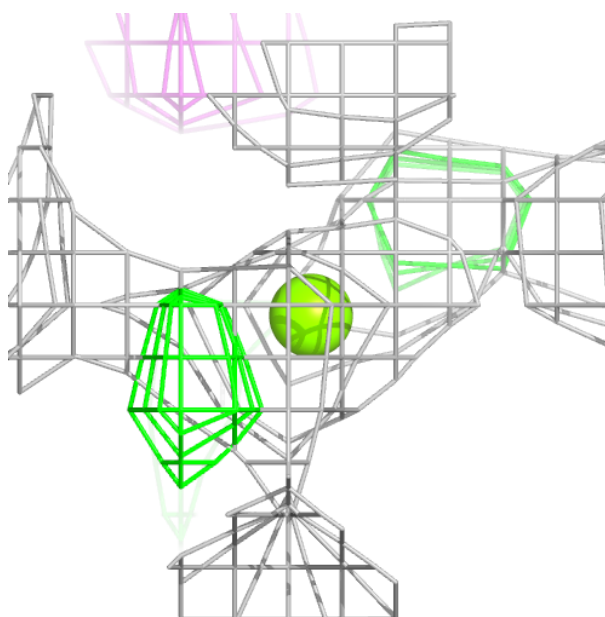
**Electron density around GOL B 614:**

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



**Electron density around MG B 609:**

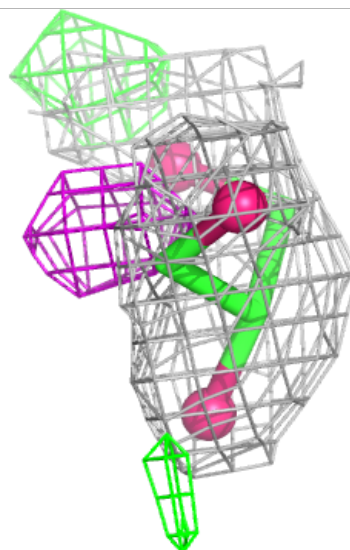
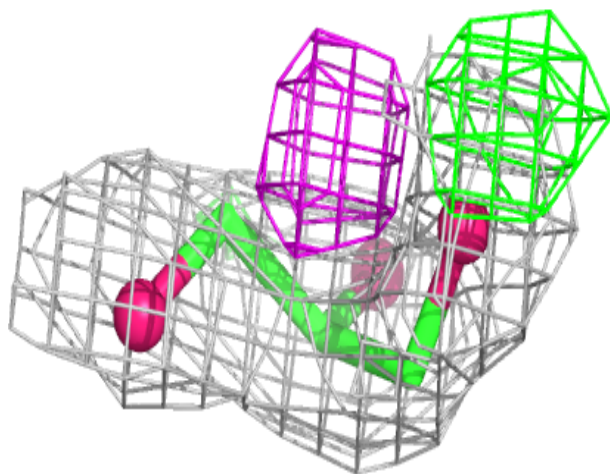
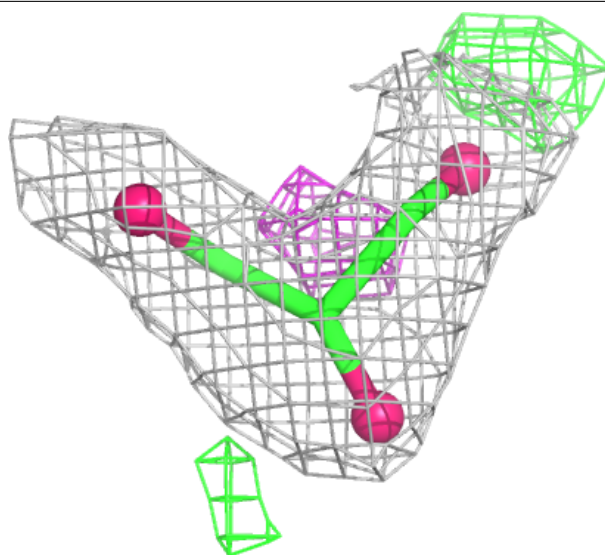
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)





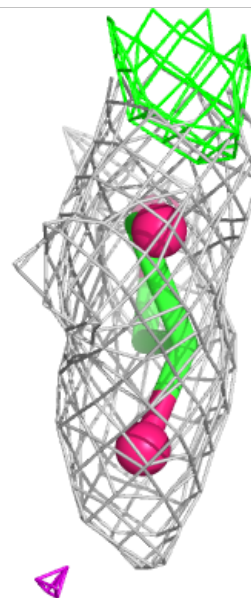
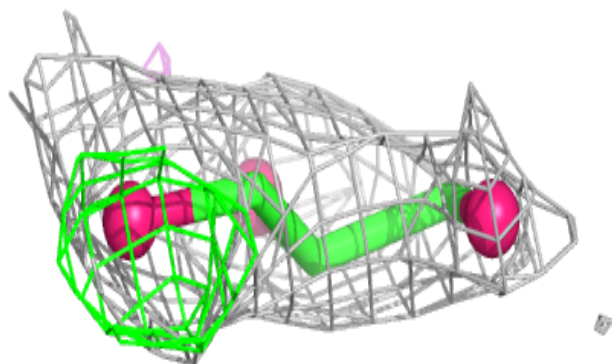
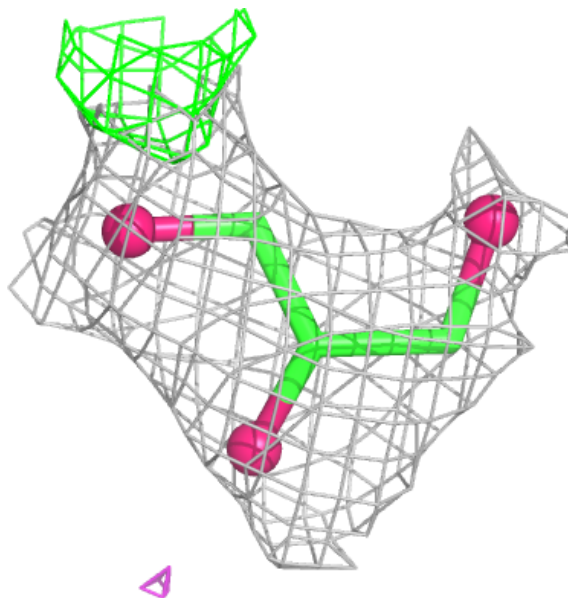
**Electron density around GOL A 612:**

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



**Electron density around GOL C 616:**

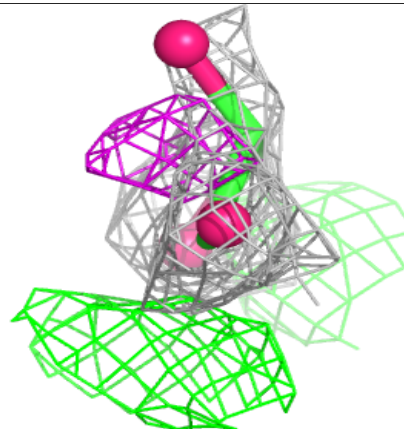
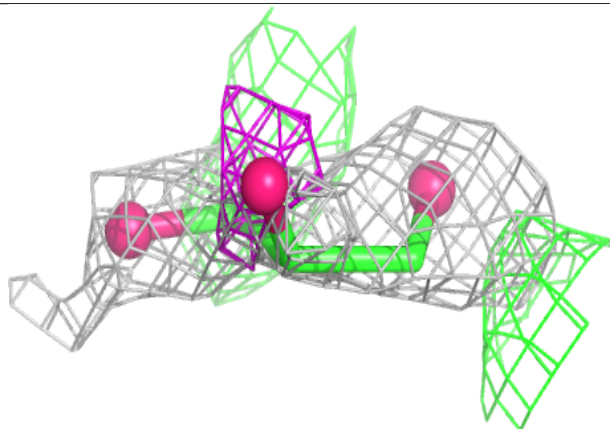
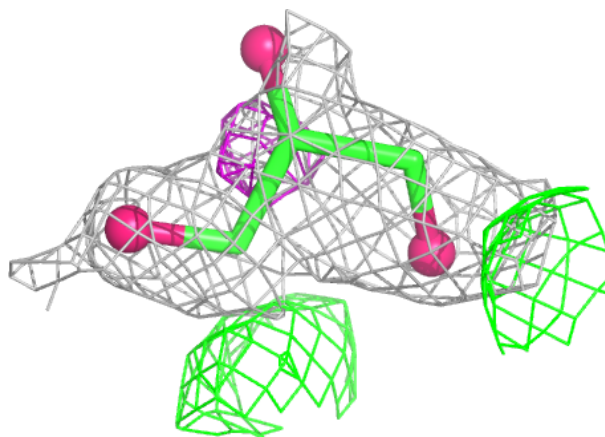
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)



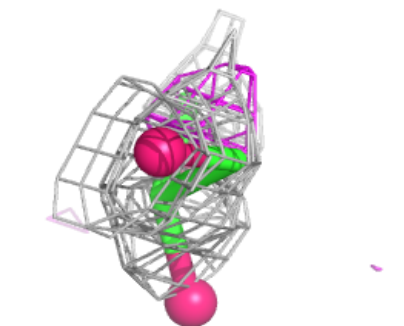
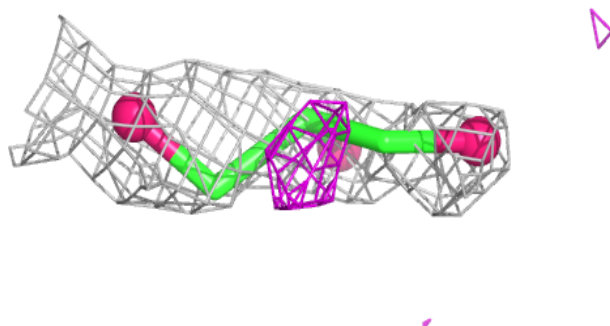
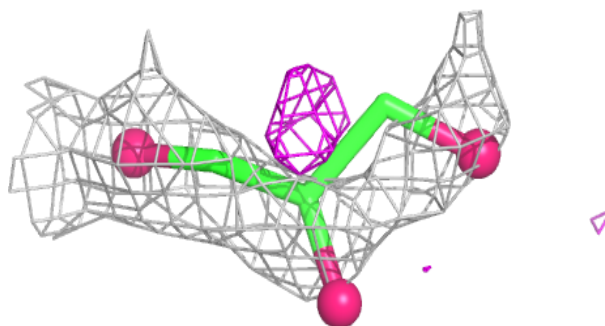


**Electron density around GOL A 610:**

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

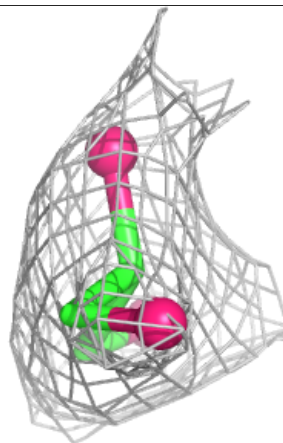
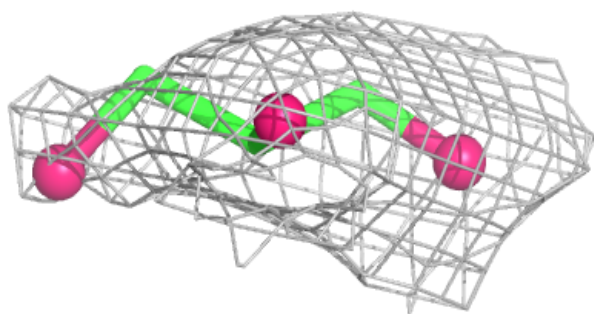
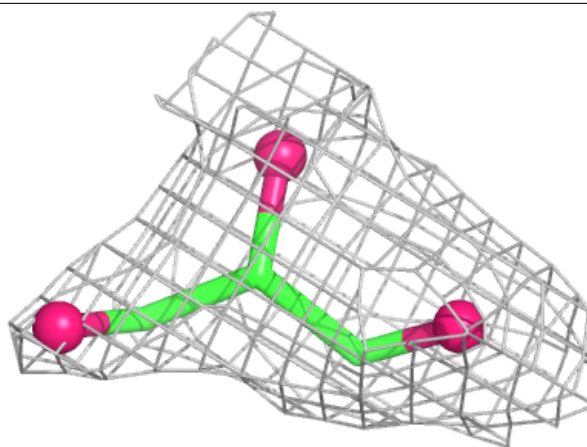
**Electron density around GOL B 612:**

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



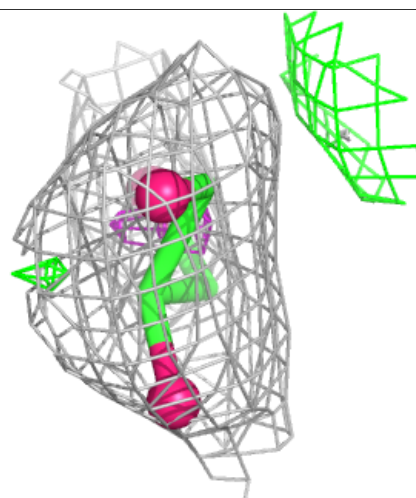
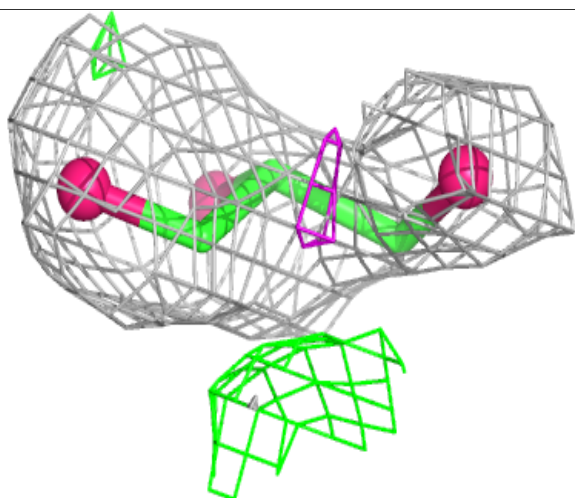
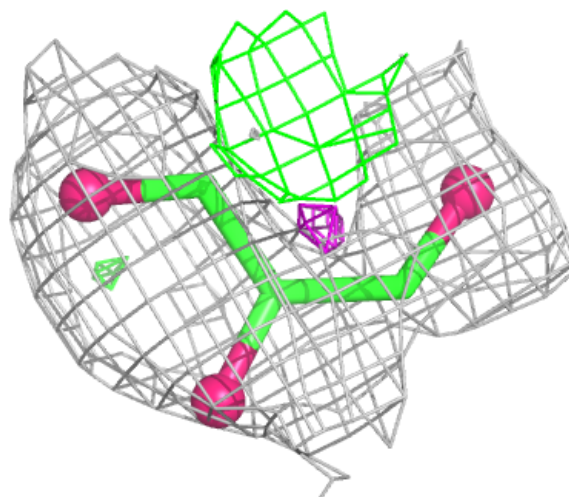
**Electron density around GOL C 615:**

