



wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 1, 2026 – 02:53 AM UTC

PDB ID : 7DLK / pdb_00007dlk
Title : Crystal Structure of veratryl alcohol bound Dye Decolorizing peroxidase from *Bacillus subtilis*
Authors : Dhankhar, P.; Dalal, V.; Kumar, P.
Deposited on : 2020-11-27
Resolution : 2.10 Å(reported)

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

We welcome your comments at validation@mail.wwpdb.org

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	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
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

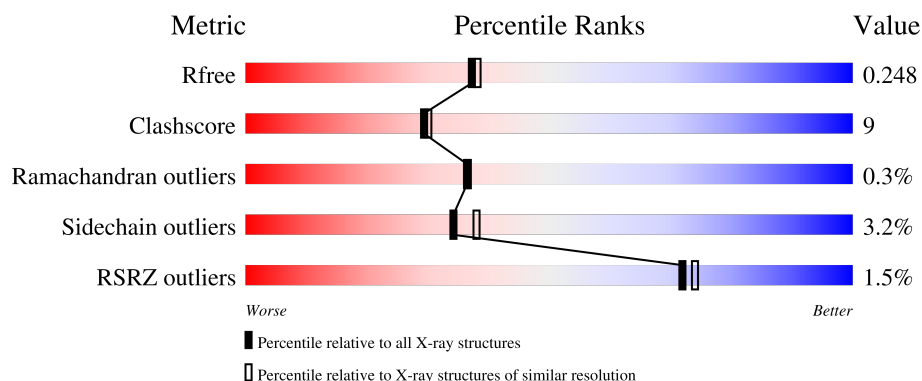
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.10 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	6658 (2.10-2.10)
Clashscore	190562	7164 (2.10-2.10)
Ramachandran outliers	187476	7099 (2.10-2.10)
Sidechain outliers	187428	7100 (2.10-2.10)
RSRZ outliers	180081	6662 (2.10-2.10)

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 ≥ 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	363	0% 83% 15% .
1	B	363	0% 86% 13% ..
1	C	363	2% 82% 16% .
1	D	363	0% 89% 10% ..
1	E	363	0% 81% 18% .

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Mol	Chain	Length	Quality of chain
1	F	363	

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
4	MPD	C	404	-	-	X	-
5	GOL	F	407	-	-	X	-
6	CL	F	408	-	-	X	-
8	VOH	C	403	-	-	X	-
8	VOH	E	403	-	X	X	-

2 Entry composition

There are 10 unique types of molecules in this entry. The entry contains 19220 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 Deferrochelataase/peroxidase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	363	Total	C	N	O	S	0	0	0
			2813	1773	481	545	14			
1	B	361	Total	C	N	O	S	0	0	0
			2803	1767	479	543	14			
1	C	363	Total	C	N	O	S	0	0	0
			2817	1775	481	547	14			
1	D	361	Total	C	N	O	S	0	1	0
			2809	1772	480	542	15			
1	E	362	Total	C	N	O	S	0	2	0
			2825	1780	483	547	15			
1	F	360	Total	C	N	O	S	0	2	0
			2814	1775	481	544	14			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MET	-	initiating methionine	UNP A0A6G2J275
A	93	SER	GLY	conflict	UNP A0A6G2J275
B	1	MET	-	initiating methionine	UNP A0A6G2J275
B	93	SER	GLY	conflict	UNP A0A6G2J275
C	1	MET	-	initiating methionine	UNP A0A6G2J275
C	93	SER	GLY	conflict	UNP A0A6G2J275
D	1	MET	-	initiating methionine	UNP A0A6G2J275
D	93	SER	GLY	conflict	UNP A0A6G2J275
E	1	MET	-	initiating methionine	UNP A0A6G2J275
E	93	SER	GLY	conflict	UNP A0A6G2J275
F	1	MET	-	initiating methionine	UNP A0A6G2J275
F	93	SER	GLY	conflict	UNP A0A6G2J275

- Molecule 2 is PROTOPORPHYRIN IX CONTAINING FE (CCD ID: HEM) (formula: $C_{34}H_{32}FeN_4O_4$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total 43	C 34	Fe 1	N 4	O 4	0	0
2	B	1	Total 43	C 34	Fe 1	N 4	O 4	0	0
2	C	1	Total 43	C 34	Fe 1	N 4	O 4	0	0
2	D	1	Total 43	C 34	Fe 1	N 4	O 4	0	0
2	E	1	Total 43	C 34	Fe 1	N 4	O 4	0	0
2	F	1	Total 43	C 34	Fe 1	N 4	O 4	0	0

- Molecule 3 is OXYGEN MOLECULE (CCD ID: OXY) (formula: O₂) (labeled as "Ligand of Interest" by depositor).



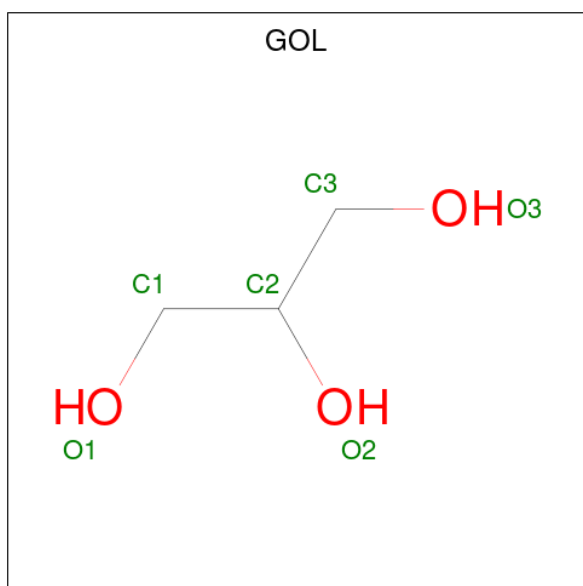
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total O 2 2	0	0
3	B	1	Total O 2 2	0	0
3	C	1	Total O 2 2	0	0
3	D	1	Total O 2 2	0	0
3	E	1	Total O 2 2	0	0
3	F	1	Total O 2 2	0	0

- Molecule 4 is (4S)-2-METHYL-2,4-PENTANEDIOL (CCD ID: MPD) (formula: C₆H₁₄O₂) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			8	6	2		
4	C	1	Total	C	O	0	0
			8	6	2		

- Molecule 5 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$) (labeled as "Ligand of Interest" by depositor).



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

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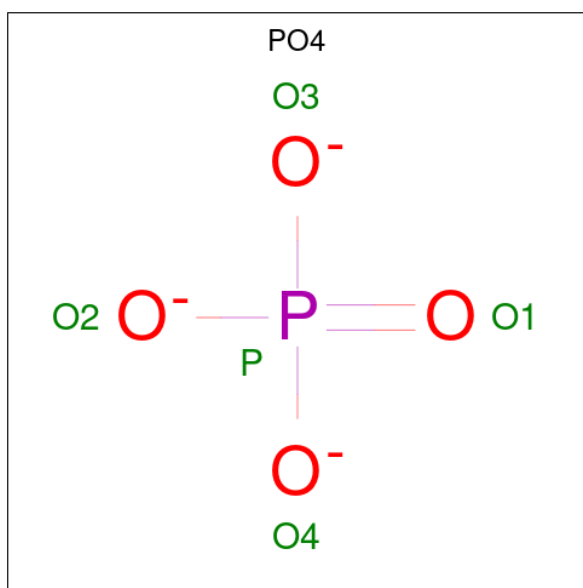
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	1	Total C O 6 3 3	0	0
5	A	1	Total C O 6 3 3	0	0
5	B	1	Total C O 6 3 3	0	0
5	C	1	Total C O 6 3 3	0	0
5	C	1	Total C O 6 3 3	0	0
5	D	1	Total C O 6 3 3	0	0
5	E	1	Total C O 6 3 3	0	0
5	E	1	Total C O 6 3 3	0	0
5	F	1	Total C O 6 3 3	0	0
5	F	1	Total C O 6 3 3	0	0
5	F	1	Total C O 6 3 3	0	0
5	F	1	Total C O 6 3 3	0	0

- Molecule 6 is CHLORIDE ION (CCD ID: CL) (formula: Cl) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	A	1	Total Cl 1 1	0	0
6	B	1	Total Cl 1 1	0	0
6	C	1	Total Cl 1 1	0	0
6	D	2	Total Cl 2 2	0	0
6	E	1	Total Cl 1 1	0	0
6	F	2	Total Cl 2 2	0	0

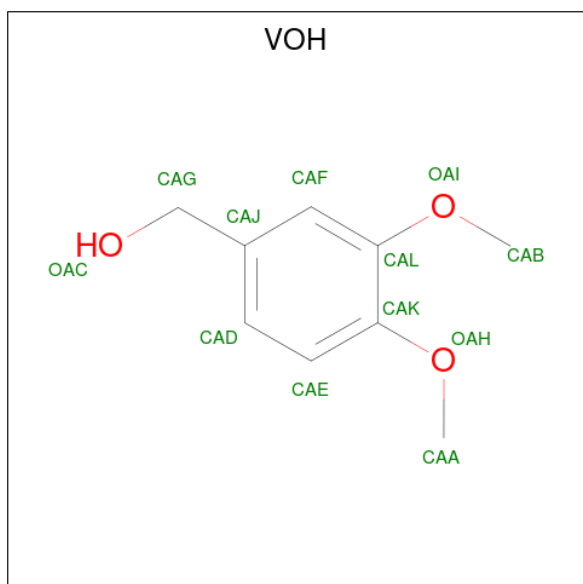
- Molecule 7 is PHOSPHATE ION (CCD ID: PO4) (formula: O₄P) (labeled as "Ligand of

Interest" by depositor).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	A	1	Total	O	P	0	0
			5	4	1		

- Molecule 8 is Veratryl alcohol (CCD ID: VOH) (formula: $C_9H_{12}O_3$) (labeled as "Ligand of Interest" by depositor).



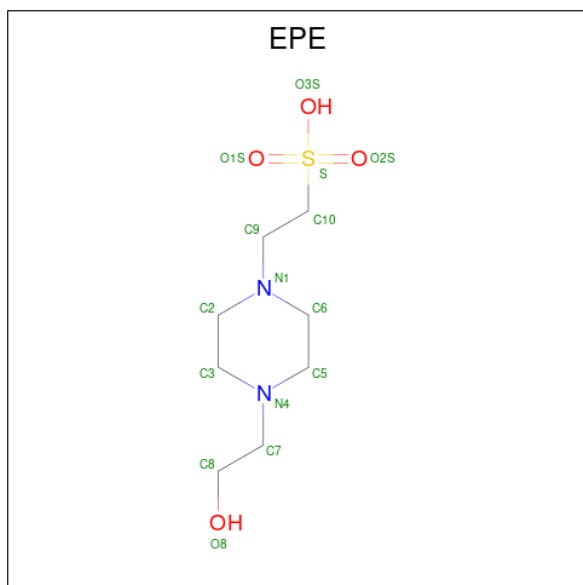
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
8	C	1	Total	C	O	0	0
			12	9	3		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
8	E	1	Total	C	O	0	0
			12	9	3		

- Molecule 9 is 4-(2-HYDROXYETHYL)-1-PIPERAZINE ETHANESULFONIC ACID (CCD ID: EPE) (formula: $C_8H_{18}N_2O_4S$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
9	F	1	Total	C	N	O	S	0	0
			15	8	2	4	1		

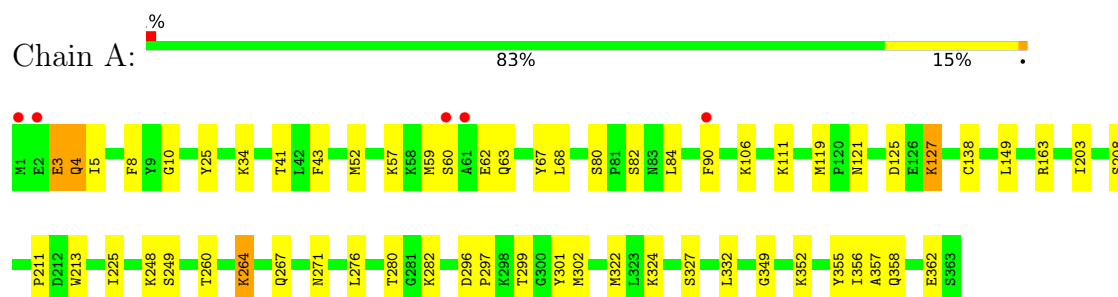
- Molecule 10 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
10	A	320	Total	O	0	0
			320	320		
10	B	334	Total	O	0	0
			334	334		
10	C	347	Total	O	0	0
			347	347		
10	D	353	Total	O	0	0
			353	353		
10	E	283	Total	O	0	0
			283	283		
10	F	286	Total	O	0	0
			286	286		

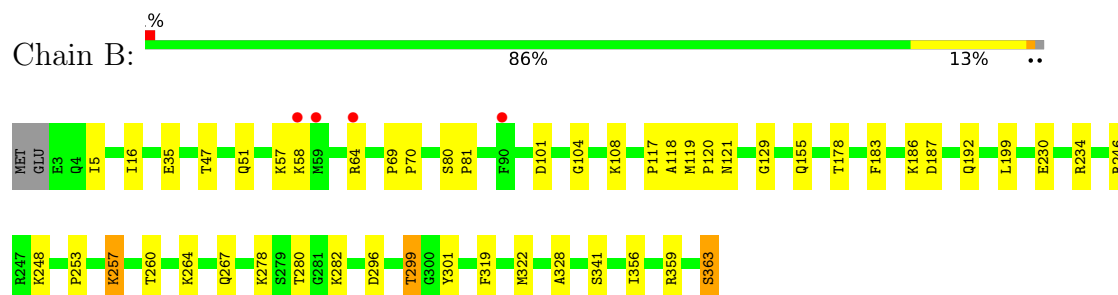
3 Residue-property plots

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.

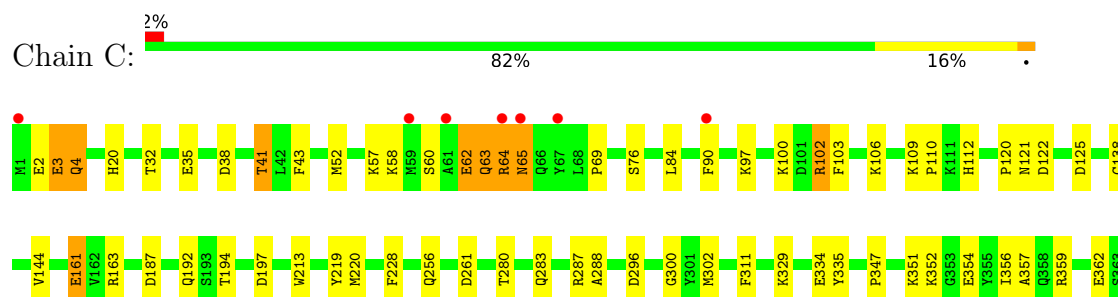
- Molecule 1: Deferrochelataase/peroxidase



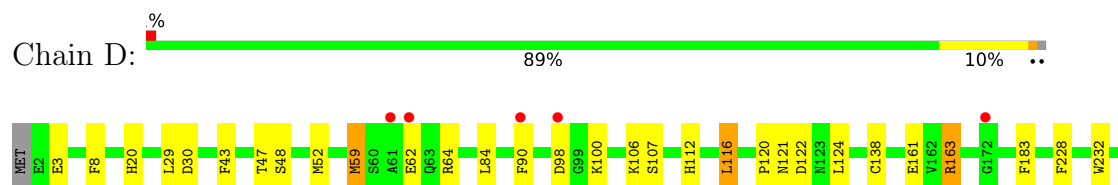
- Molecule 1: Deferrochelataase/peroxidase



- Molecule 1: Deferrochelataase/peroxidase

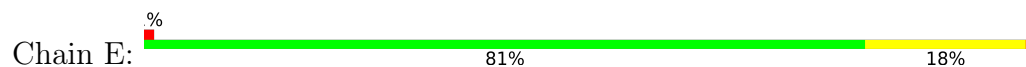


- Molecule 1: Deferrochelataase/peroxidase

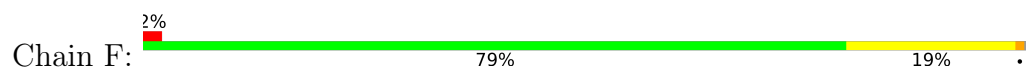




• Molecule 1: Deferrochelataase/peroxidase



• Molecule 1: Deferrochelataase/peroxidase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	79.99Å 102.11Å 105.47Å 87.84° 76.45° 83.22°	Depositor
Resolution (Å)	45.68 – 2.10 45.68 – 2.10	Depositor EDS
% Data completeness (in resolution range)	90.6 (45.68-2.10) 90.6 (45.68-2.10)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.68 (at 2.10Å)	Xtriage
Refinement program	REFMAC 5.8.0238	Depositor
R, R_{free}	0.185 , 0.249 0.186 , 0.248	Depositor DCC
R_{free} test set	8575 reflections (4.56%)	wwPDB-VP
Wilson B-factor (Å ²)	39.5	Xtriage
Anisotropy	0.204	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 47.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	19220	wwPDB-VP
Average B, all atoms (Å ²)	49.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.27% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical 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, VOH, EPE, PO4, HEM, OXY, CL, MPD

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	0.65	0/2873	1.20	5/3873 (0.1%)
1	B	0.66	0/2863	1.17	7/3859 (0.2%)
1	C	0.67	0/2877	1.18	8/3878 (0.2%)
1	D	0.66	0/2869	1.16	1/3868 (0.0%)
1	E	0.62	0/2885	1.18	3/3888 (0.1%)
1	F	0.65	1/2874 (0.0%)	1.21	8/3874 (0.2%)
All	All	0.65	1/17241 (0.0%)	1.18	32/23240 (0.1%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	F	149	LEU	C-N	-8.17	1.23	1.33

The worst 5 of 32 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	F	149	LEU	CA-C-N	7.45	130.59	120.54
1	F	149	LEU	C-N-CA	7.45	130.59	120.54
1	C	194	THR	CA-CB-OG1	-6.82	99.38	109.60
1	C	122	ASP	CA-CB-CG	6.76	119.36	112.60
1	B	69	PRO	CB-CA-C	-6.44	103.06	110.92

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	2813	0	2753	48	0
1	B	2803	0	2746	38	0
1	C	2817	0	2757	56	0
1	D	2809	0	2751	31	0
1	E	2825	0	2763	56	0
1	F	2814	0	2758	49	0
2	A	43	0	30	3	0
2	B	43	0	30	2	0
2	C	43	0	30	4	0
2	D	43	0	30	3	0
2	E	43	0	30	0	0
2	F	43	0	30	4	0
3	A	2	0	0	0	0
3	B	2	0	0	0	0
3	C	2	0	0	0	0
3	D	2	0	0	0	0
3	E	2	0	0	0	0
3	F	2	0	0	0	0
4	A	8	0	14	5	0
4	C	8	0	14	8	0
5	A	18	0	24	4	0
5	B	6	0	8	0	0
5	C	12	0	16	4	0
5	D	6	0	8	1	0
5	E	12	0	16	0	0
5	F	24	0	32	7	0
6	A	1	0	0	1	0
6	B	1	0	0	0	0
6	C	1	0	0	0	0
6	D	2	0	0	1	0
6	E	1	0	0	0	0
6	F	2	0	0	2	0
7	A	5	0	0	0	0
8	C	12	0	0	7	0
8	E	12	0	0	6	0
9	F	15	0	18	2	0
10	A	320	0	0	12	0
10	B	334	0	0	5	0
10	C	347	0	0	11	0
10	D	353	0	0	10	0
10	E	283	0	0	7	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
10	F	286	0	0	10	0
All	All	19220	0	16858	294	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 9.

The worst 5 of 294 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:C:102:ARG:HD3	10:C:699:HOH:O	1.45	1.15
1:F:70:PRO:HA	5:F:407:GOL:H32	1.05	1.00
1:A:119:MET:HE1	1:A:324:LYS:HA	1.39	1.00
4:A:403:MPD:H51	1:E:302:MET:H	1.25	1.00
1:F:70:PRO:CA	5:F:407:GOL:H32	1.93	0.99

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	361/363 (99%)	349 (97%)	12 (3%)	0	100	100
1	B	359/363 (99%)	351 (98%)	8 (2%)	0	100	100
1	C	361/363 (99%)	345 (96%)	12 (3%)	4 (1%)	11	8
1	D	360/363 (99%)	347 (96%)	12 (3%)	1 (0%)	36	36
1	E	362/363 (100%)	349 (96%)	12 (3%)	1 (0%)	36	36
1	F	360/363 (99%)	340 (94%)	19 (5%)	1 (0%)	36	36
All	All	2163/2178 (99%)	2081 (96%)	75 (4%)	7 (0%)	36	36

5 of 7 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	63	GLN
1	E	362	GLU
1	C	65	ASN
1	C	60	SER
1	D	62	GLU

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	302/305 (99%)	288 (95%)	14 (5%)	24	25
1	B	302/305 (99%)	297 (98%)	5 (2%)	53	62
1	C	303/305 (99%)	292 (96%)	11 (4%)	31	34
1	D	302/305 (99%)	295 (98%)	7 (2%)	44	51
1	E	304/305 (100%)	295 (97%)	9 (3%)	36	41
1	F	303/305 (99%)	291 (96%)	12 (4%)	28	29
All	All	1816/1830 (99%)	1758 (97%)	58 (3%)	34	38

5 of 58 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	352	LYS
1	F	326	LEU
1	D	161	GLU
1	F	313	LYS
1	F	109	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 29 such sidechains are listed below:

Mol	Chain	Res	Type
1	D	256	GLN
1	F	207	GLN
1	E	51	GLN
1	F	71	GLN

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Mol	Chain	Res	Type
1	E	20	HIS

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 39 ligands modelled in this entry, 8 are monoatomic - leaving 31 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	D	403	-	5,5,5	0.26	0	5,5,5	0.70	0
5	GOL	F	406	-	5,5,5	0.07	0	5,5,5	0.35	0
5	GOL	E	404	-	5,5,5	0.09	0	5,5,5	0.35	0
2	HEM	D	401	1	50,50,50	1.43	5 (10%)	67,82,82	1.96	19 (28%)
9	EPE	F	403	-	15,15,15	2.19	1 (6%)	19,20,20	1.82	7 (36%)
2	HEM	F	401	1	50,50,50	1.52	5 (10%)	67,82,82	1.97	18 (26%)
3	OXY	B	402	-	1,1,1	0.12	0	-	-	-
3	OXY	D	402	-	1,1,1	0.11	0	-	-	-
4	MPD	A	403	-	7,7,7	0.16	0	9,10,10	0.61	0
8	VOH	E	403	-	12,12,12	1.92	4 (33%)	15,15,15	3.04	9 (60%)
5	GOL	A	405	-	5,5,5	0.08	0	5,5,5	0.32	0
3	OXY	F	402	-	1,1,1	0.02	0	-	-	-