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



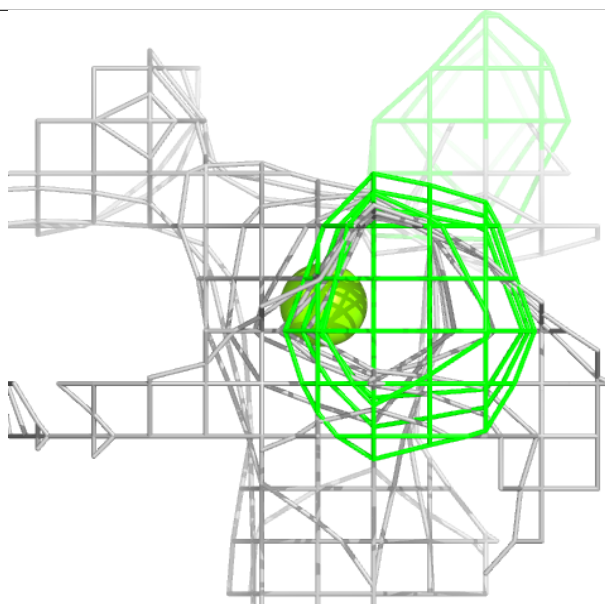
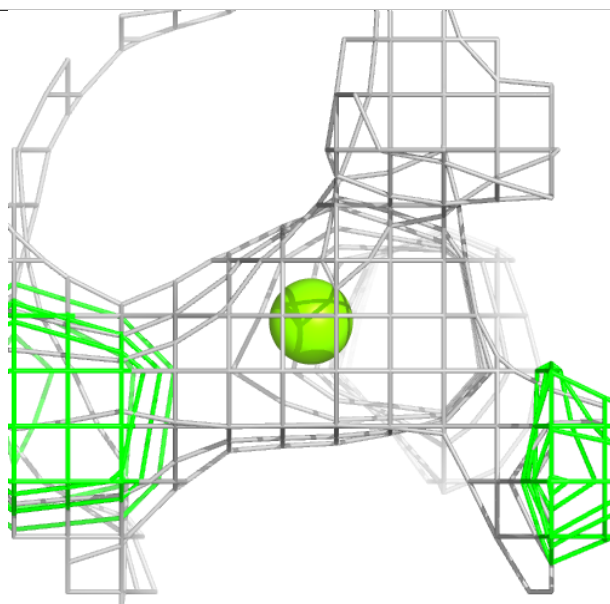
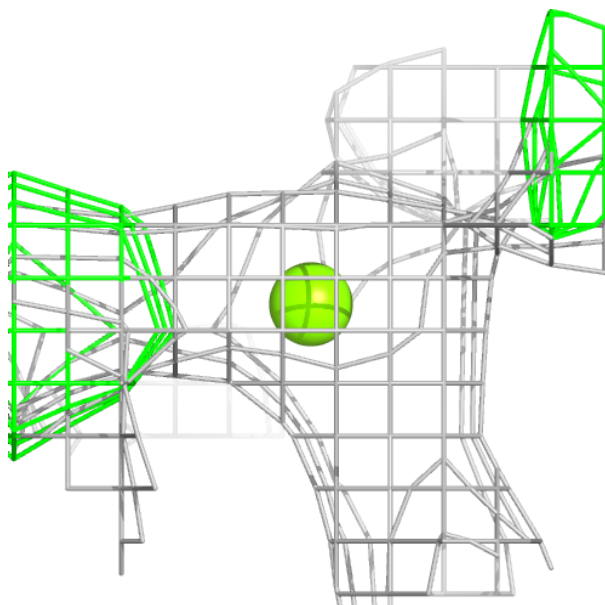
**Electron density around GOL C 613:**

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



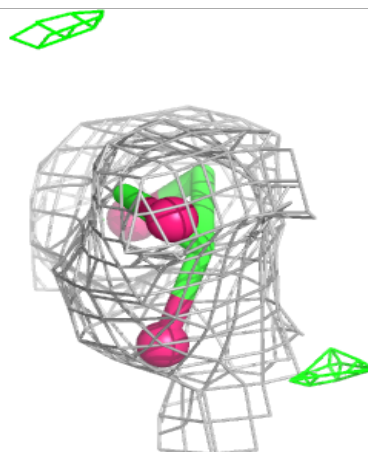
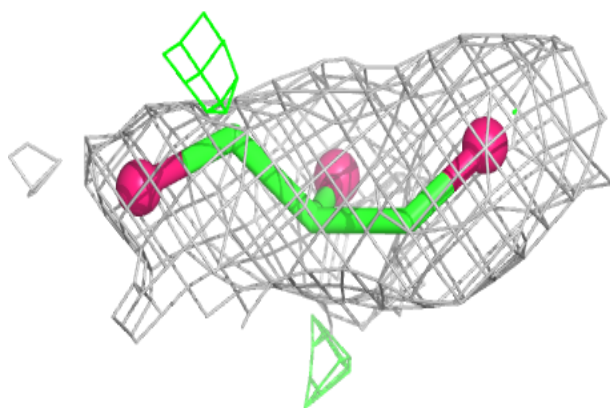
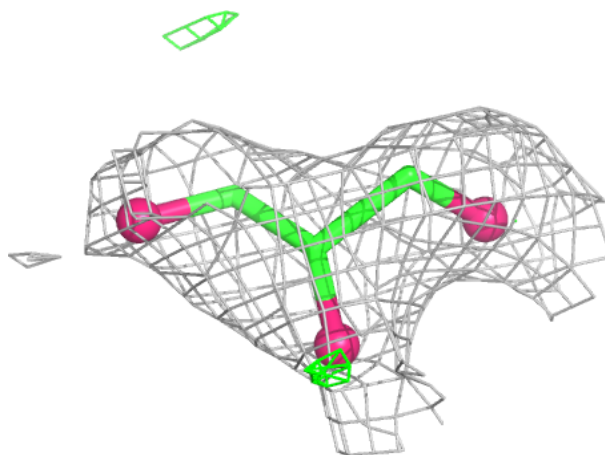
**Electron density around MG C 611:**

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



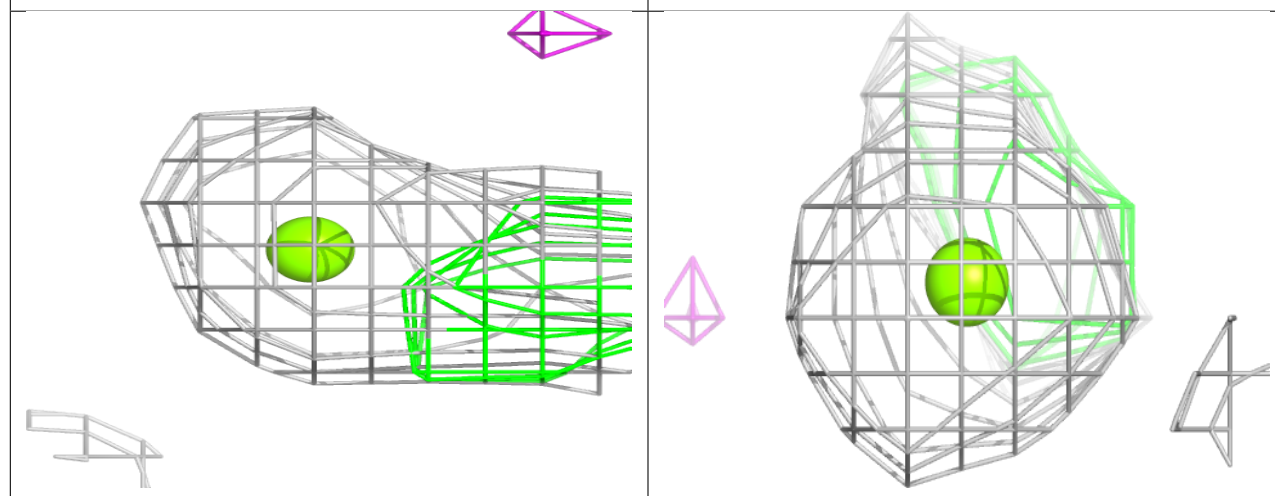
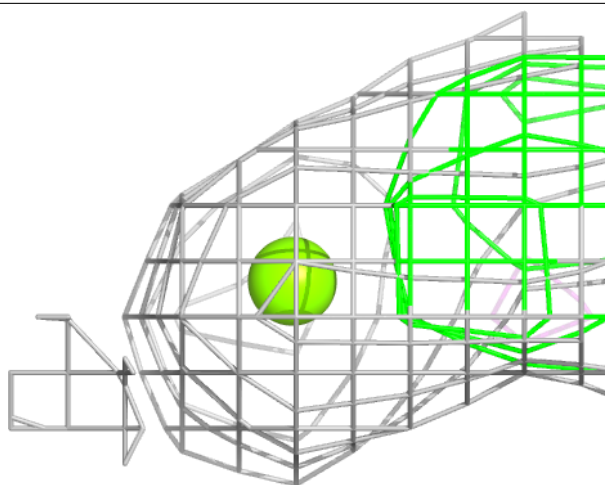
**Electron density around GOL B 613:**

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



**Electron density around MG A 607:**

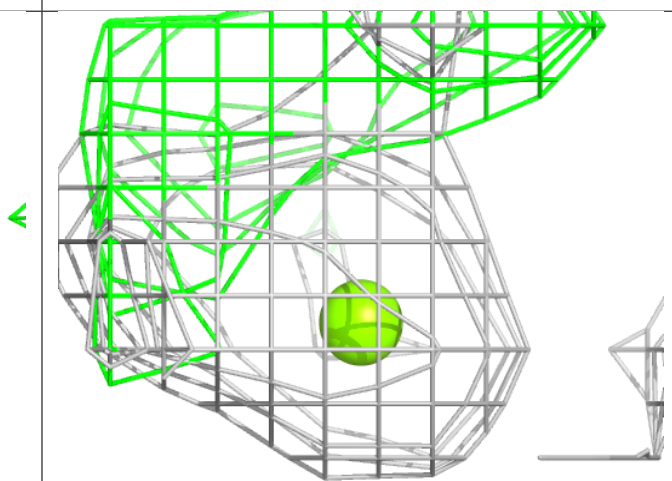
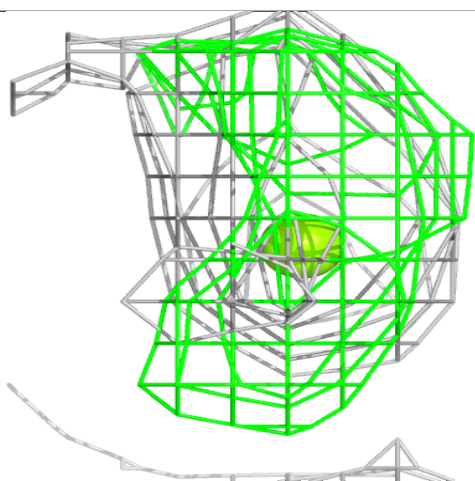
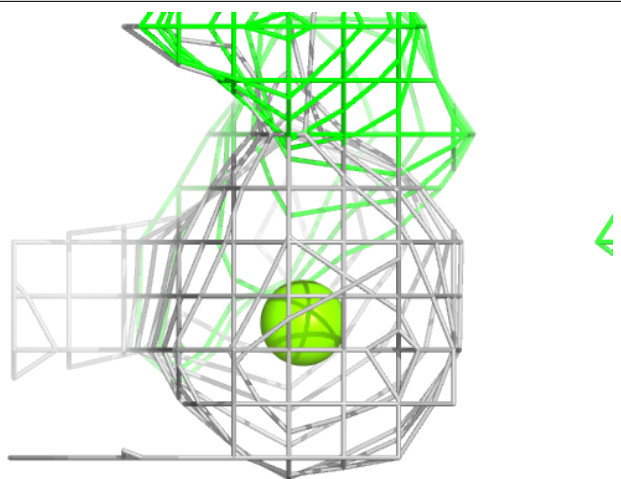
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)





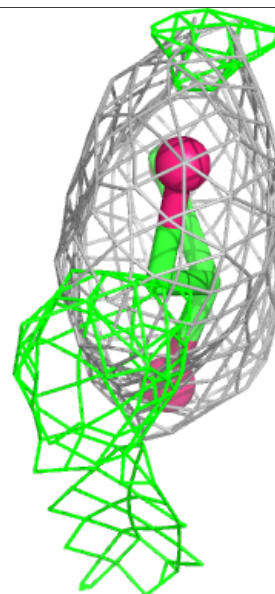
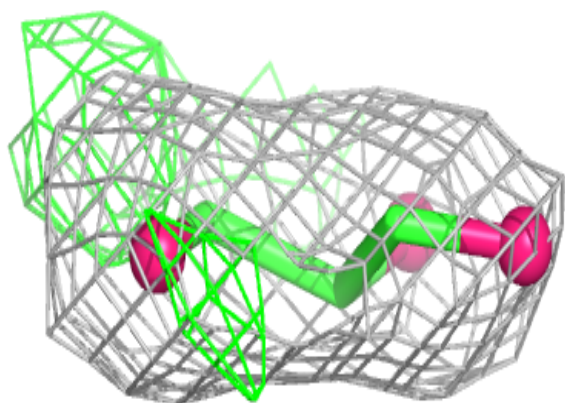
**Electron density around MG B 610:**

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



**Electron density around GOL A 611:**

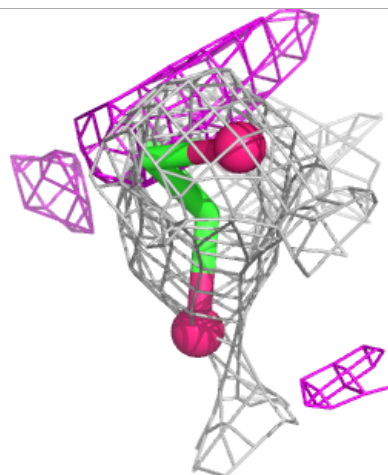
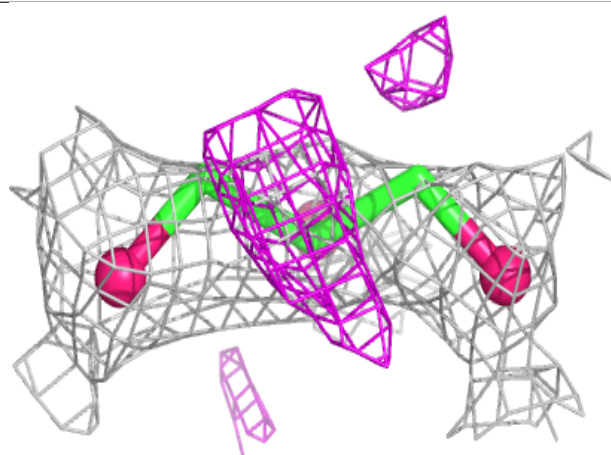
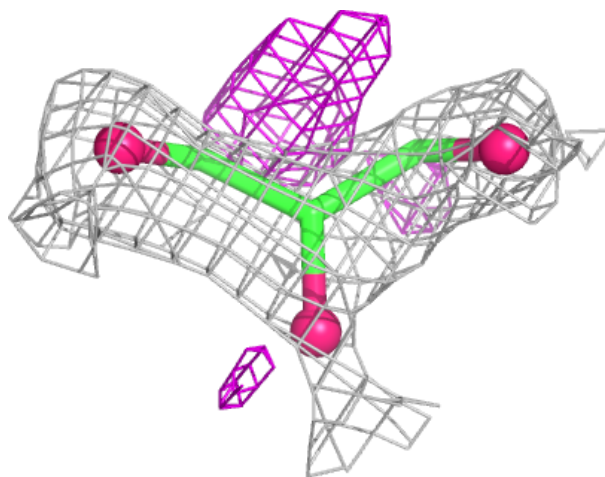
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)





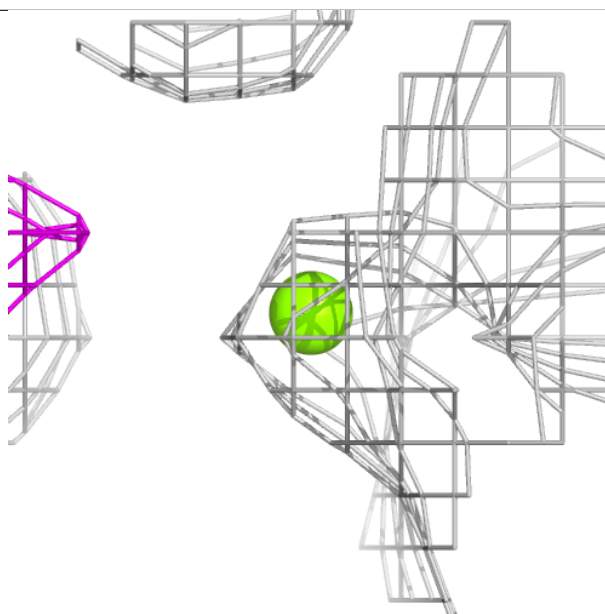
**Electron density around GOL C 614:**

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



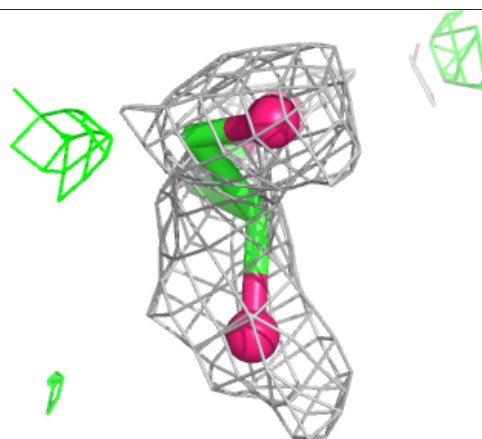
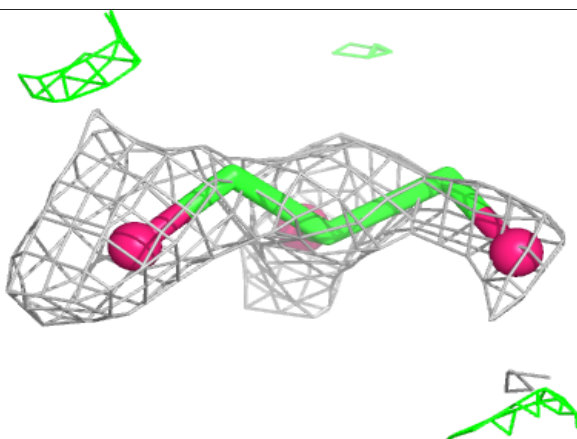
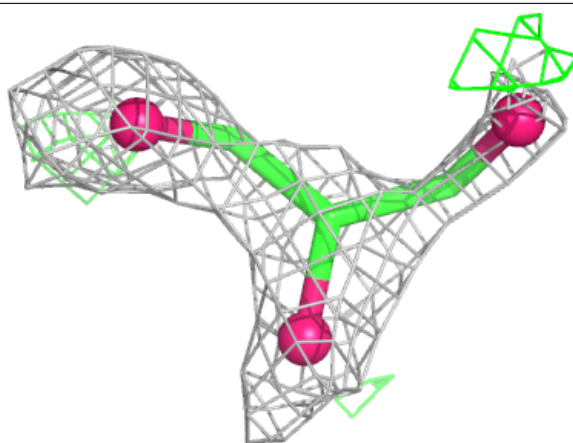
**Electron density around MG C 604:**

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



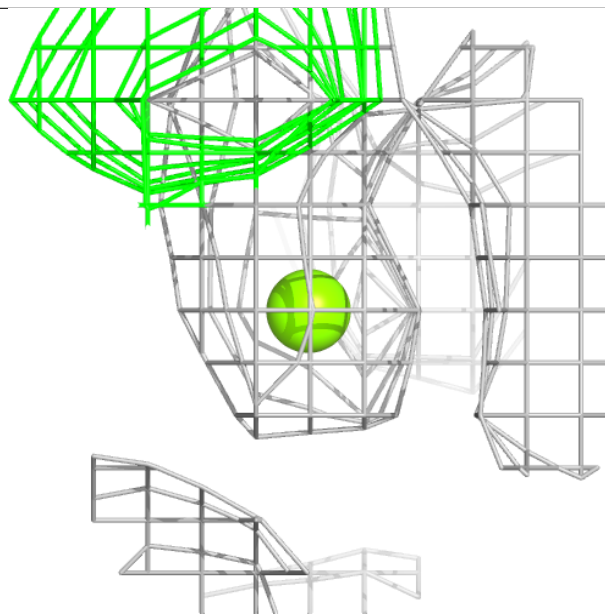
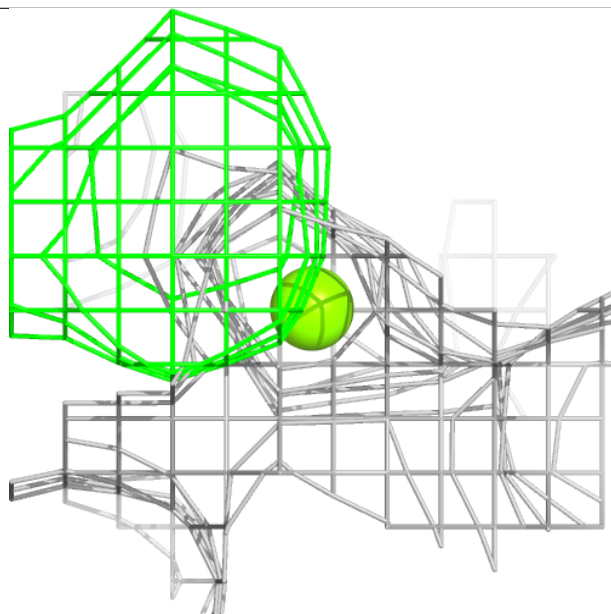
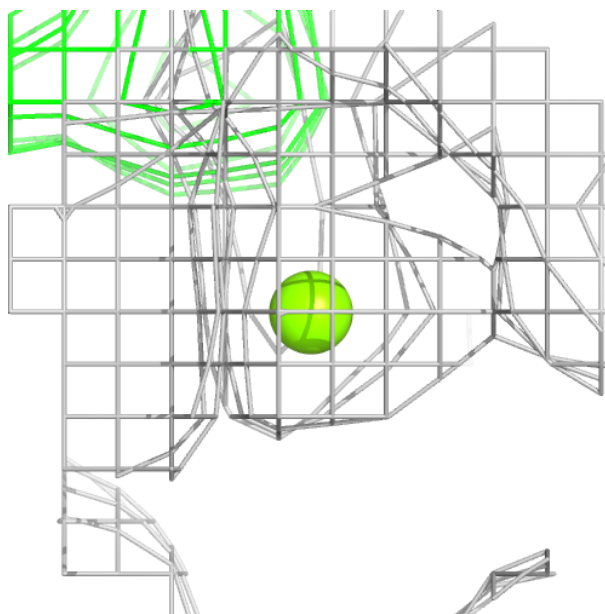
**Electron density around GOL C 619:**

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



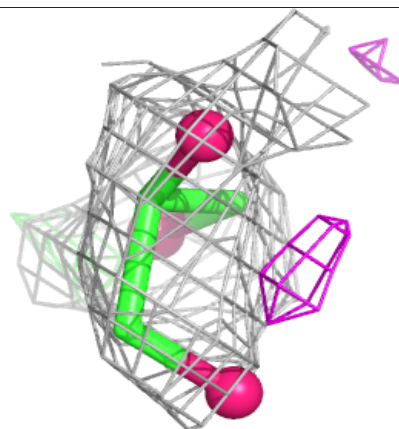
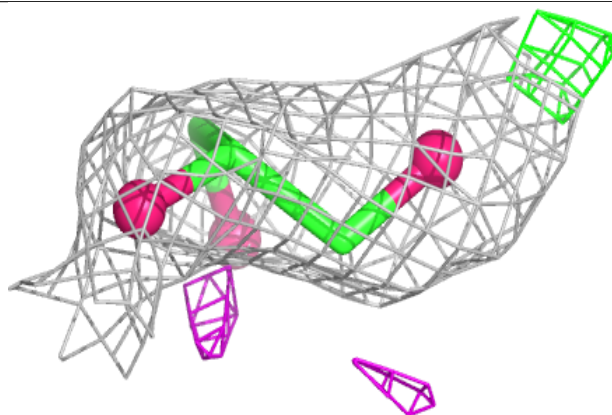
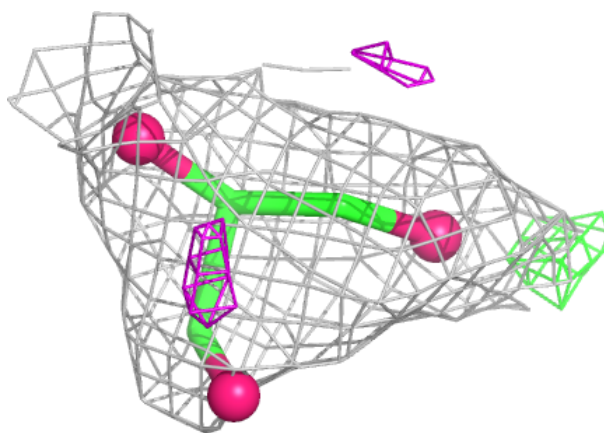
**Electron density around MG B 611:**

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



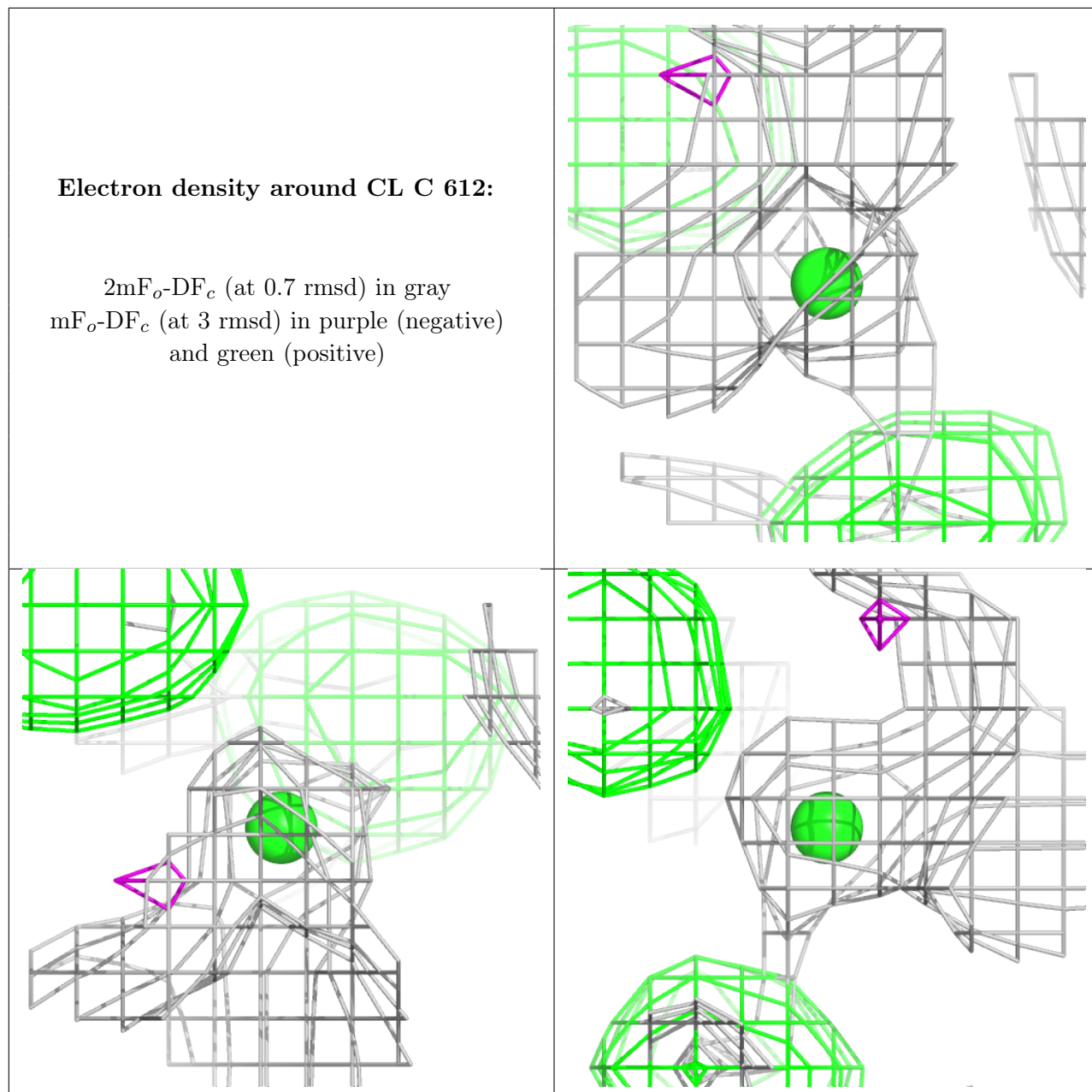
**Electron density around GOL C 617:**

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



**Electron density around CL C 612:**

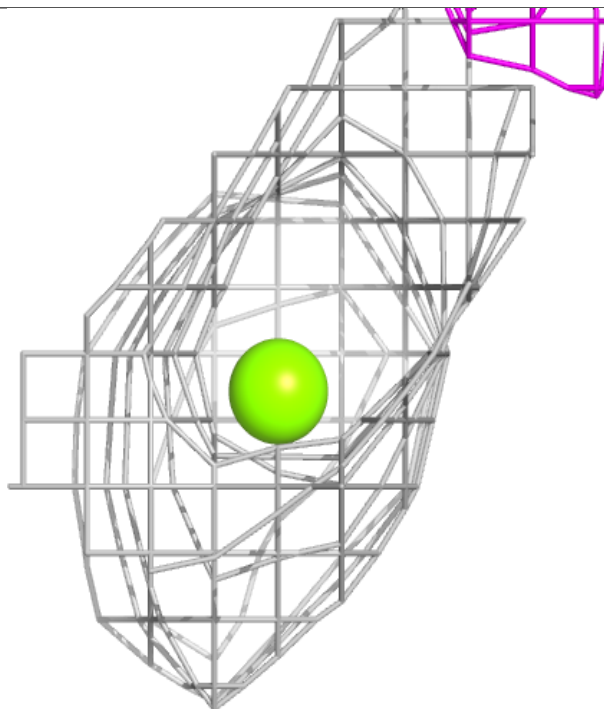
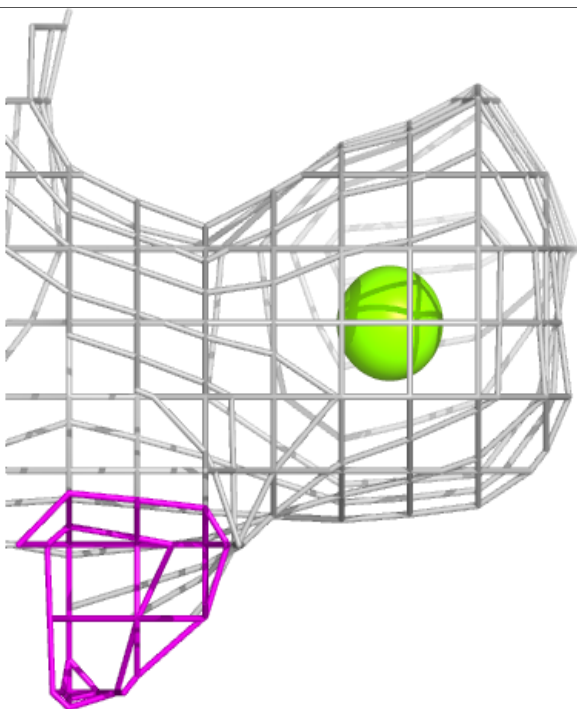
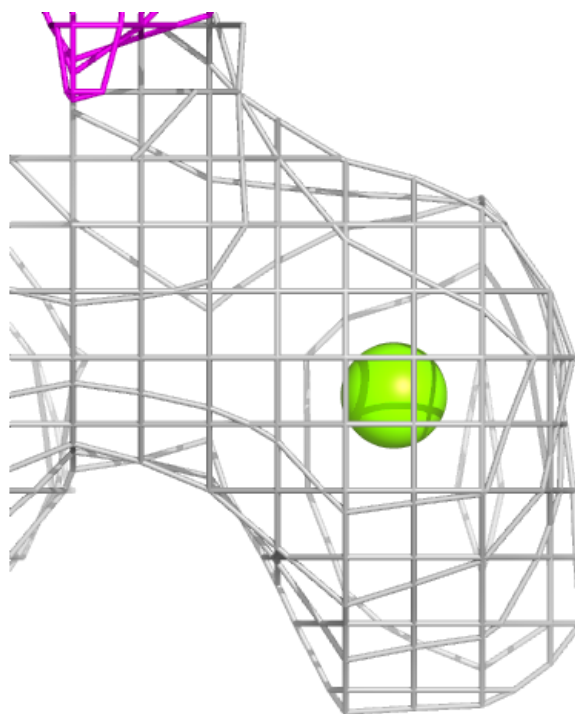
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)