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	OXY	E	402	-	1,1,1	0.16	0	-		
2	HEM	E	401	1	50,50,50	1.49	8 (16%)	67,82,82	1.76	17 (25%)
4	MPD	C	404	-	7,7,7	0.27	0	9,10,10	0.67	0
5	GOL	A	404	-	5,5,5	0.10	0	5,5,5	0.34	0
7	PO4	A	408	-	4,4,4	0.61	0	6,6,6	0.54	0
2	HEM	B	401	1	50,50,50	1.46	5 (10%)	67,82,82	1.83	15 (22%)
2	HEM	C	401	1,3	50,50,50	1.58	9 (18%)	67,82,82	2.06	16 (23%)
5	GOL	A	406	-	5,5,5	0.15	0	5,5,5	0.21	0
8	VOH	C	403	-	12,12,12	1.80	3 (25%)	15,15,15	3.17	5 (33%)
5	GOL	F	407	-	5,5,5	0.41	0	5,5,5	1.06	0
3	OXY	A	402	-	1,1,1	0.19	0	-		
5	GOL	C	405	-	5,5,5	0.17	0	5,5,5	0.53	0
2	HEM	A	401	1	50,50,50	1.55	5 (10%)	67,82,82	1.84	16 (23%)
5	GOL	B	403	-	5,5,5	0.20	0	5,5,5	0.52	0
5	GOL	F	404	-	5,5,5	0.12	0	5,5,5	0.45	0
5	GOL	F	405	-	5,5,5	0.13	0	5,5,5	0.62	0
5	GOL	E	405	-	5,5,5	0.20	0	5,5,5	0.42	0
5	GOL	C	406	-	5,5,5	0.16	0	5,5,5	0.58	0
3	OXY	C	402	2	1,1,1	0.01	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	D	403	-	-	2/4/4/4	-
5	GOL	F	406	-	-	0/4/4/4	-
5	GOL	E	404	-	-	4/4/4/4	-
2	HEM	D	401	1	-	4/14/54/54	-
9	EPE	F	403	-	-	2/9/19/19	0/1/1/1
2	HEM	F	401	1	-	5/14/54/54	-
8	VOH	E	403	-	-	2/6/6/6	0/1/1/1
4	MPD	A	403	-	-	3/5/5/5	-
5	GOL	A	405	-	-	1/4/4/4	-
2	HEM	E	401	1	-	4/14/54/54	-
4	MPD	C	404	-	-	3/5/5/5	-
5	GOL	A	404	-	-	4/4/4/4	-
2	HEM	B	401	1	-	4/14/54/54	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	HEM	C	401	1,3	-	4/14/54/54	-
5	GOL	A	406	-	-	0/4/4/4	-
8	VOH	C	403	-	-	2/6/6/6	0/1/1/1
5	GOL	F	407	-	-	2/4/4/4	-
5	GOL	C	405	-	-	2/4/4/4	-
2	HEM	A	401	1	-	4/14/54/54	-
5	GOL	B	403	-	-	3/4/4/4	-
5	GOL	F	404	-	-	2/4/4/4	-
5	GOL	F	405	-	-	4/4/4/4	-
5	GOL	E	405	-	-	4/4/4/4	-
5	GOL	C	406	-	-	2/4/4/4	-

The worst 5 of 45 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
9	F	403	EPE	C10-S	-7.95	1.66	1.77
2	F	401	HEM	FE-NB	6.09	2.13	1.94
2	A	401	HEM	C1B-NB	-5.05	1.31	1.40
2	D	401	HEM	FE-NC	4.80	2.11	1.95
8	C	403	VOH	CAG-CAJ	-4.78	1.35	1.51

The worst 5 of 122 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
8	C	403	VOH	CAA-OAH-CAK	-9.14	104.10	117.51
2	C	401	HEM	C1B-NB-C4B	6.42	112.81	105.21
2	C	401	HEM	CHC-C4B-NB	6.41	131.32	124.42
2	A	401	HEM	CHC-C4B-NB	6.27	131.17	124.42
2	D	401	HEM	CHC-C4B-NB	6.22	131.12	124.42

There are no chirality outliers.

5 of 67 torsion outliers are listed below:

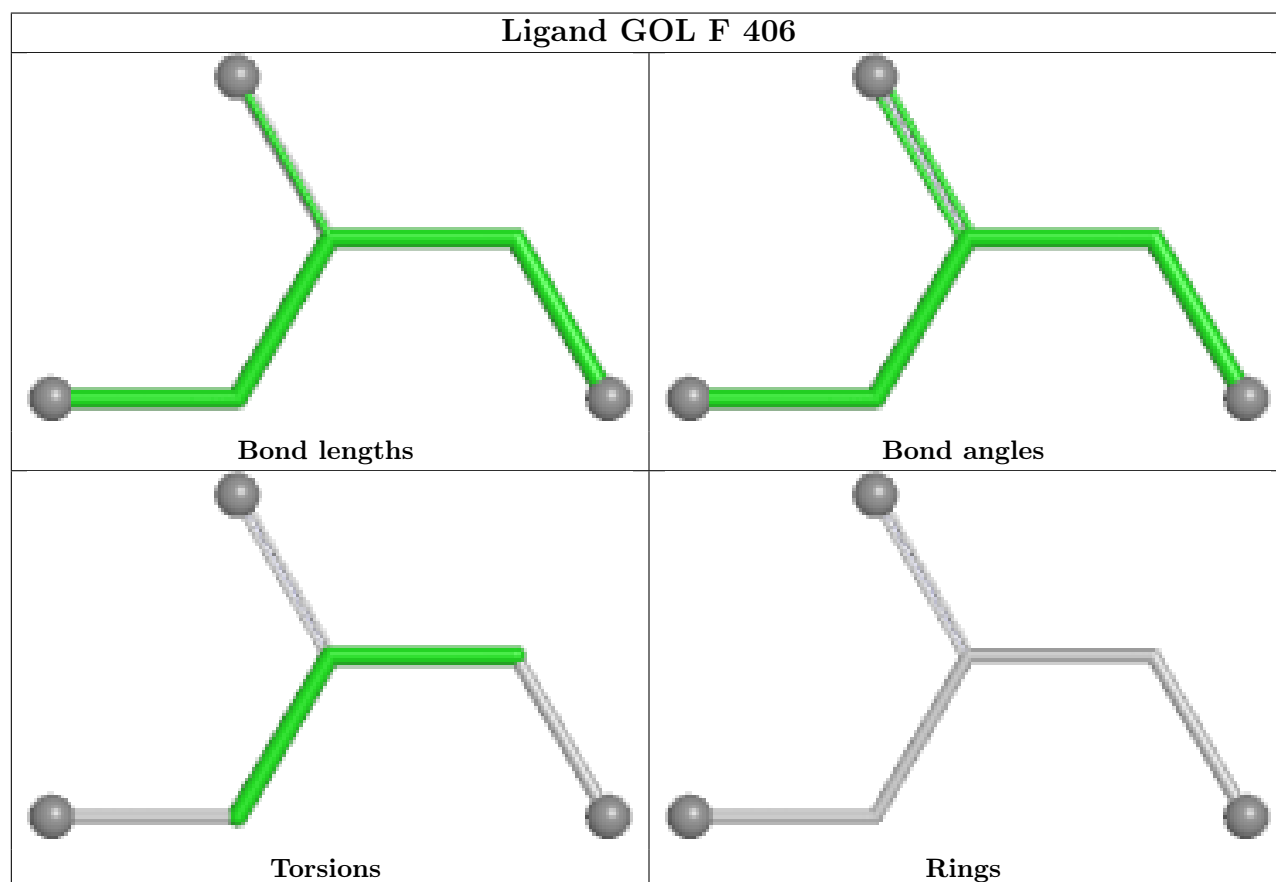
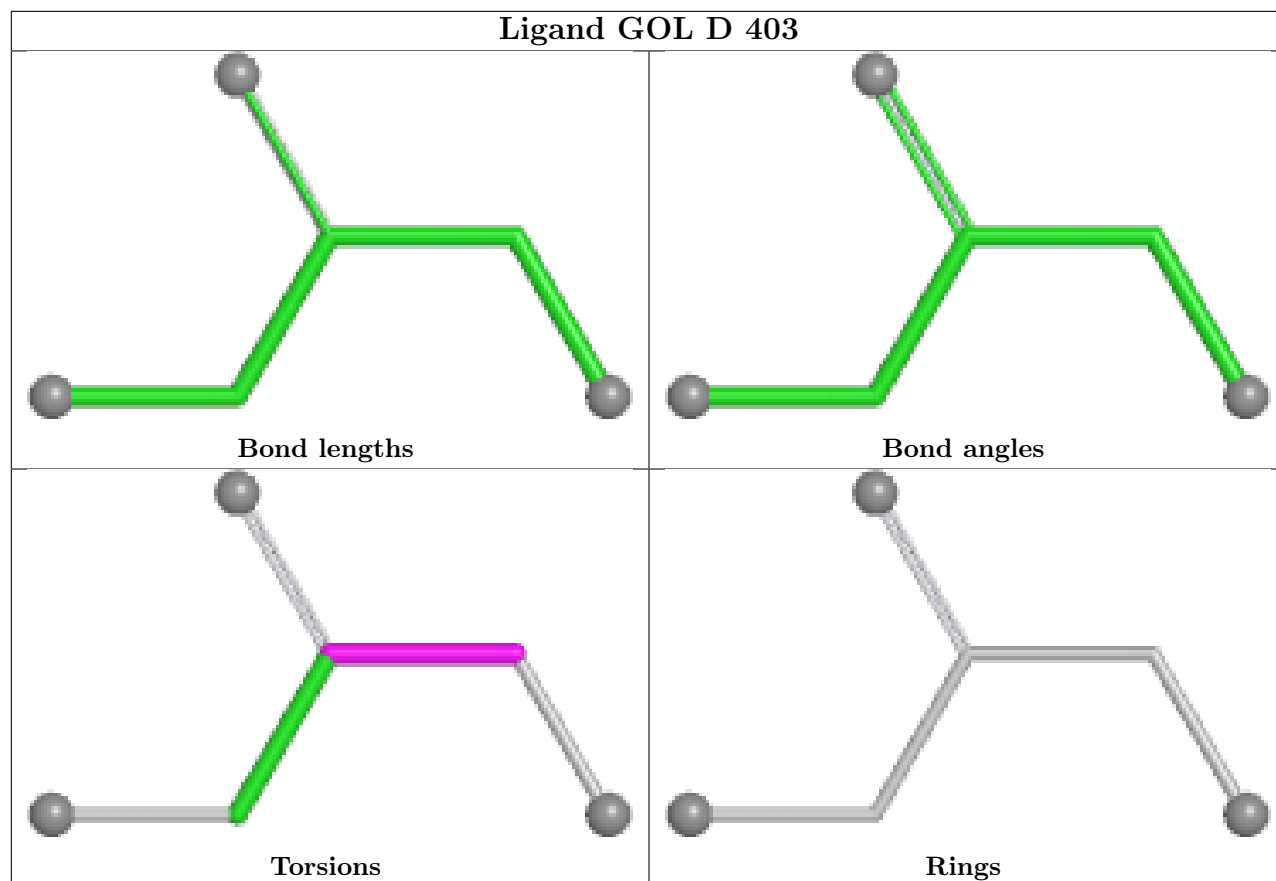
Mol	Chain	Res	Type	Atoms
4	A	403	MPD	CM-C2-C3-C4
5	A	404	GOL	O1-C1-C2-O2
5	A	404	GOL	O1-C1-C2-C3
5	A	404	GOL	C1-C2-C3-O3
5	A	404	GOL	O2-C2-C3-O3