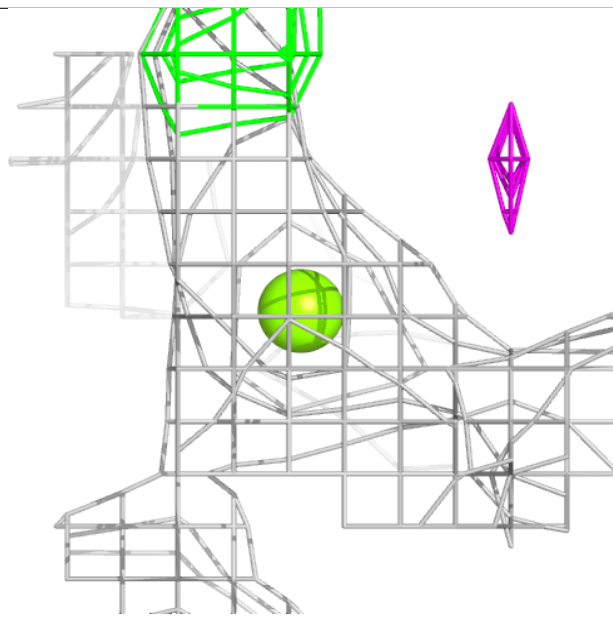
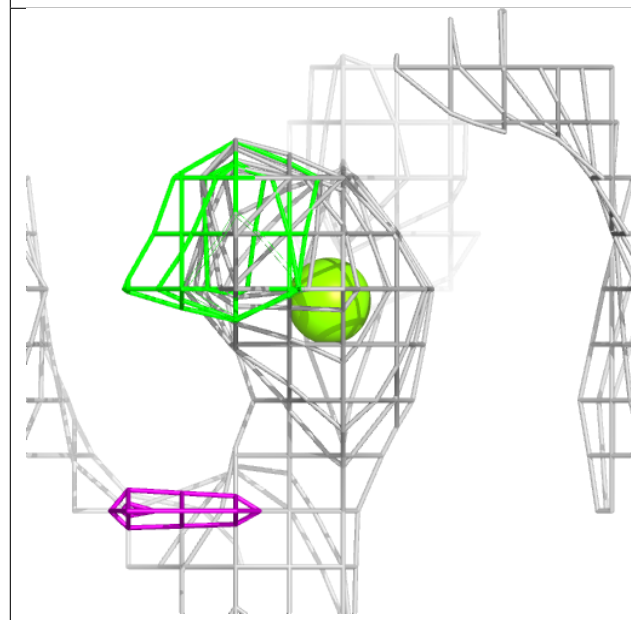
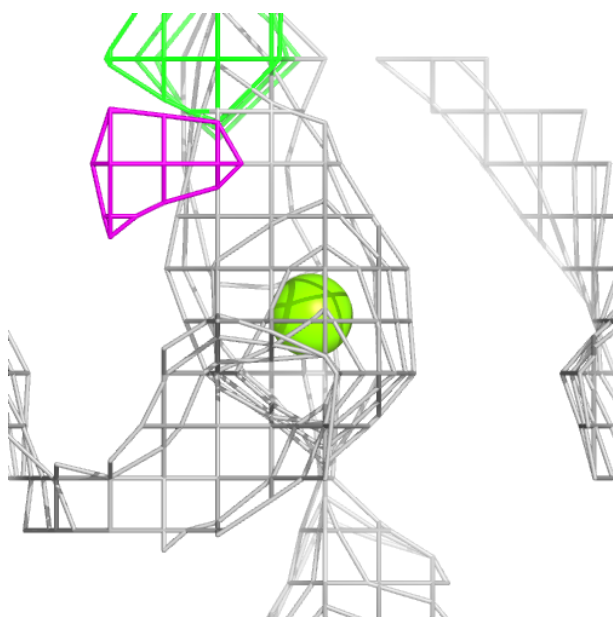
**Electron density around MG A 608:**

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



**Electron density around MG A 602:**

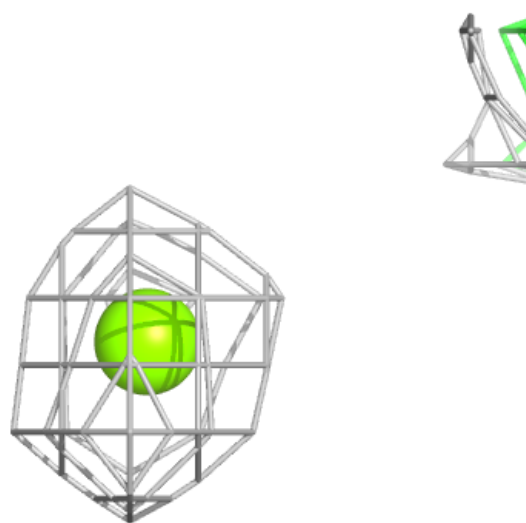
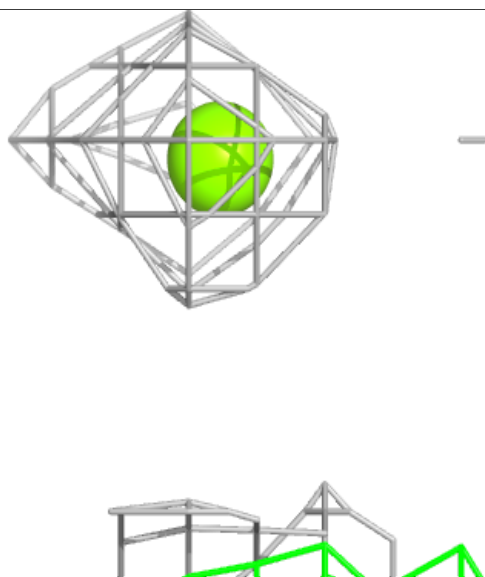
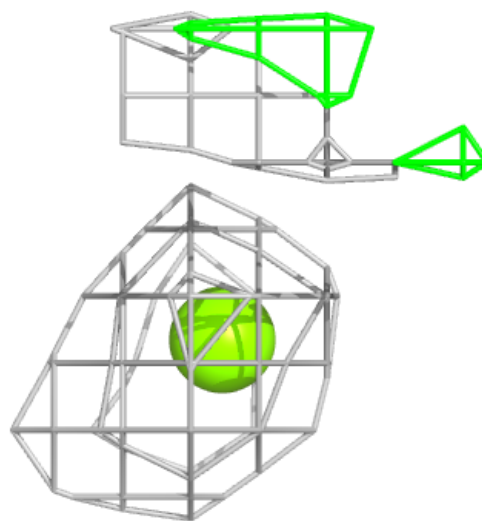
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)





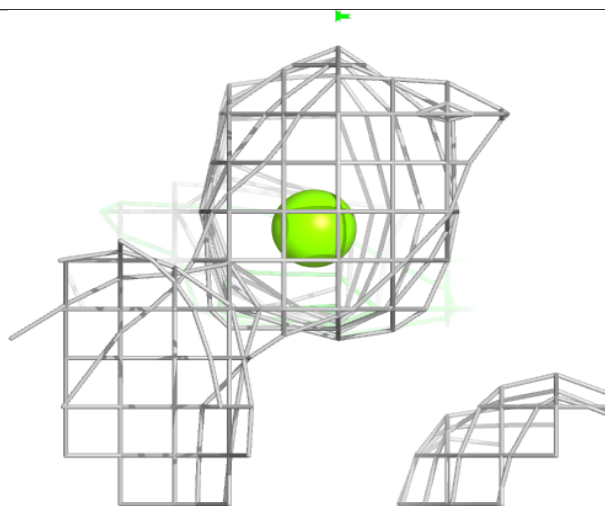
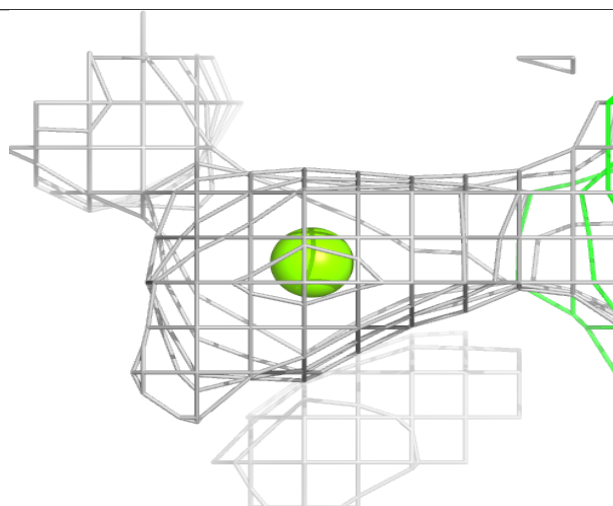
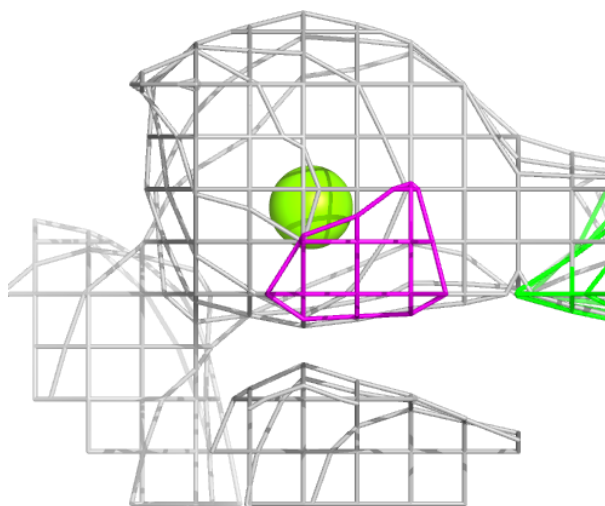
**Electron density around MG B 607:**

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



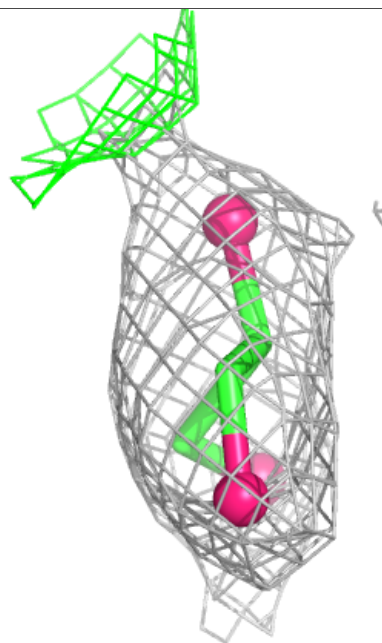
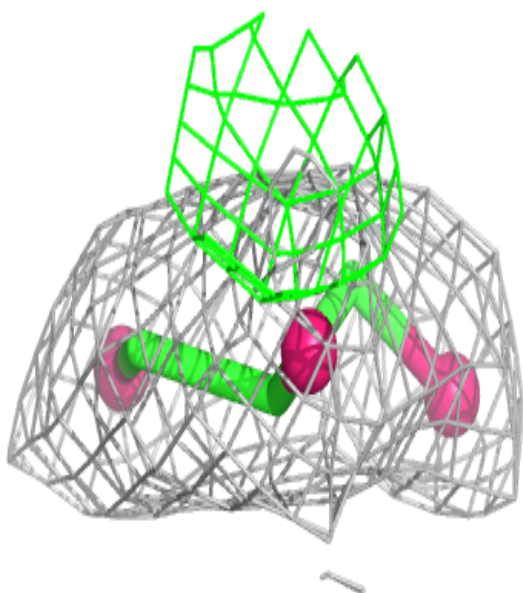
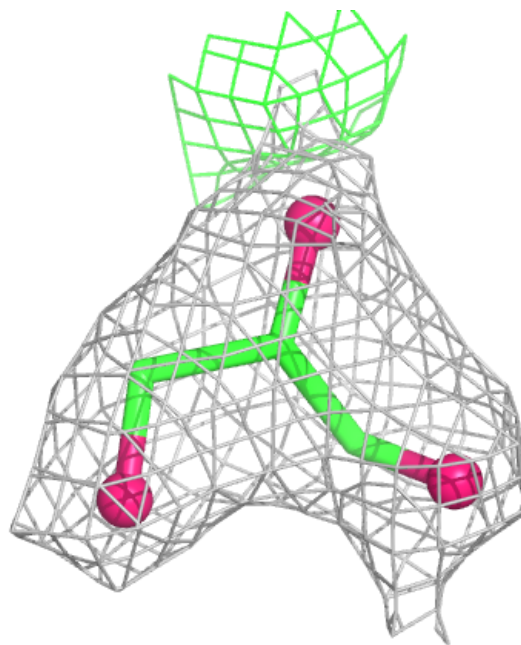
**Electron density around MG A 601:**

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



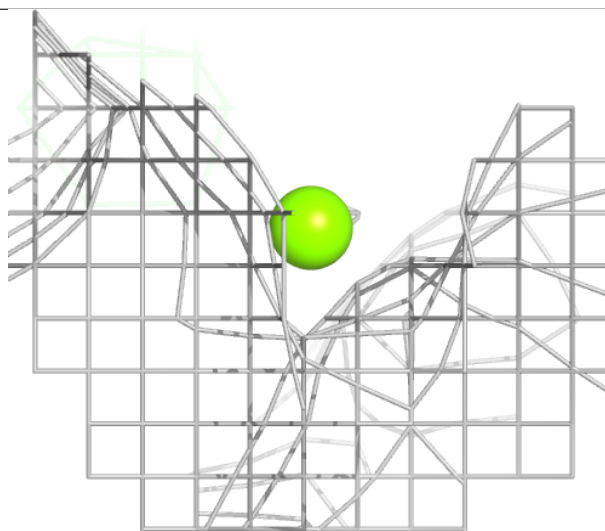
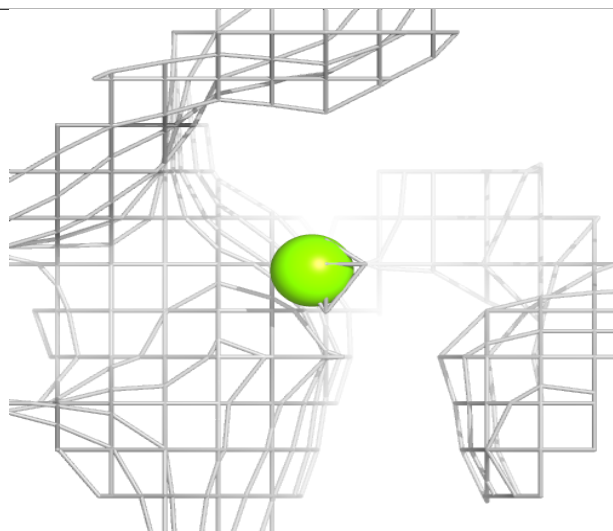
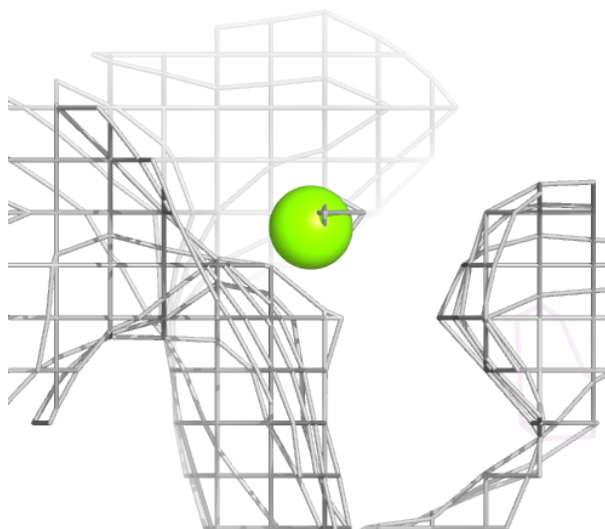
**Electron density around GOL A 609:**

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



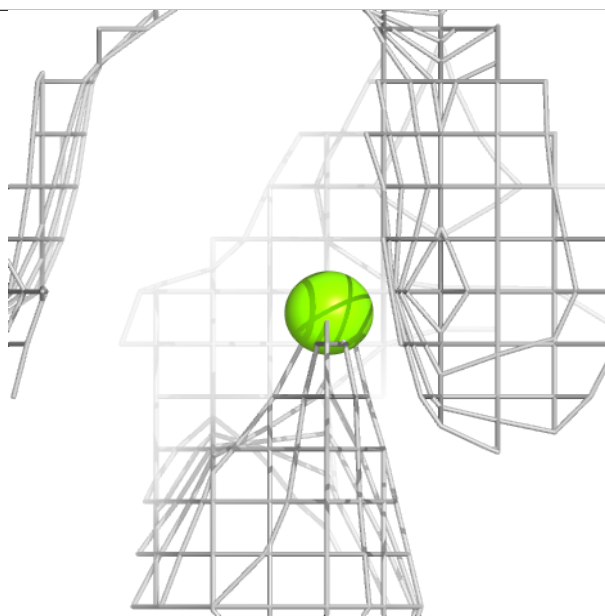
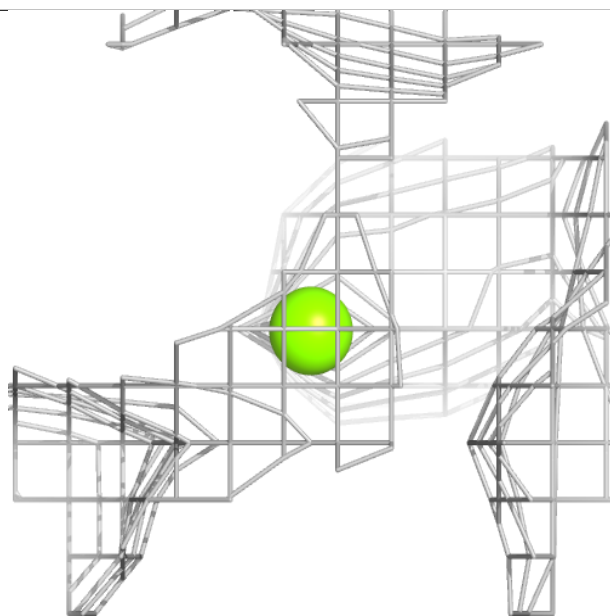
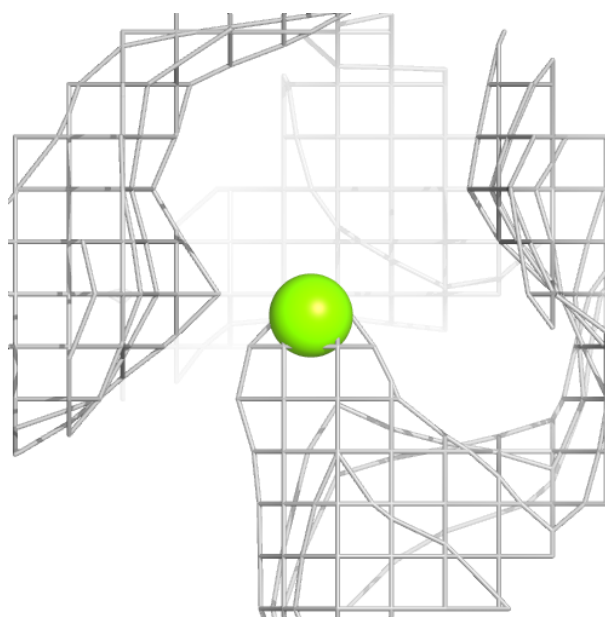
**Electron density around MG B 601:**

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



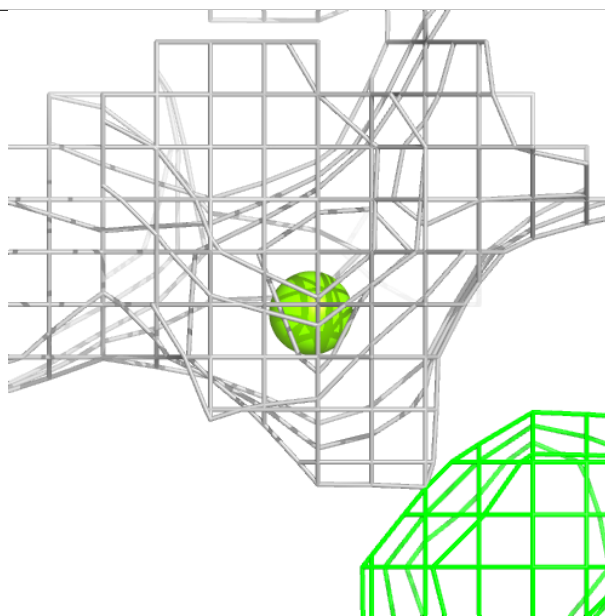
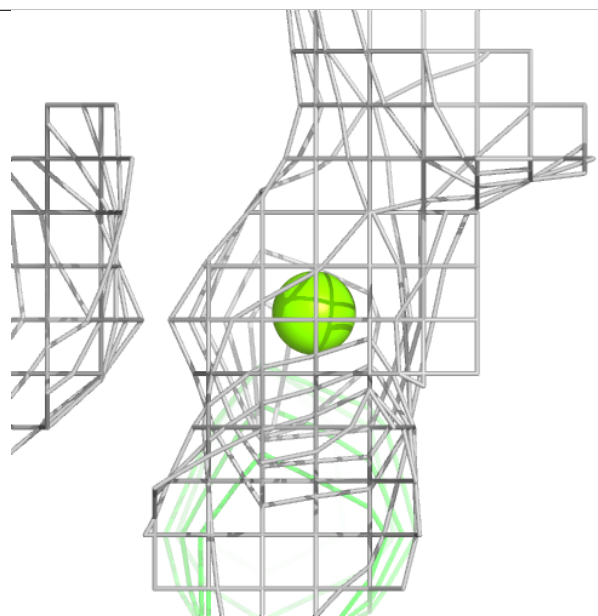
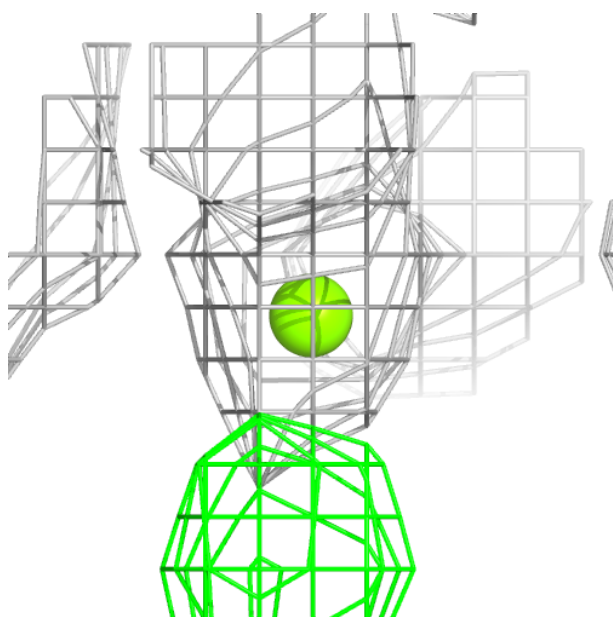
**Electron density around MG B 604:**

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



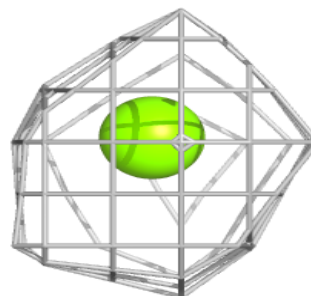
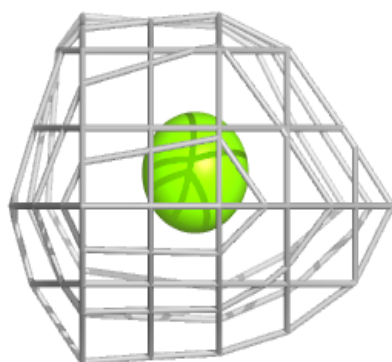
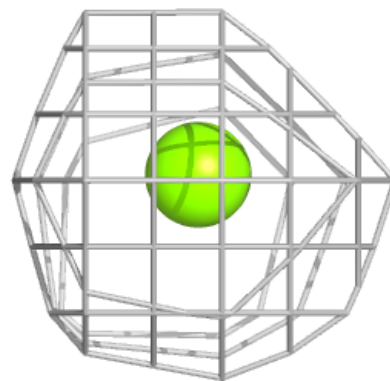
**Electron density around MG D 602:**

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



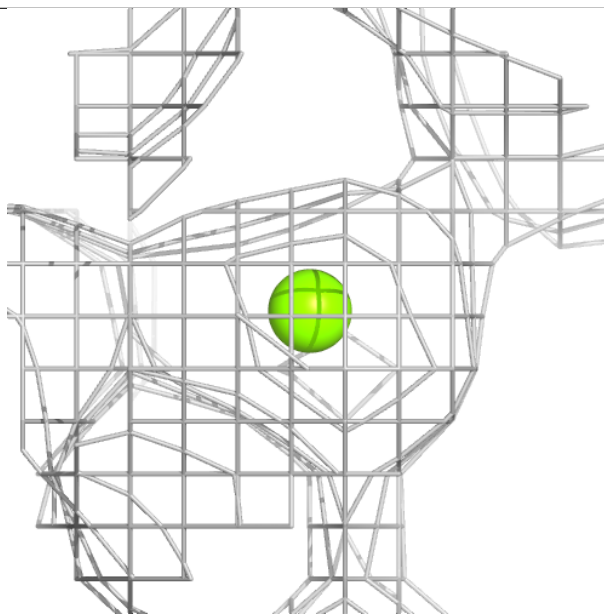
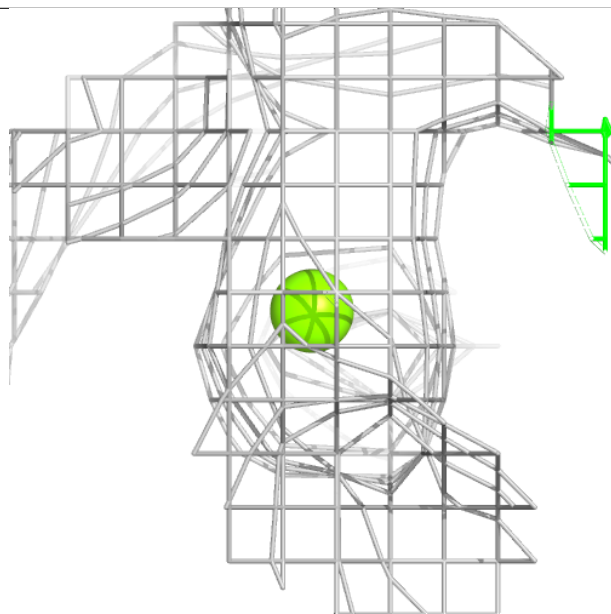
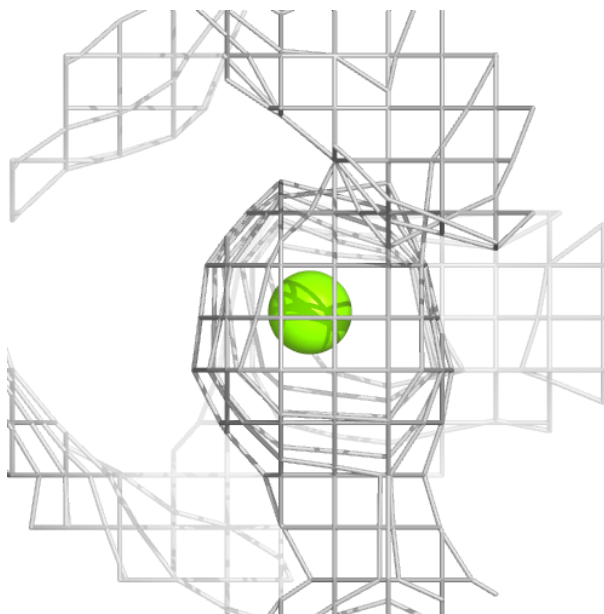
**Electron density around MG C 605:**

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



**Electron density around MG A 603:**

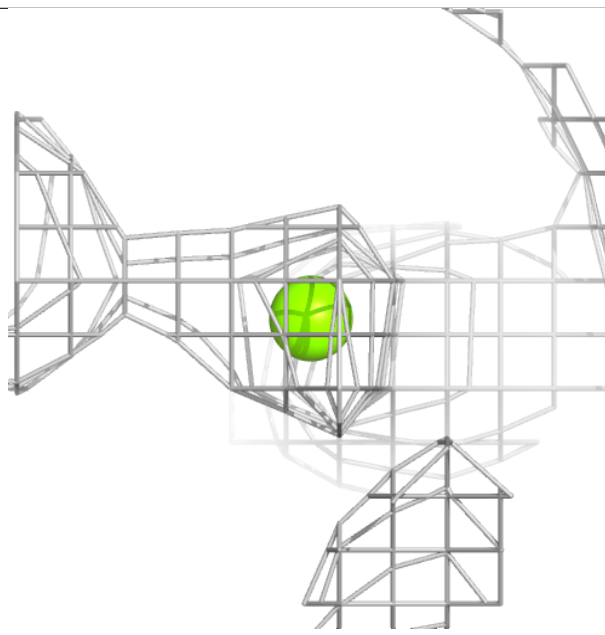
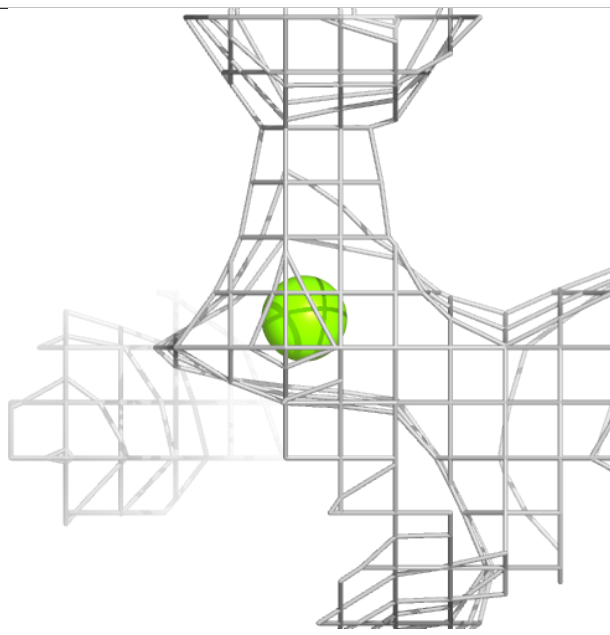
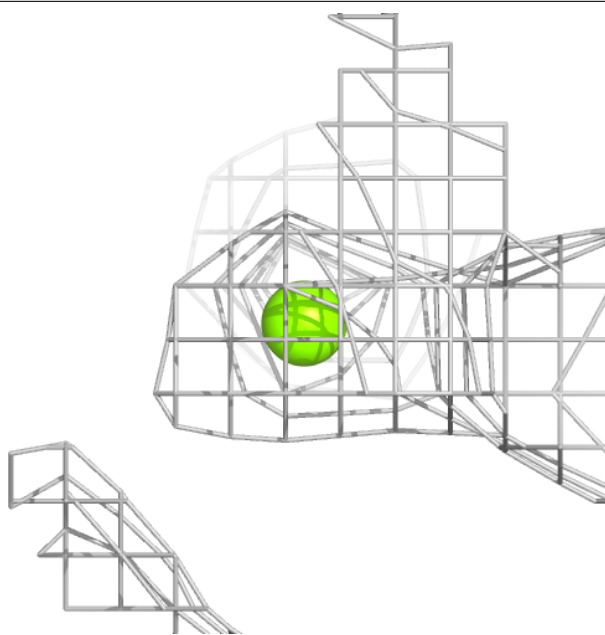
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)