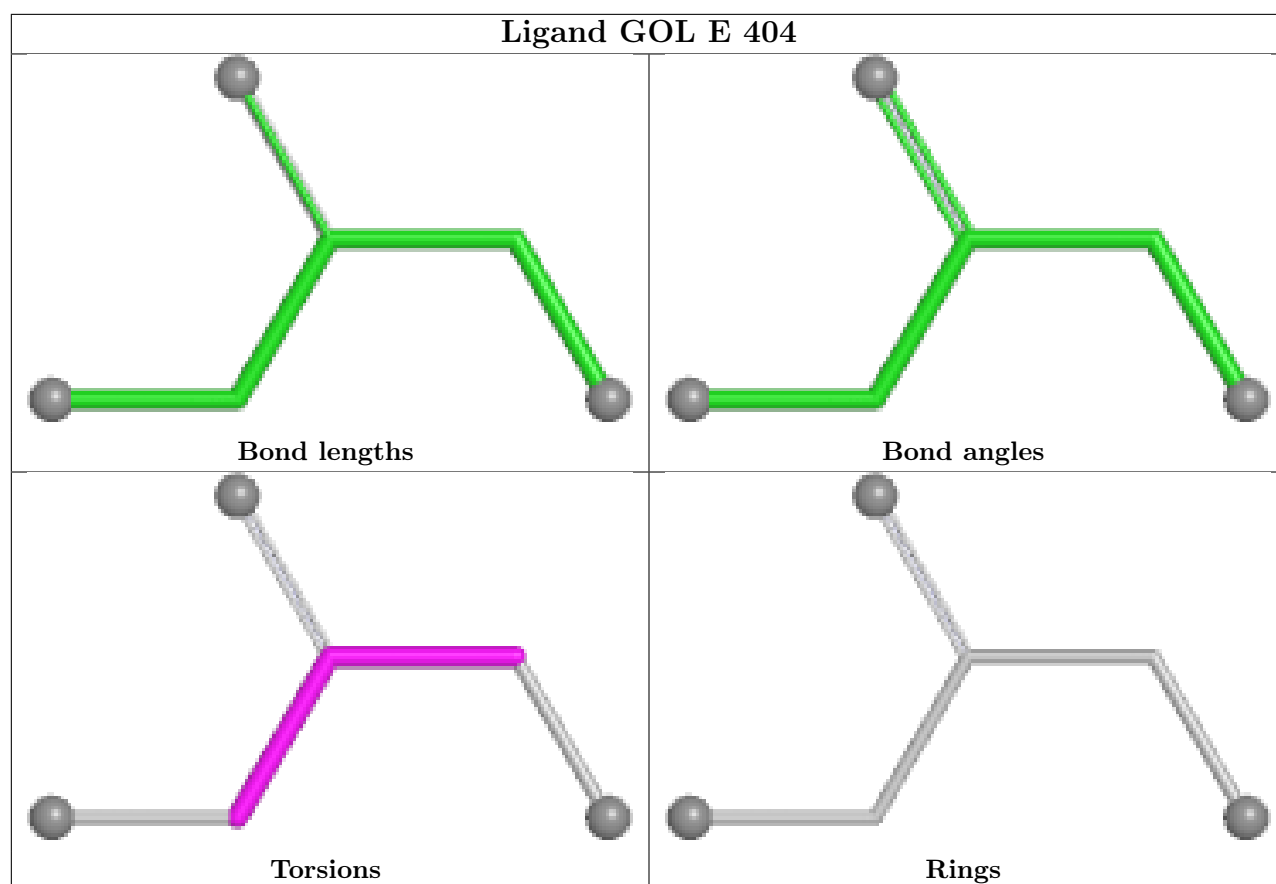
There are no ring outliers.

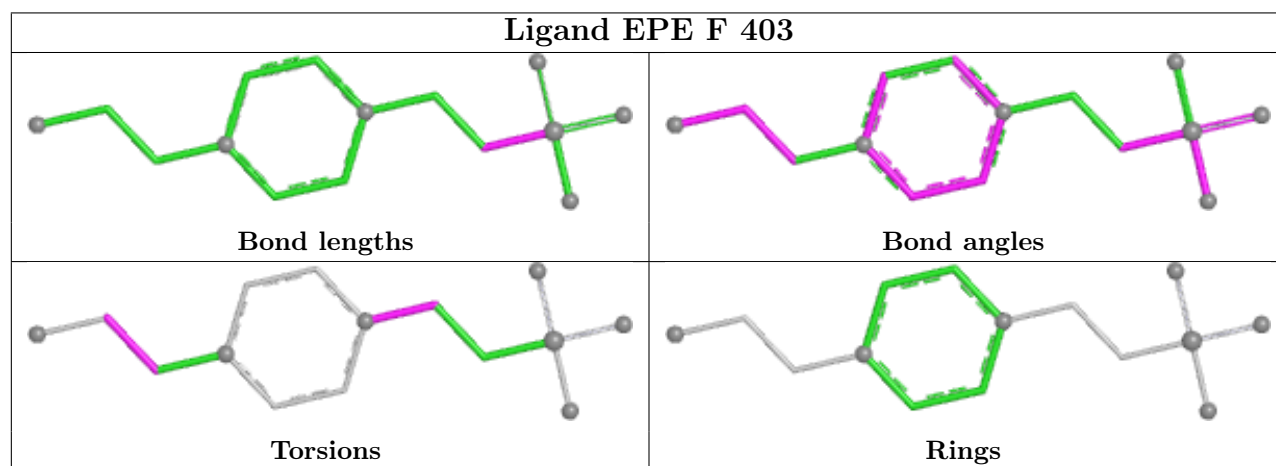
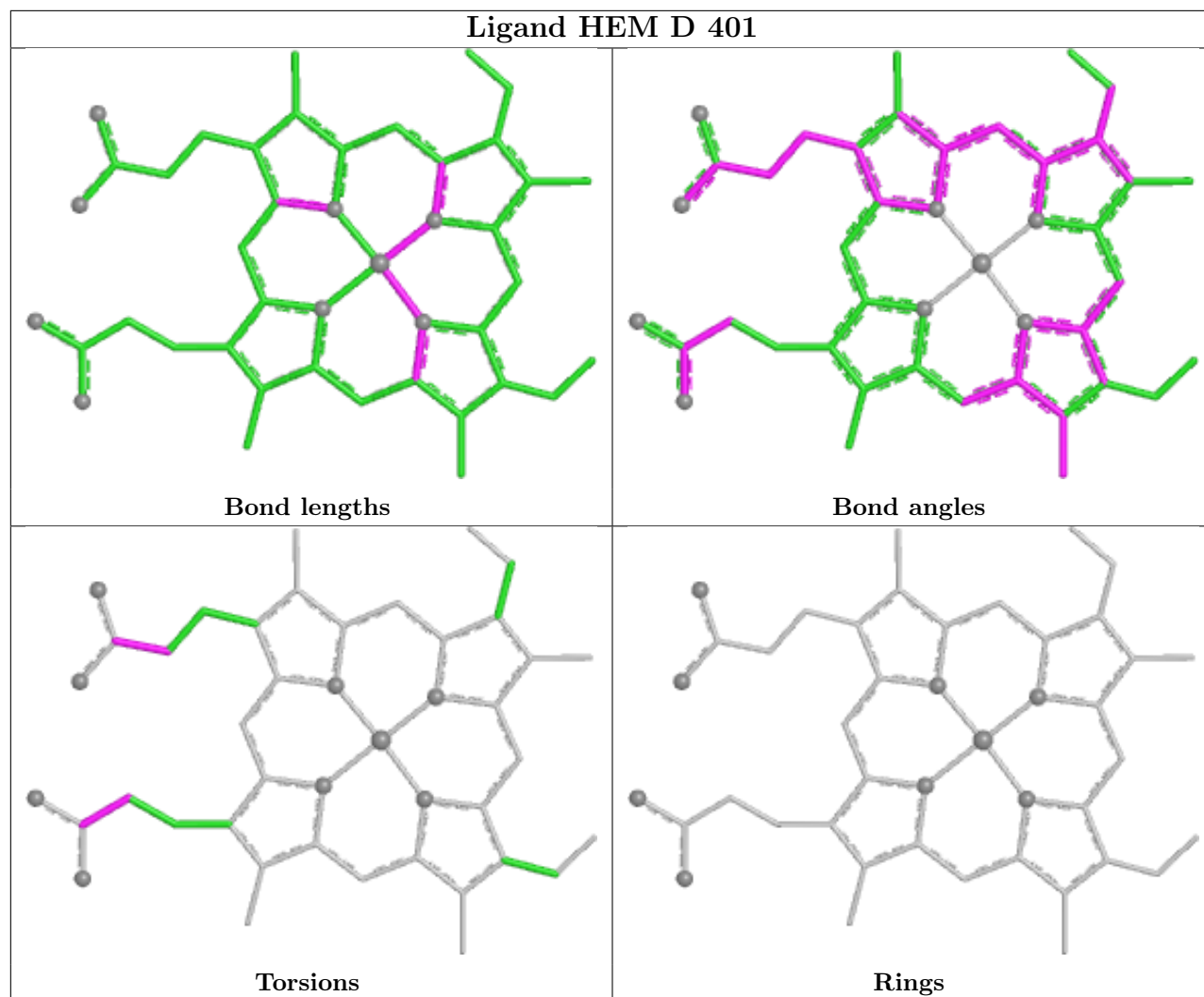
17 monomers are involved in 60 short contacts:

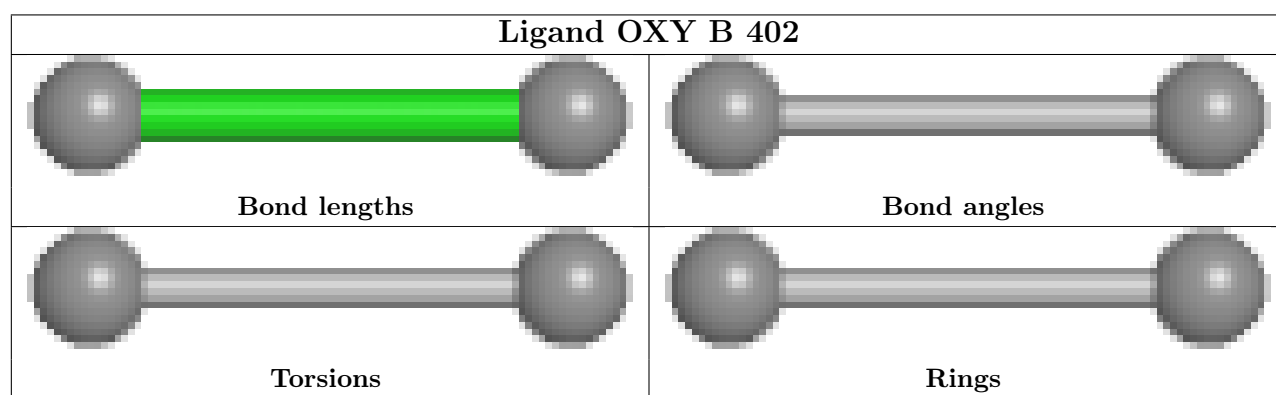
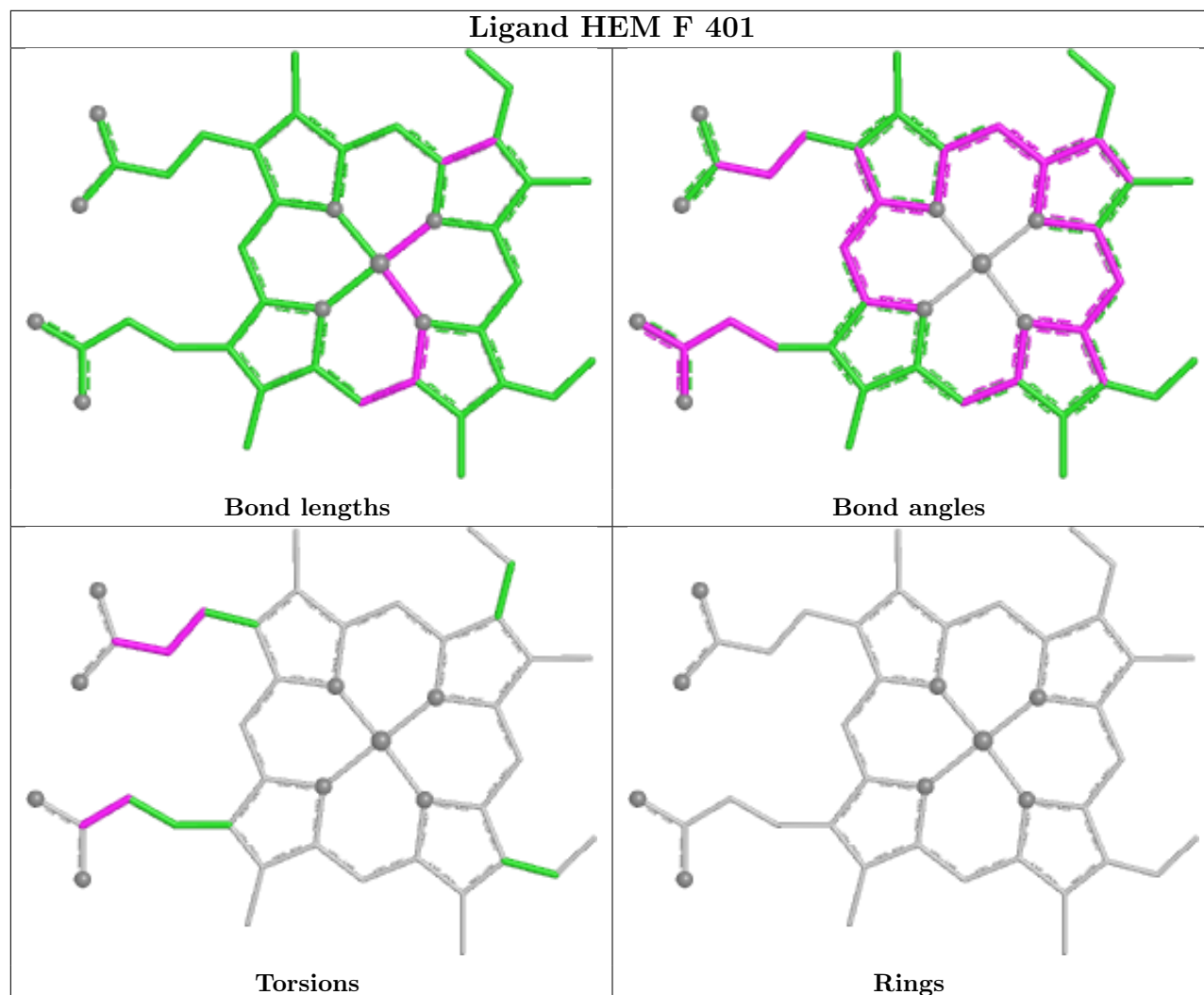
Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	D	403	GOL	1	0
2	D	401	HEM	3	0
9	F	403	EPE	2	0
2	F	401	HEM	4	0
4	A	403	MPD	5	0
8	E	403	VOH	6	0
5	A	405	GOL	3	0
4	C	404	MPD	8	0
5	A	404	GOL	1	0
2	B	401	HEM	2	0
2	C	401	HEM	4	0
8	C	403	VOH	7	0
5	F	407	GOL	6	0
5	C	405	GOL	2	0
2	A	401	HEM	3	0
5	F	405	GOL	1	0
5	C	406	GOL	2	0