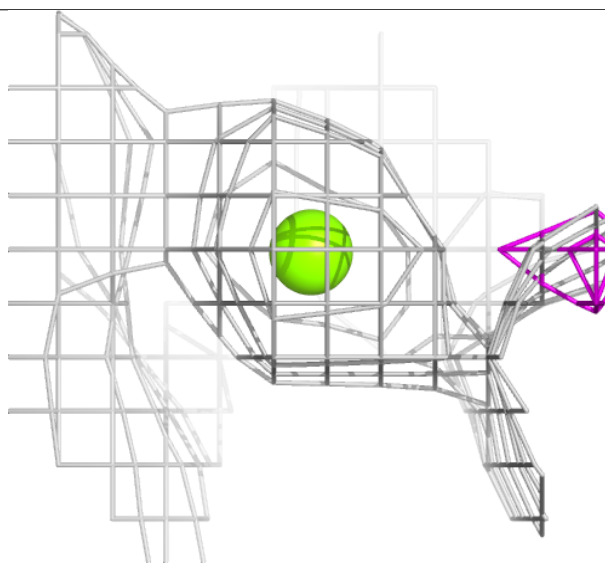
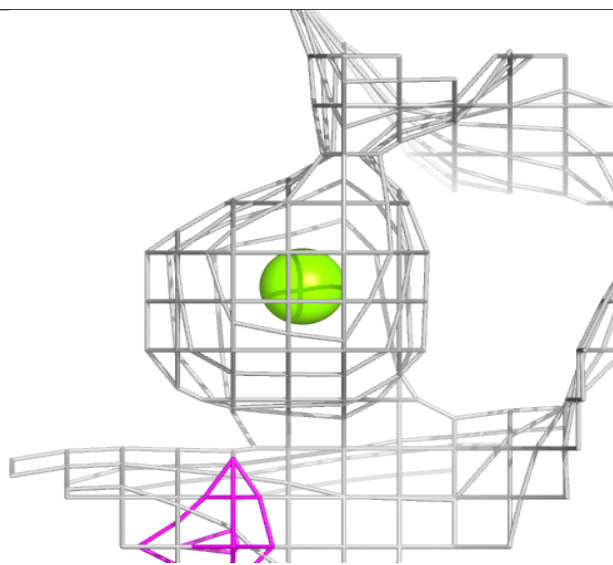
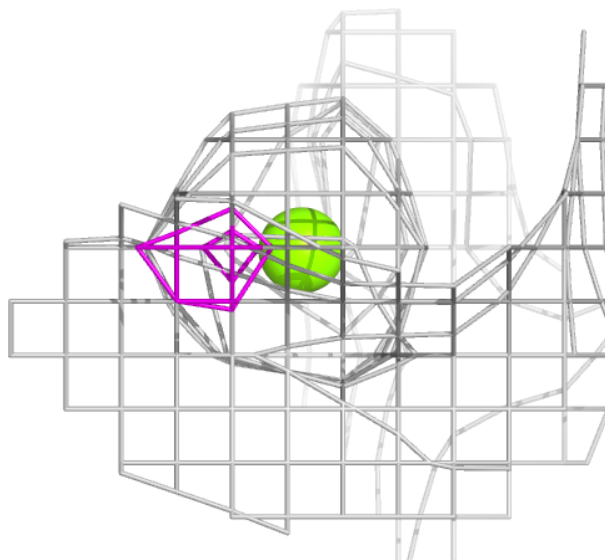
**Electron density around MG B 605:**

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



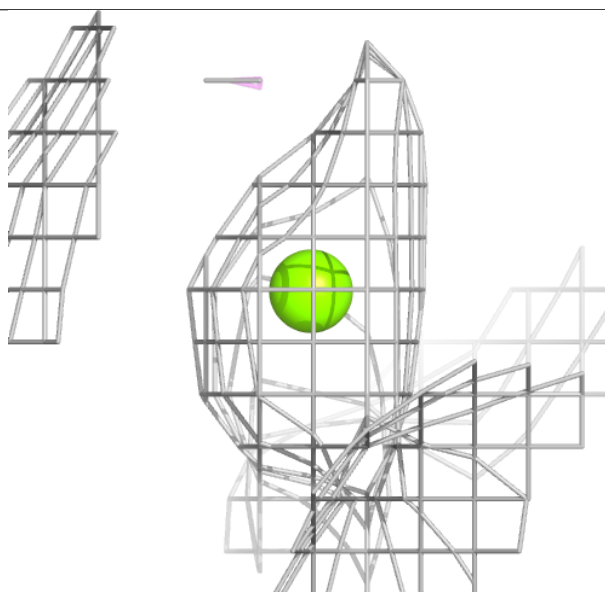
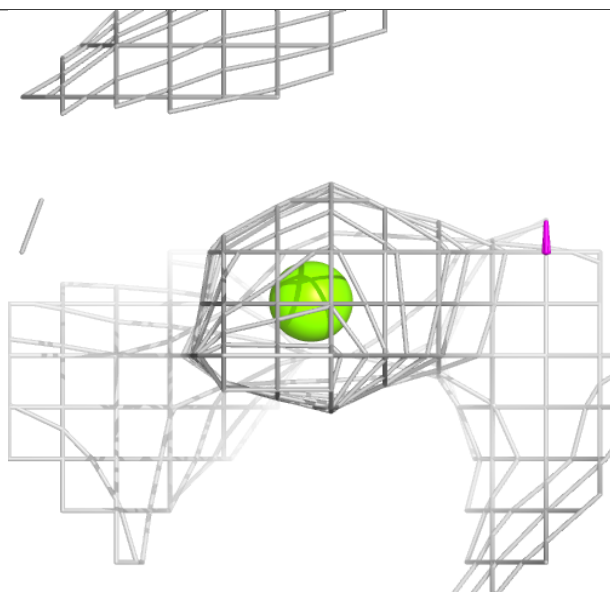
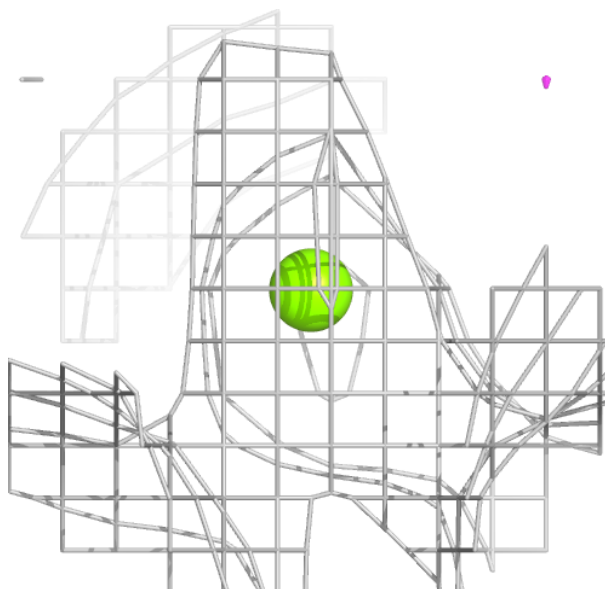
**Electron density around MG C 607:**

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



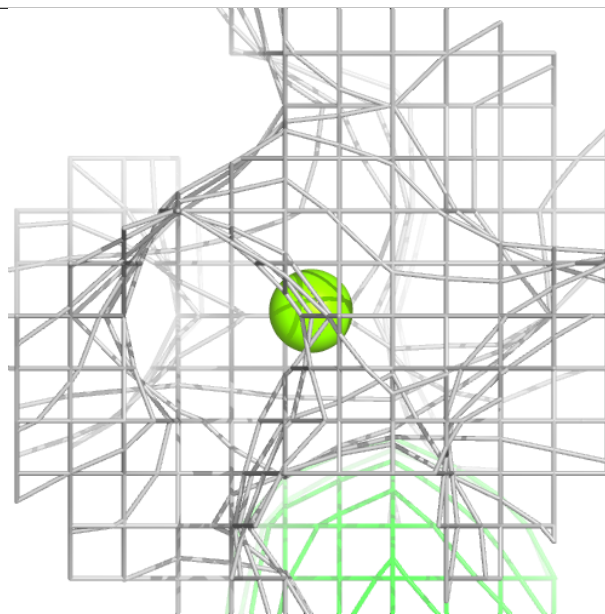
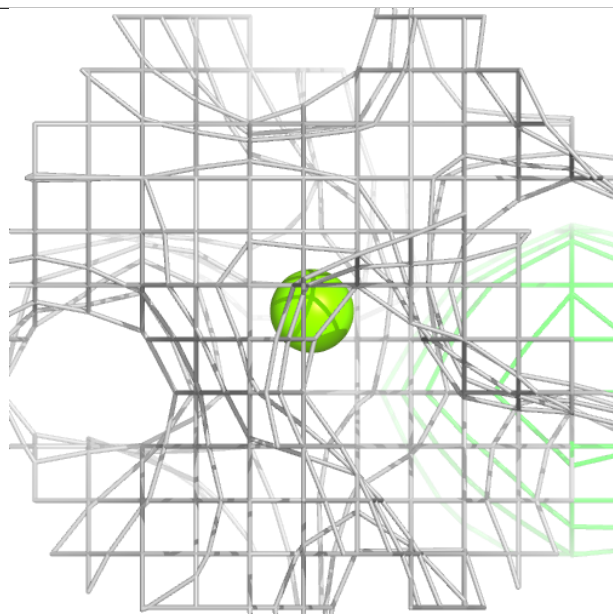
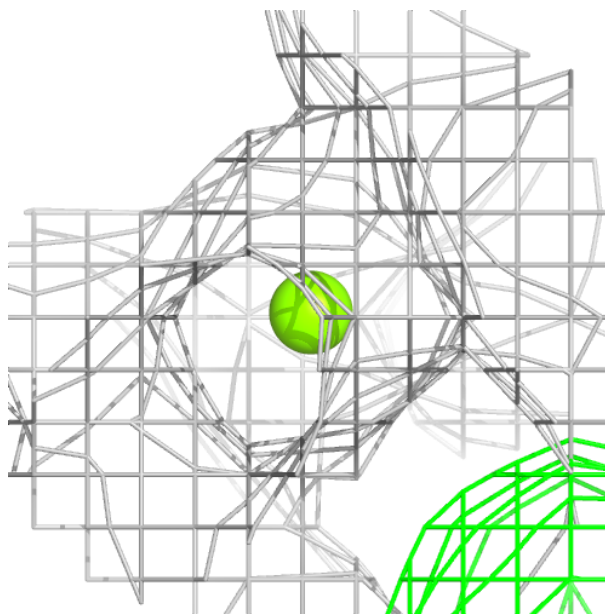
**Electron density around MG B 602:**

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



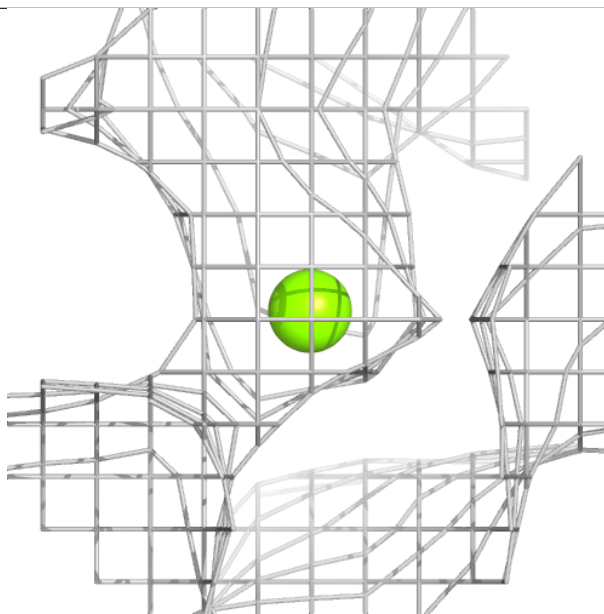
**Electron density around MG B 608:**

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



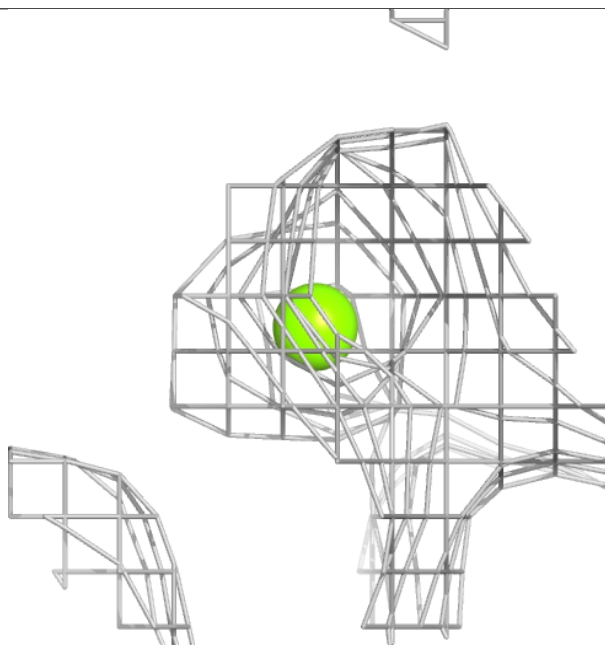
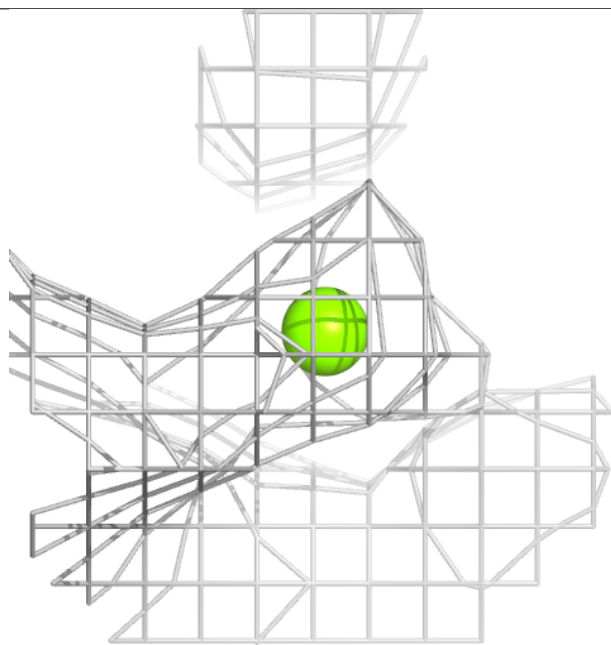
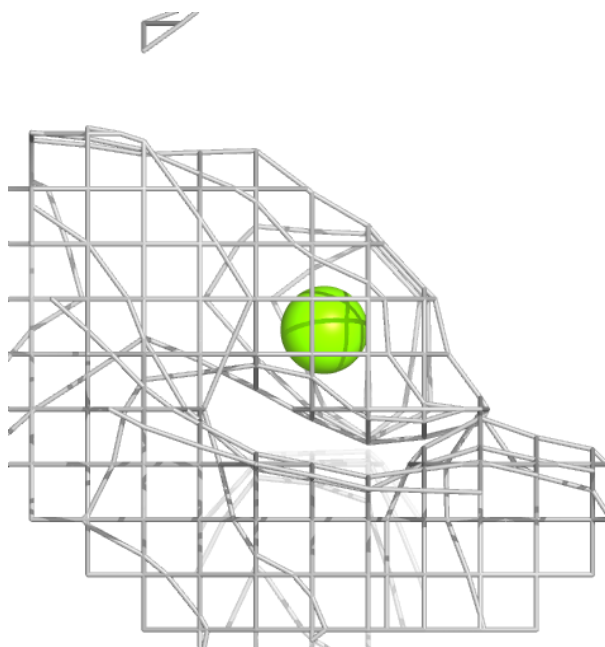
**Electron density around MG D 601:**

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



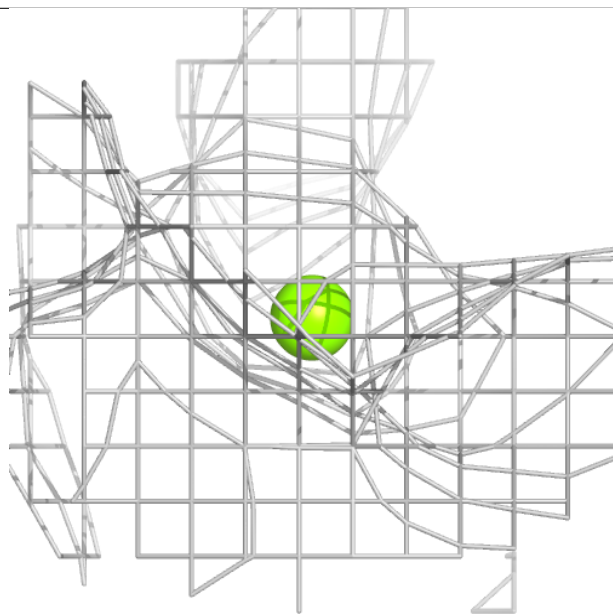
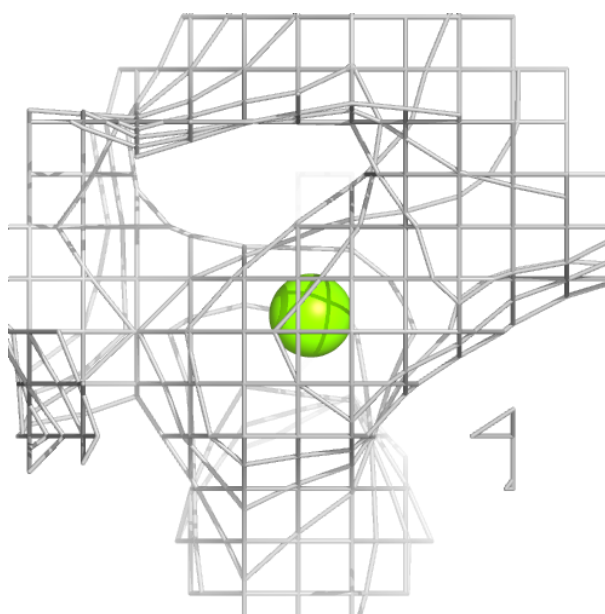
**Electron density around MG B 603:**

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



**Electron density around MG C 601:**

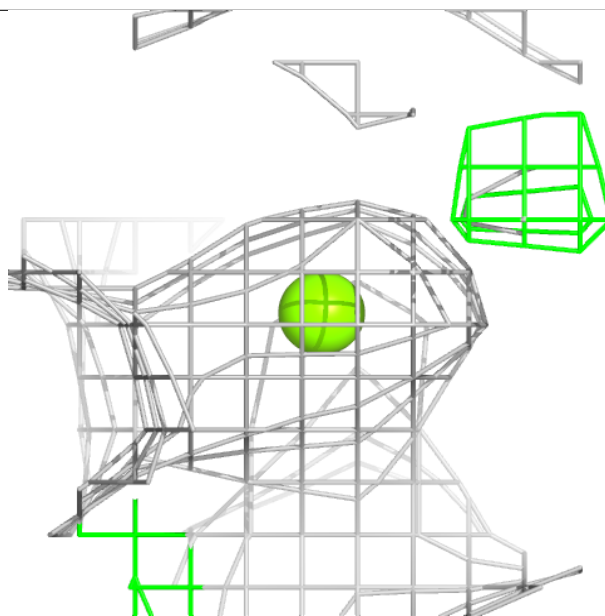
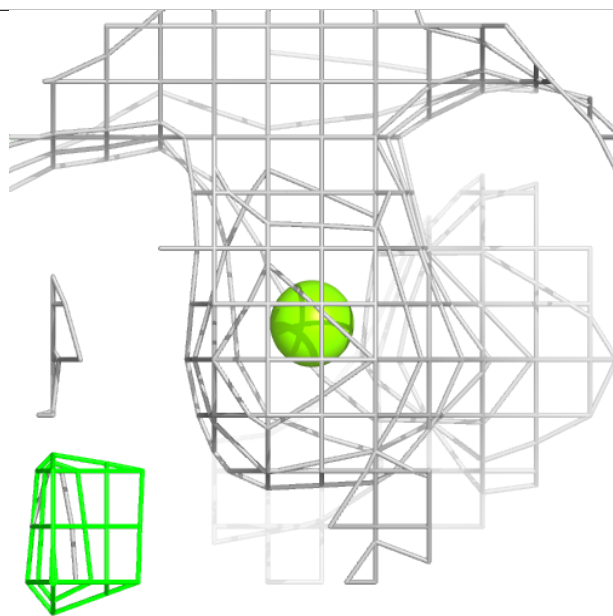
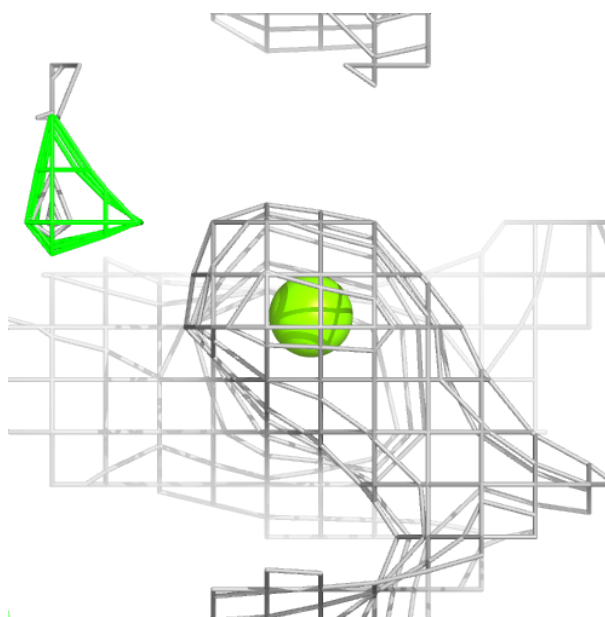
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)





**Electron density around MG C 609:**

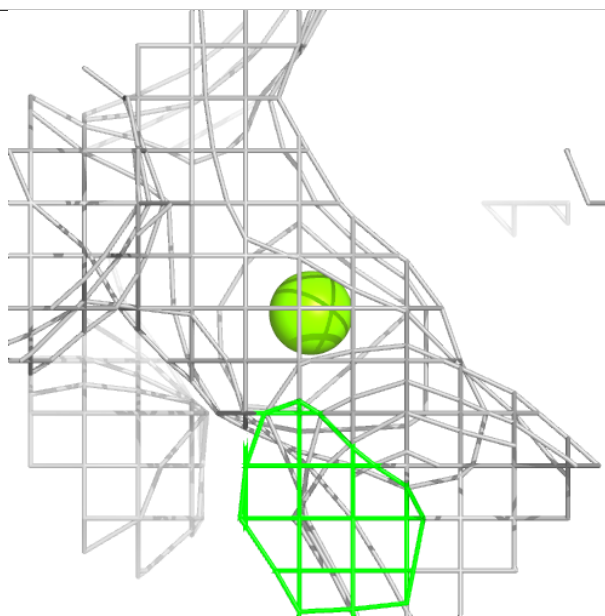
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)





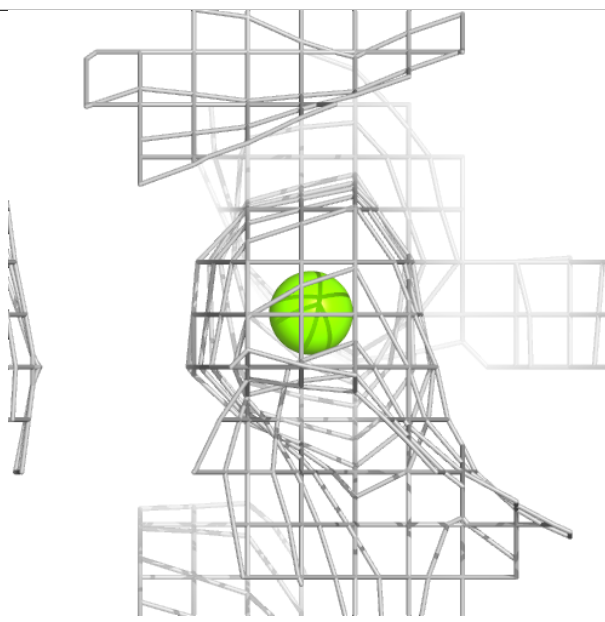
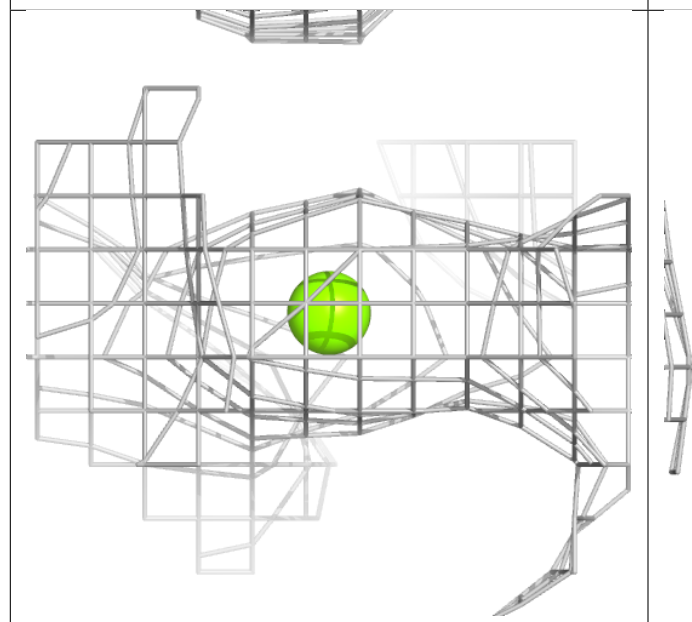
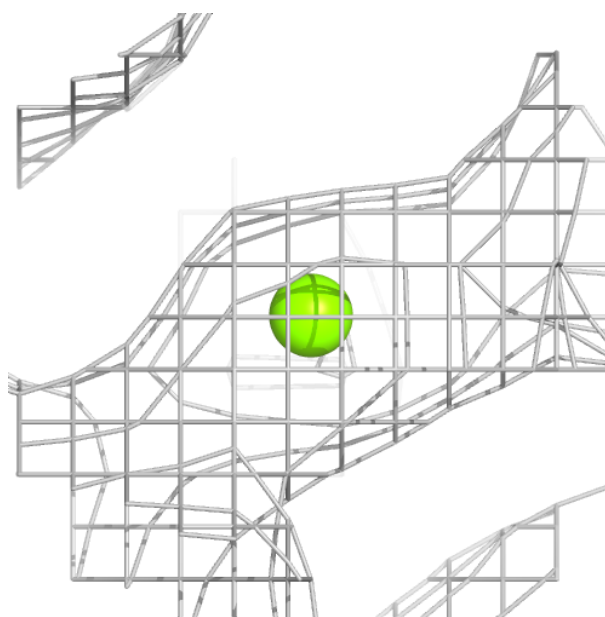
**Electron density around MG C 610:**

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



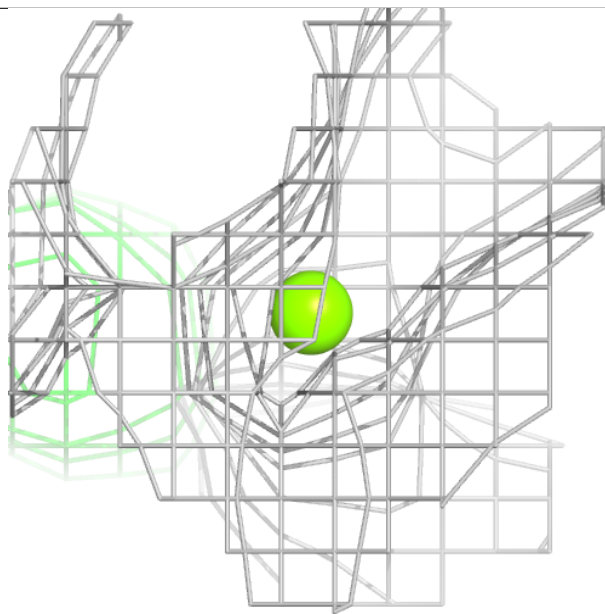
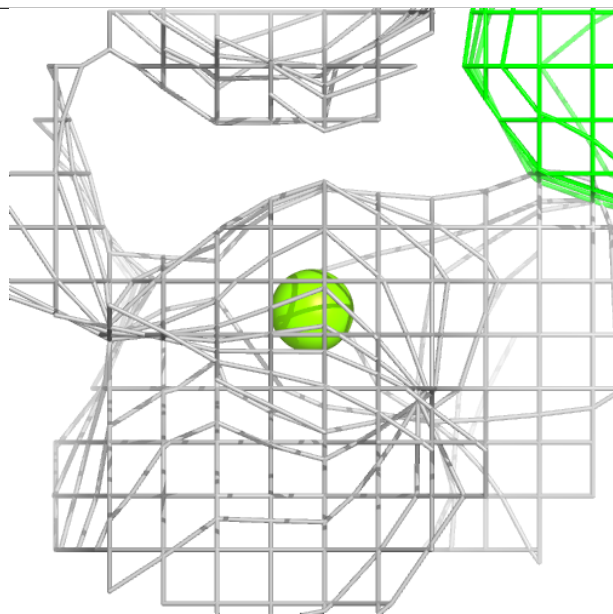
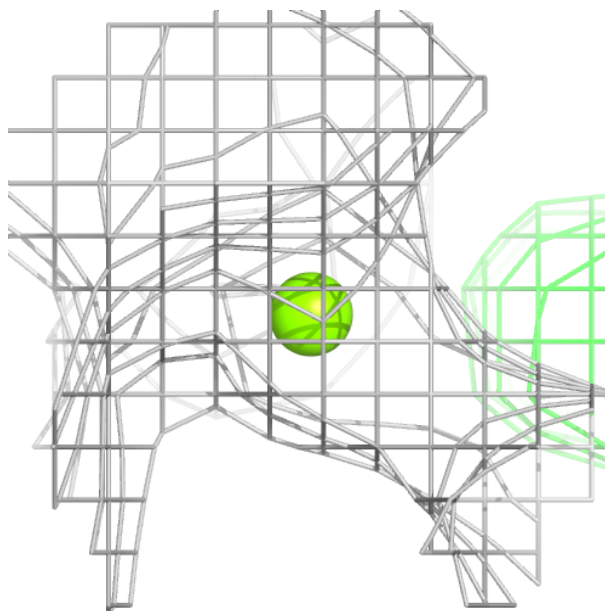
**Electron density around MG A 605:**

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



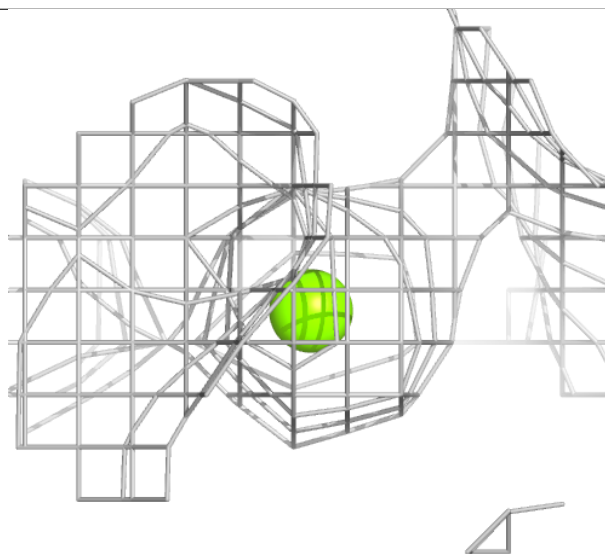
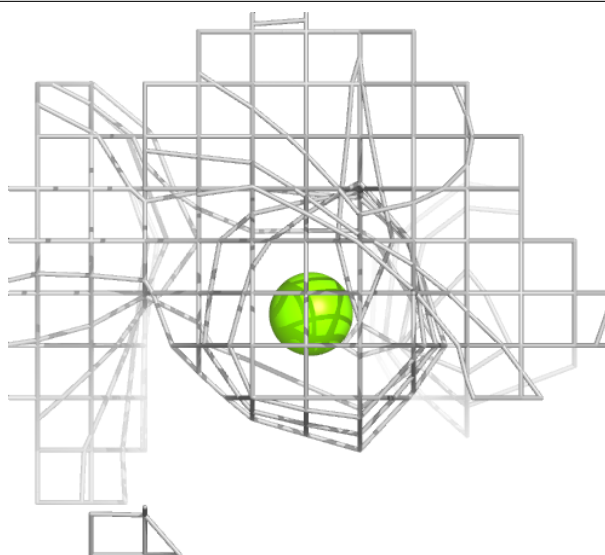
**Electron density around MG B 606:**