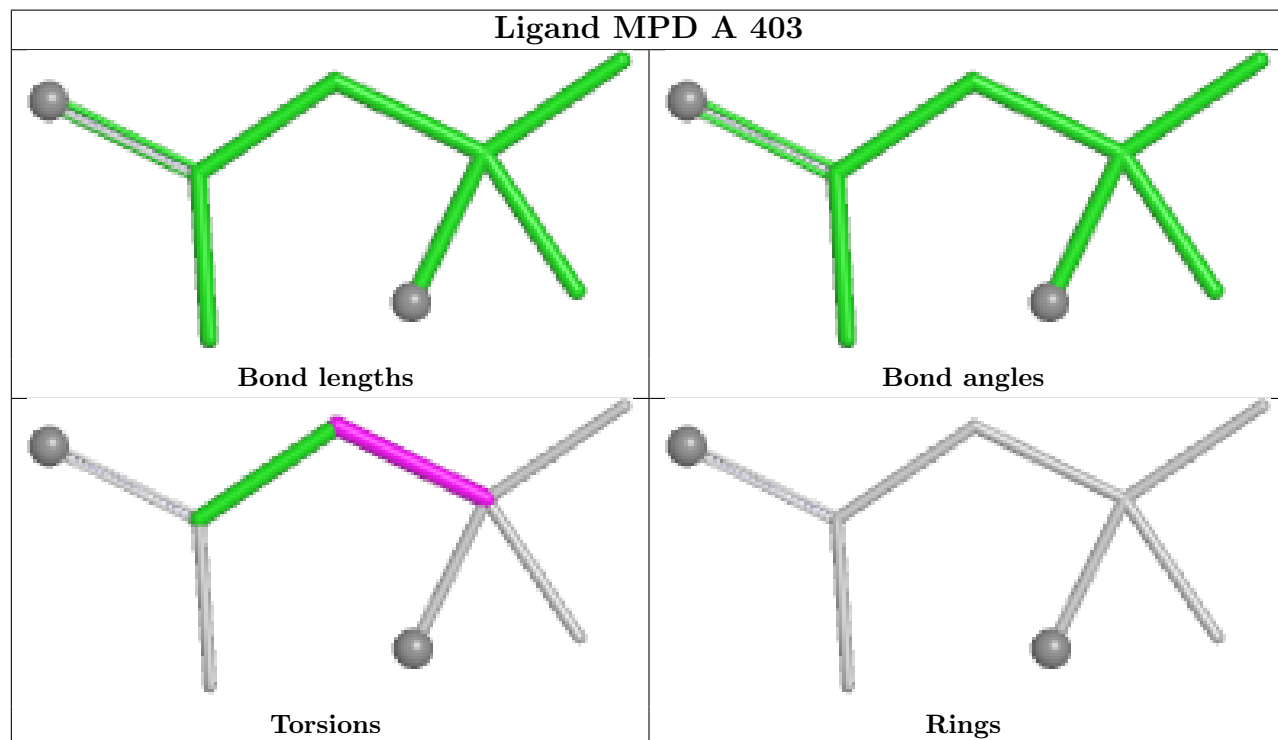
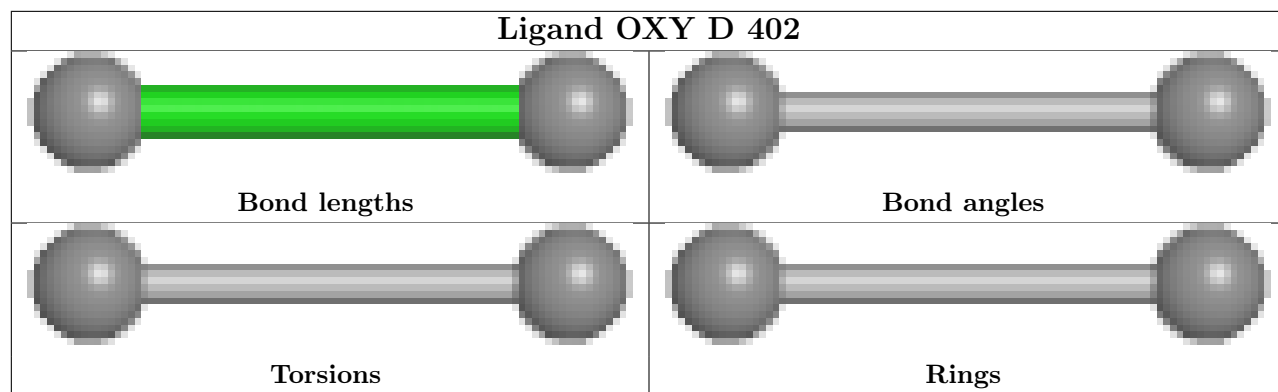
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.

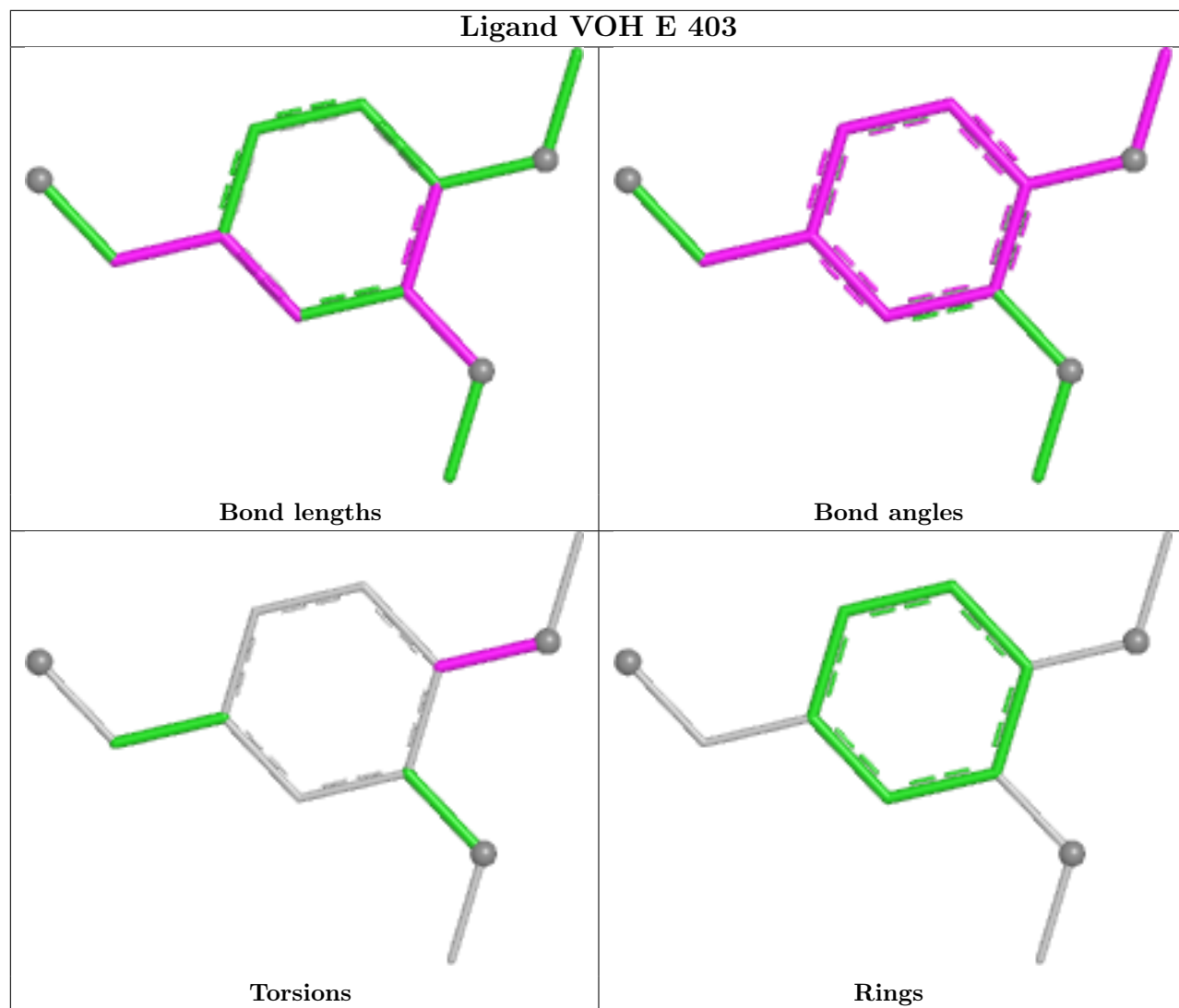


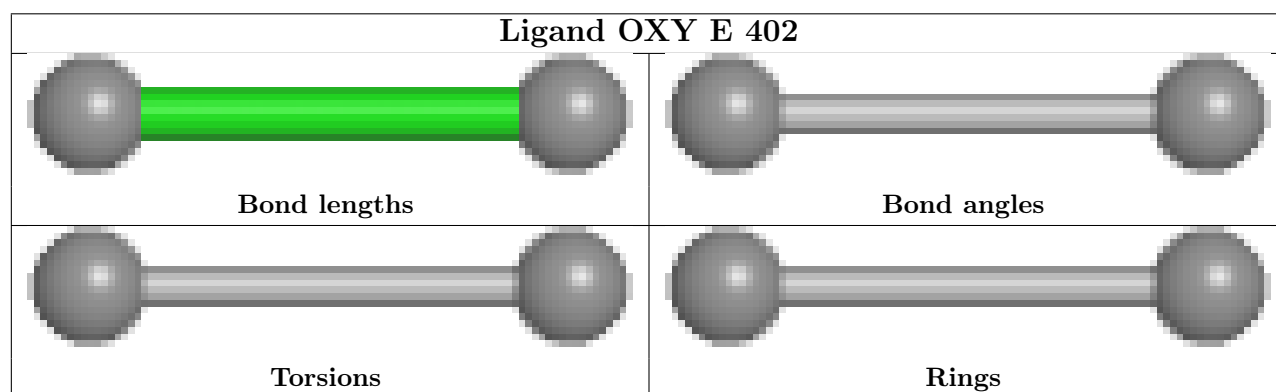
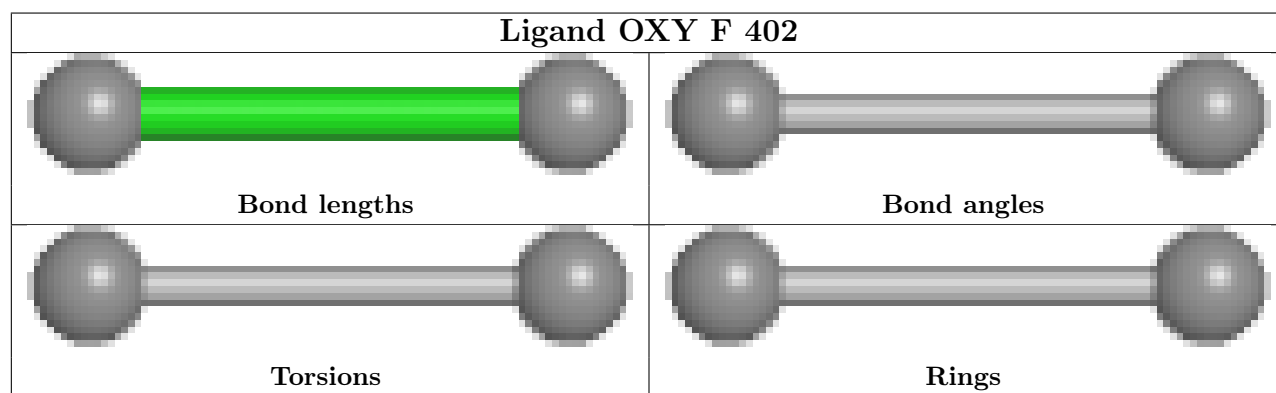
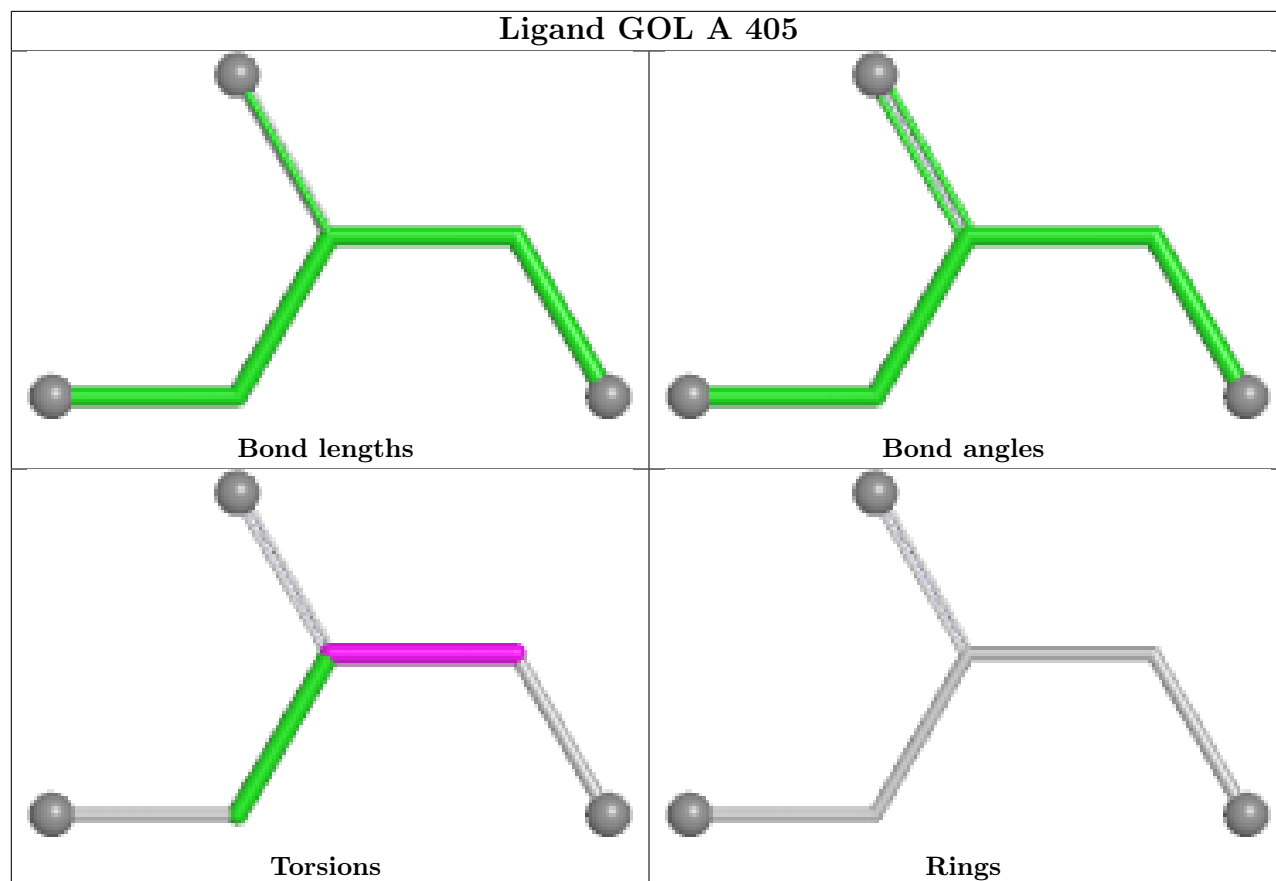


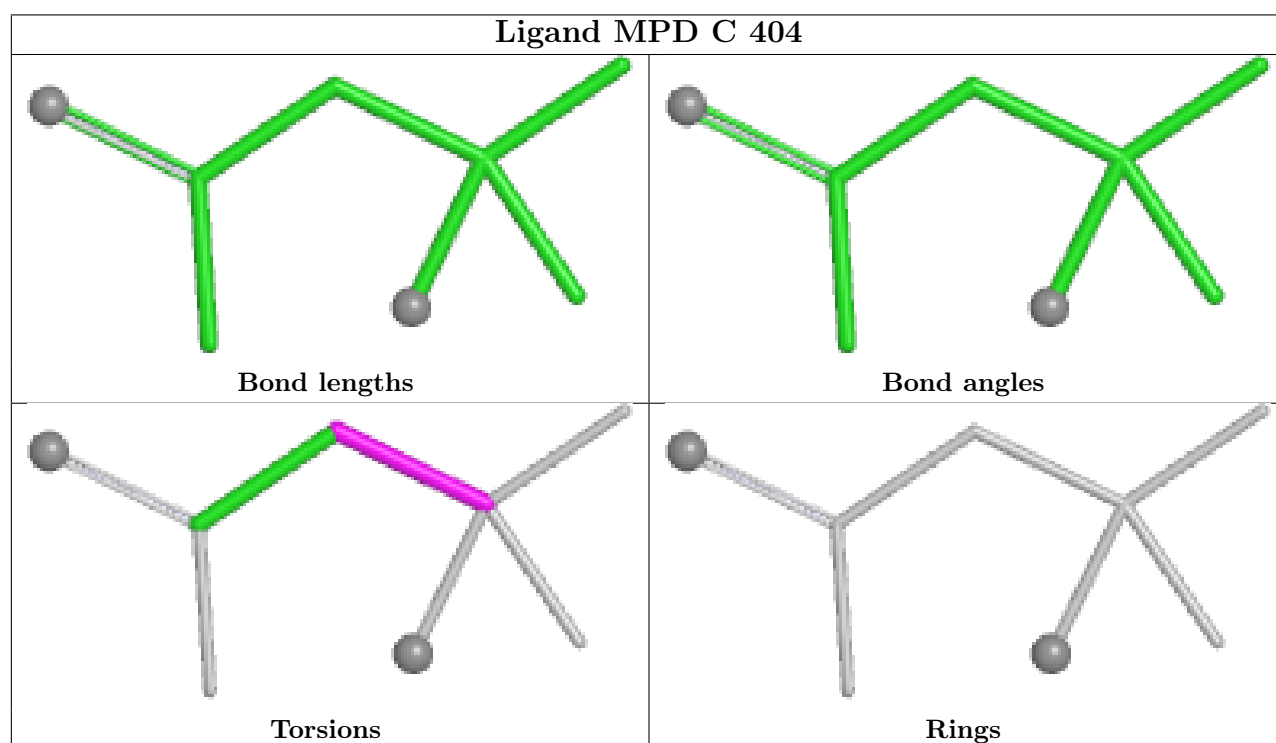
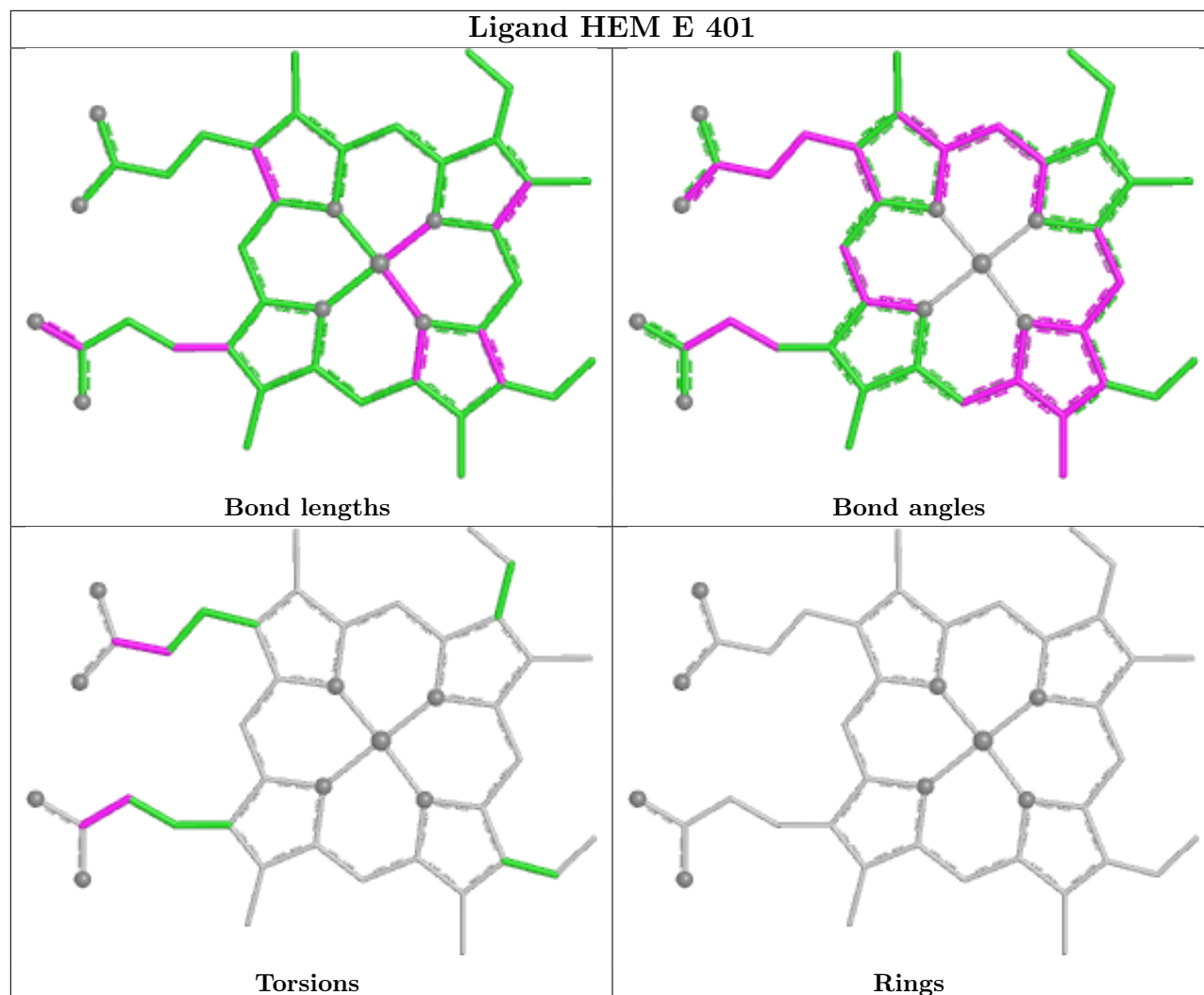