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



**Electron density around MG C 602:**

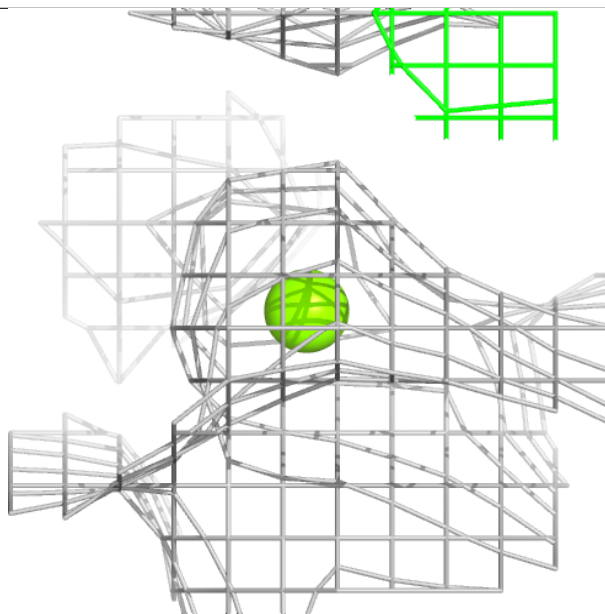
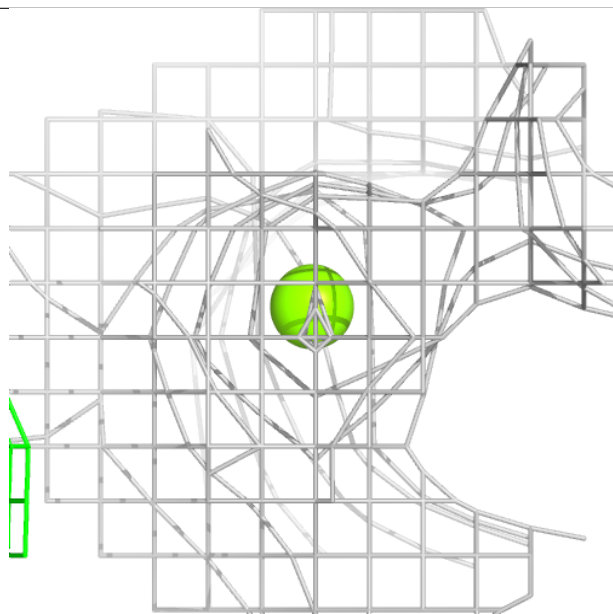
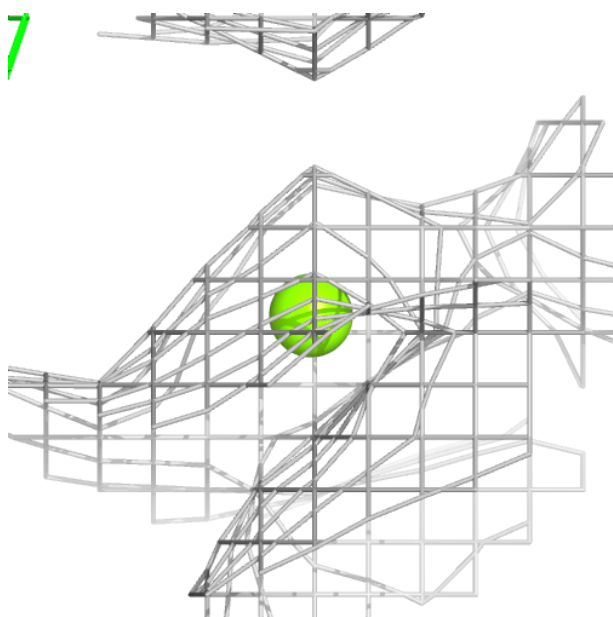
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)



**Electron density around MG C 608:**

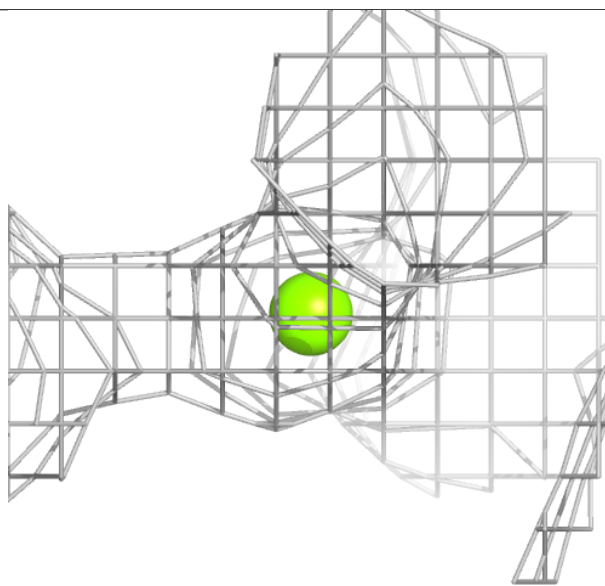
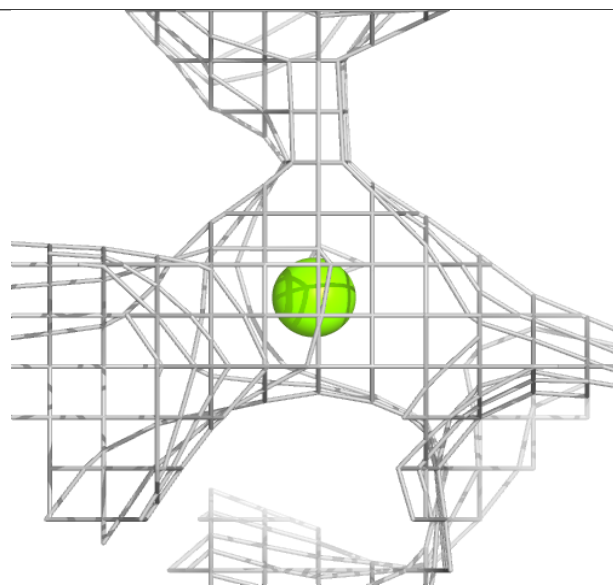
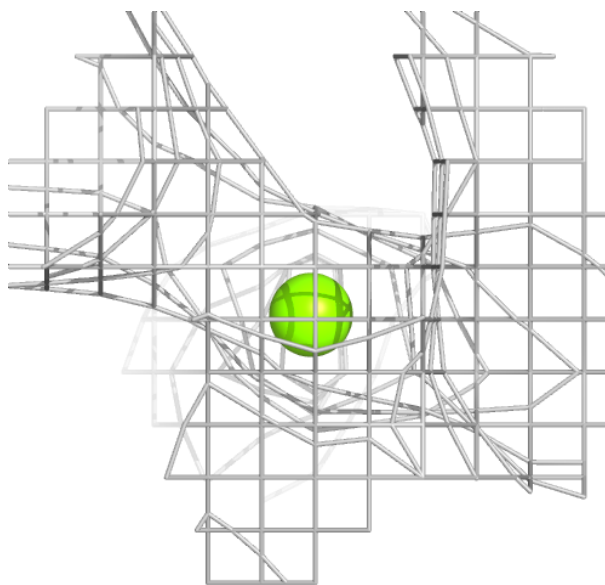
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

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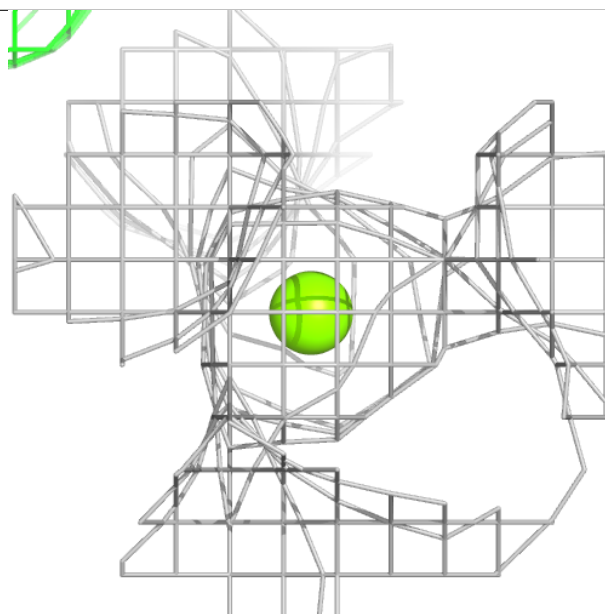
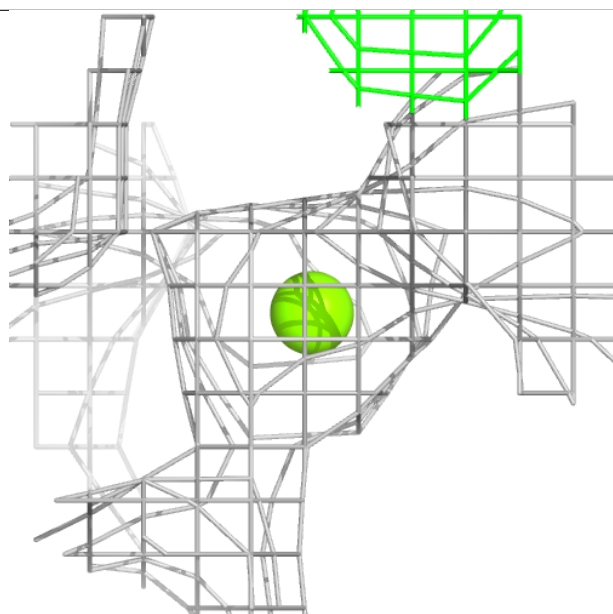
**Electron density around MG C 603:**

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



**Electron density around MG C 606:**

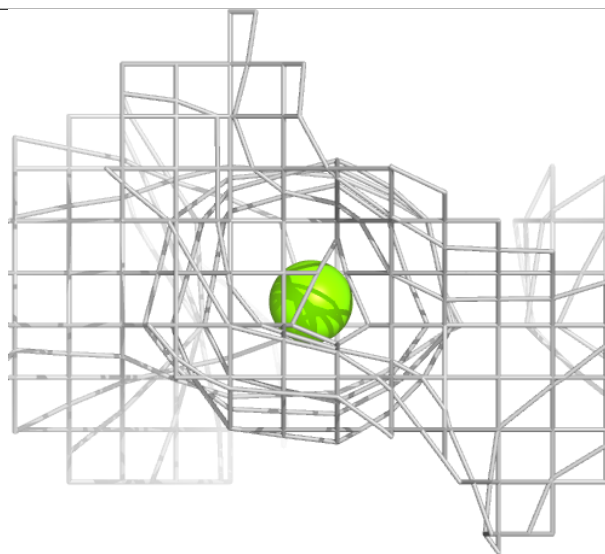
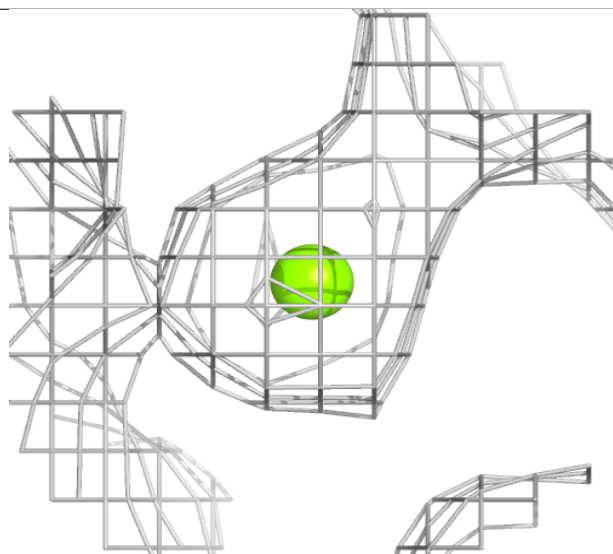
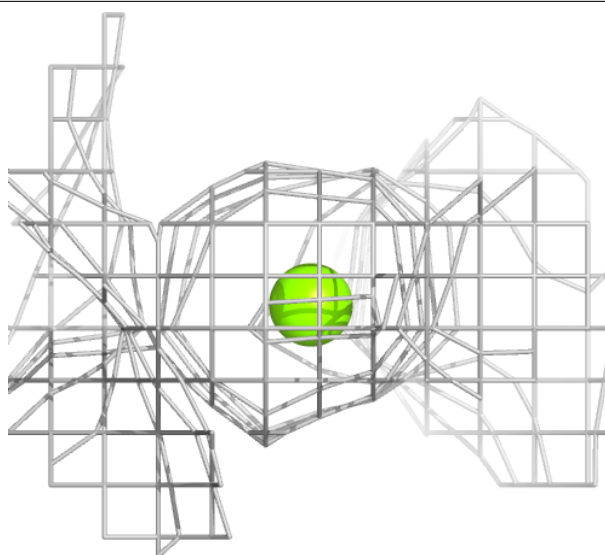
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)





**Electron density around MG A 606:**

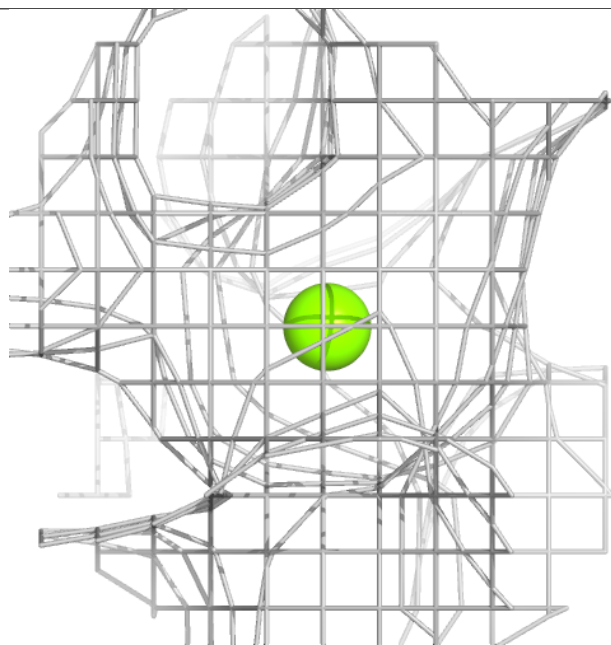
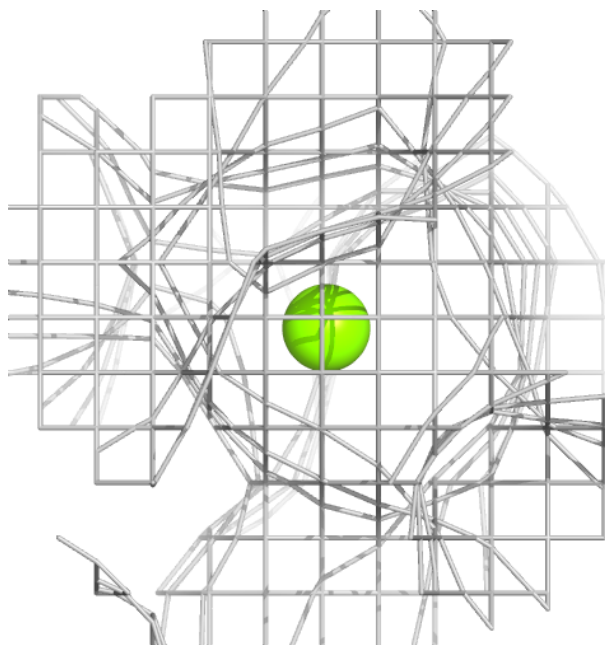
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)





**Electron density around MG A 604:**

$2mF_o-DF_c$  (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.