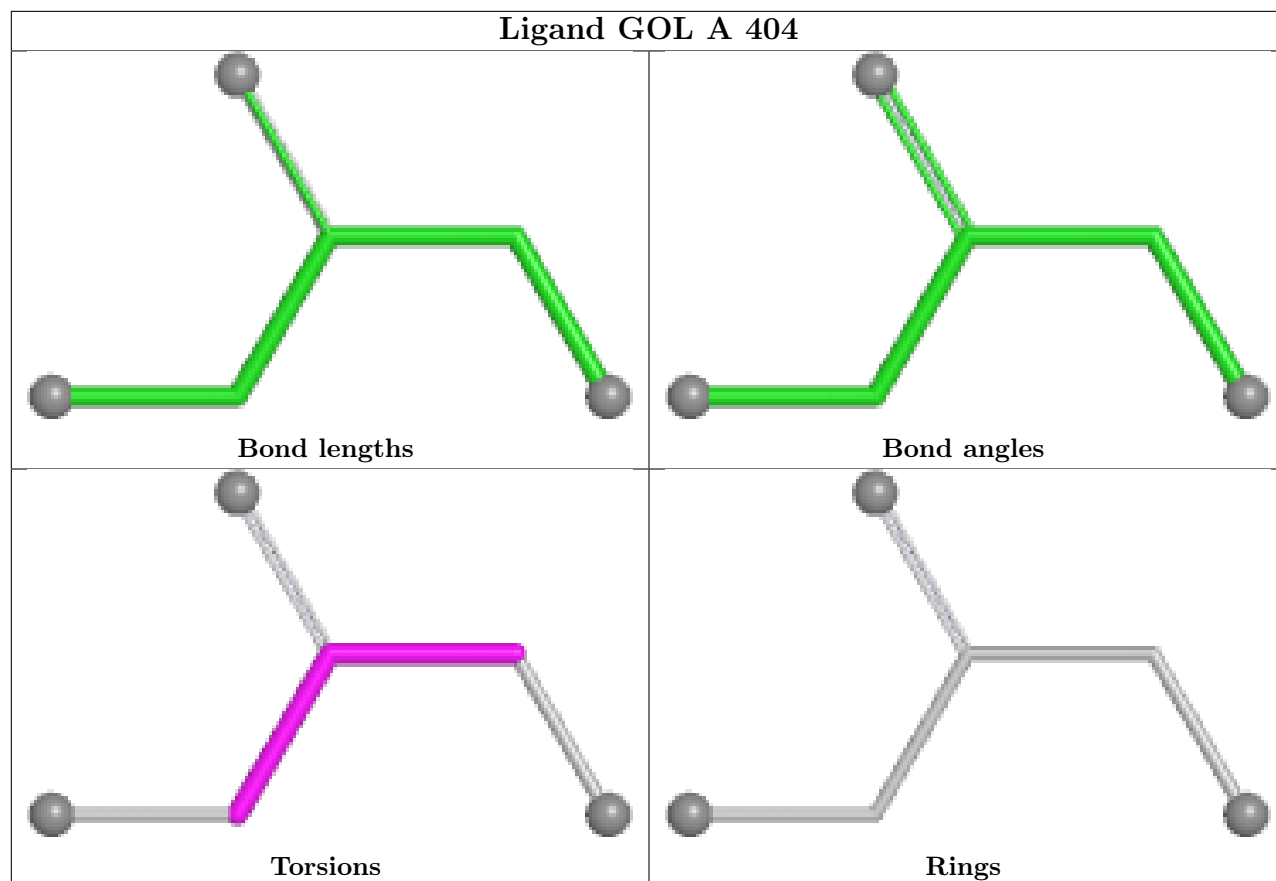


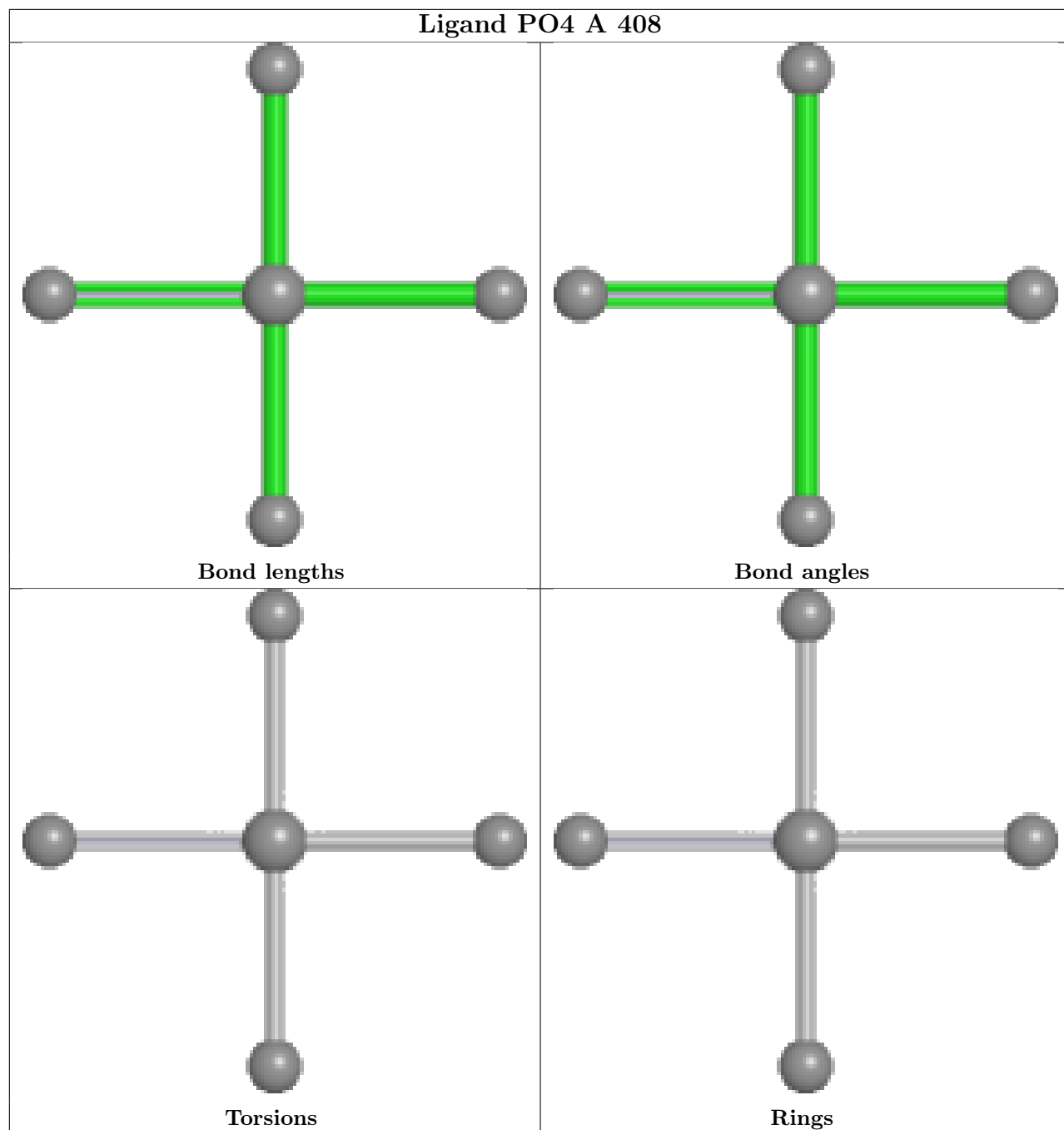


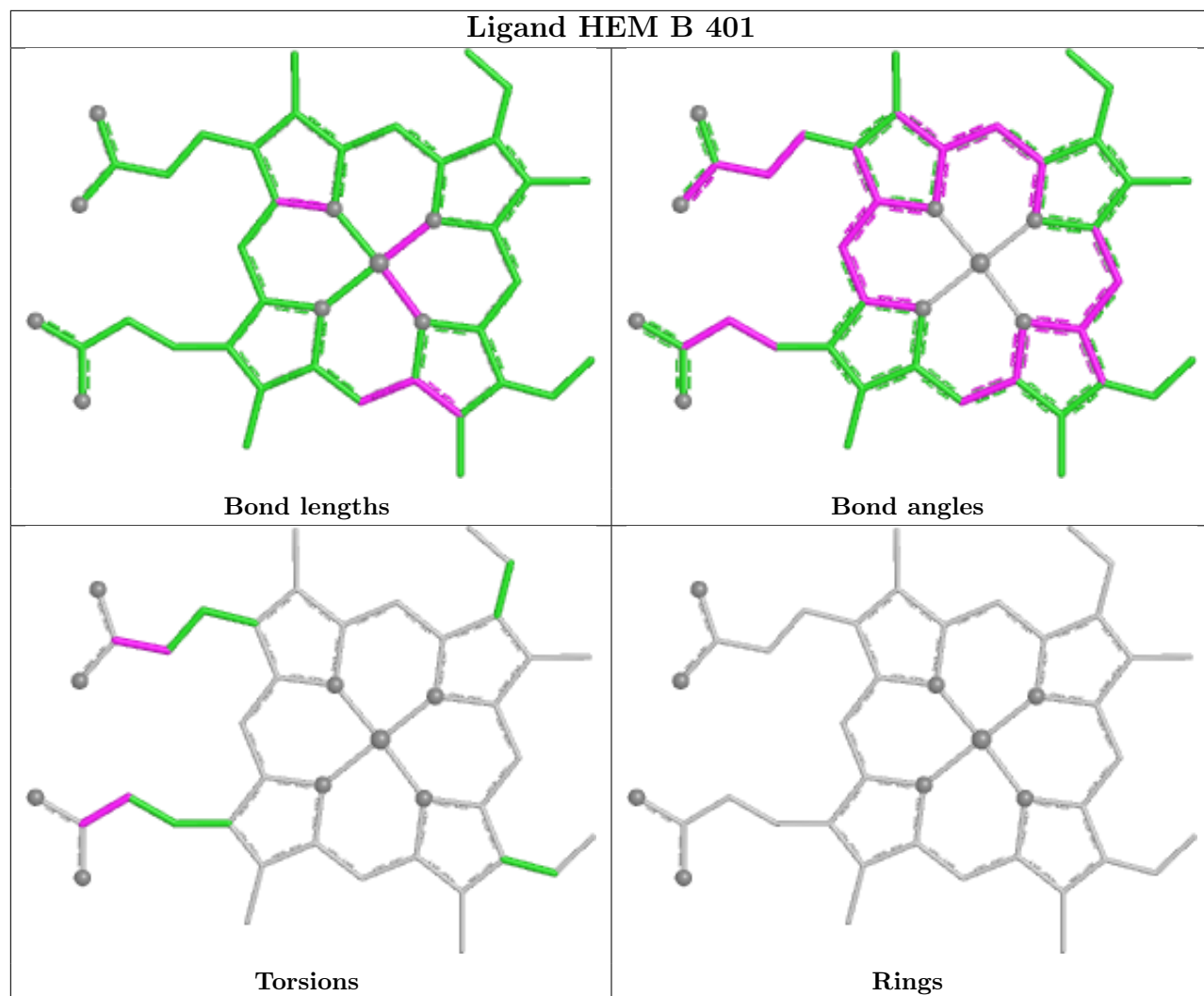


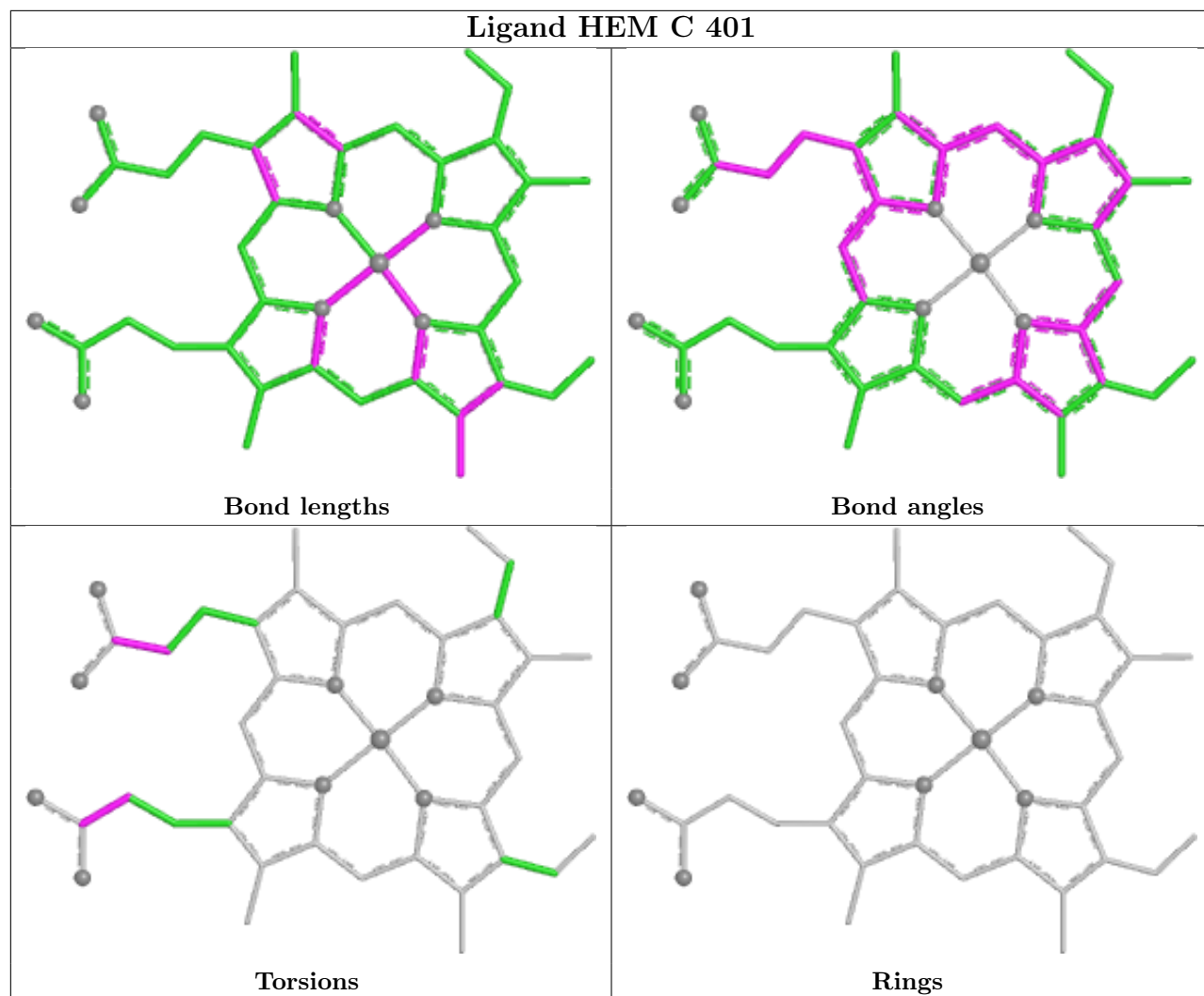


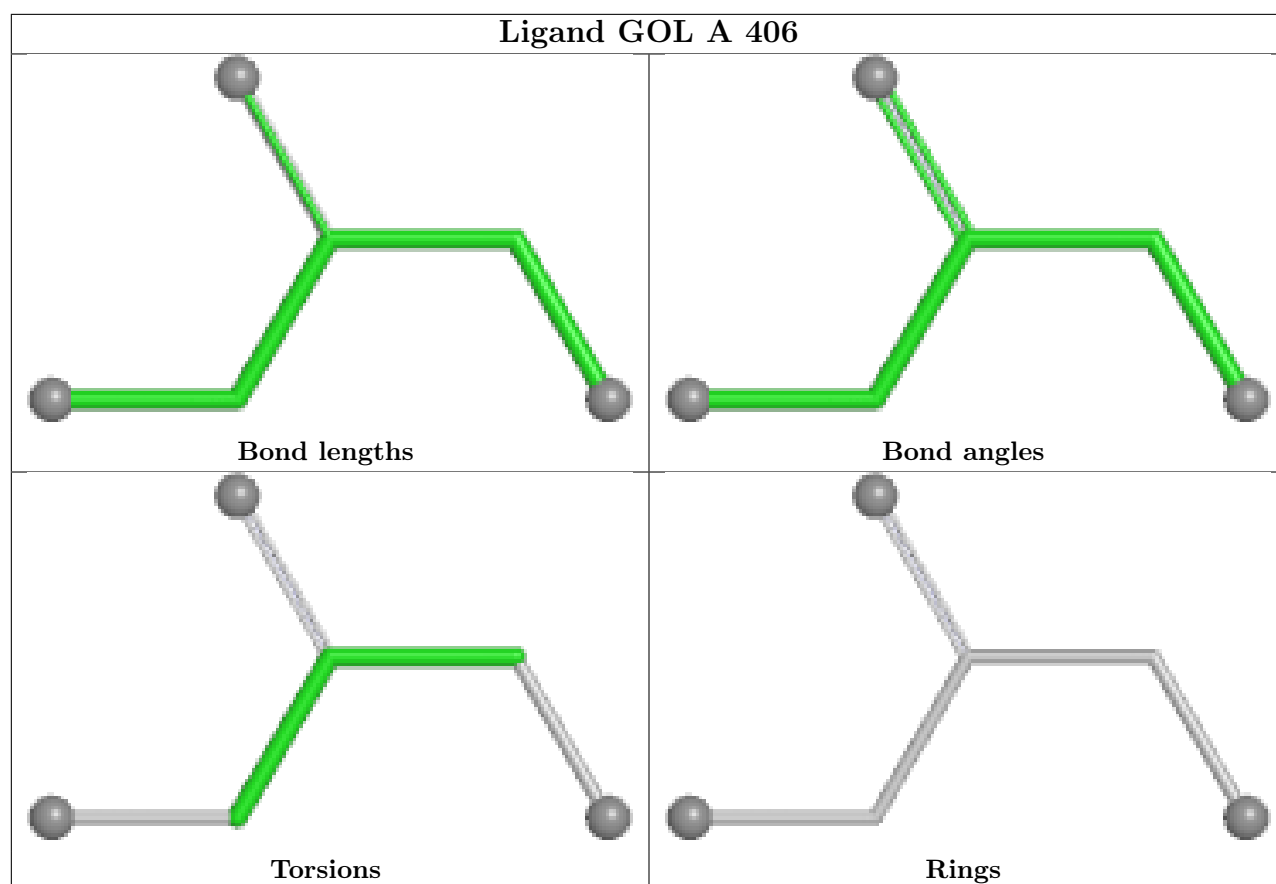


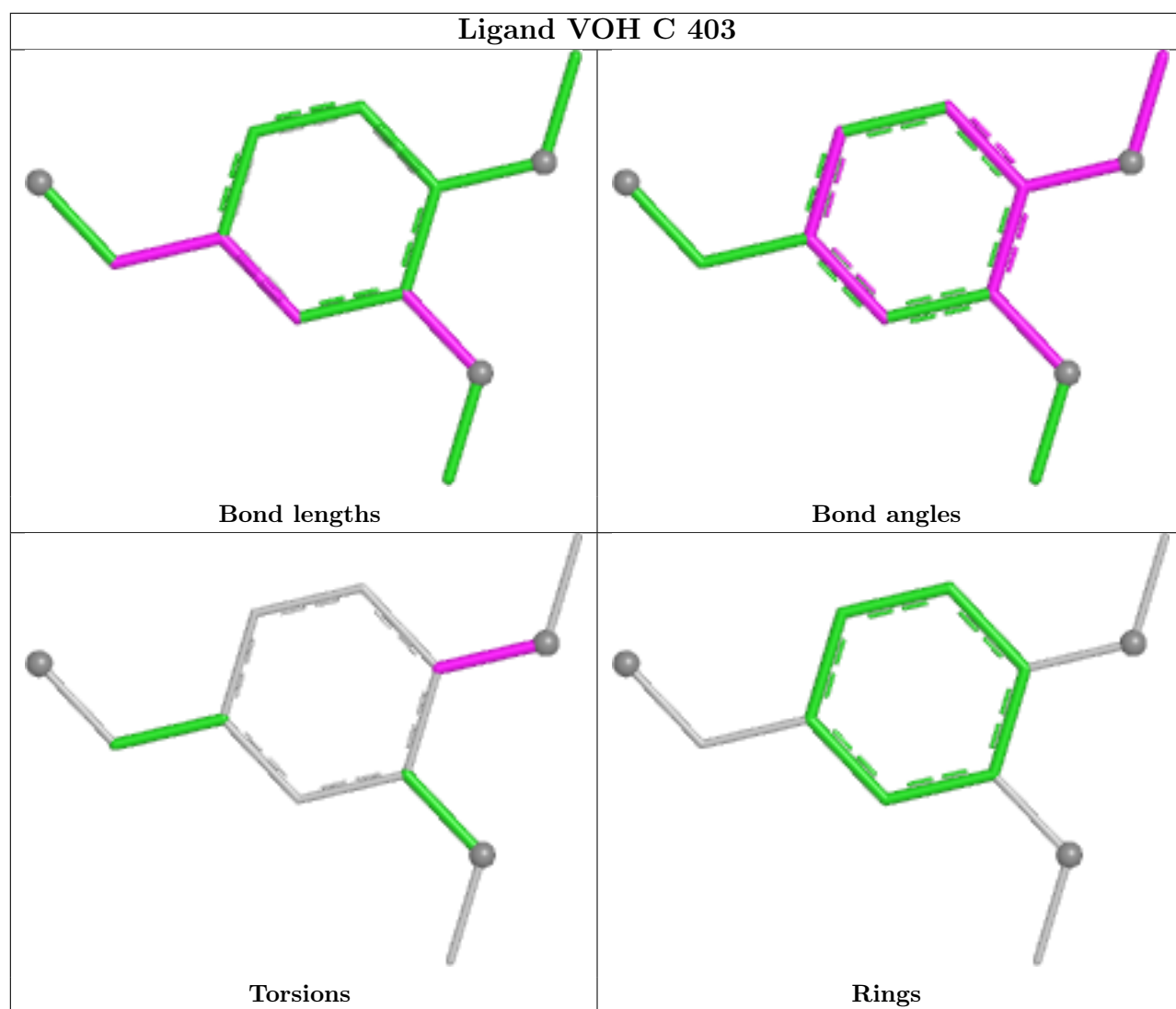


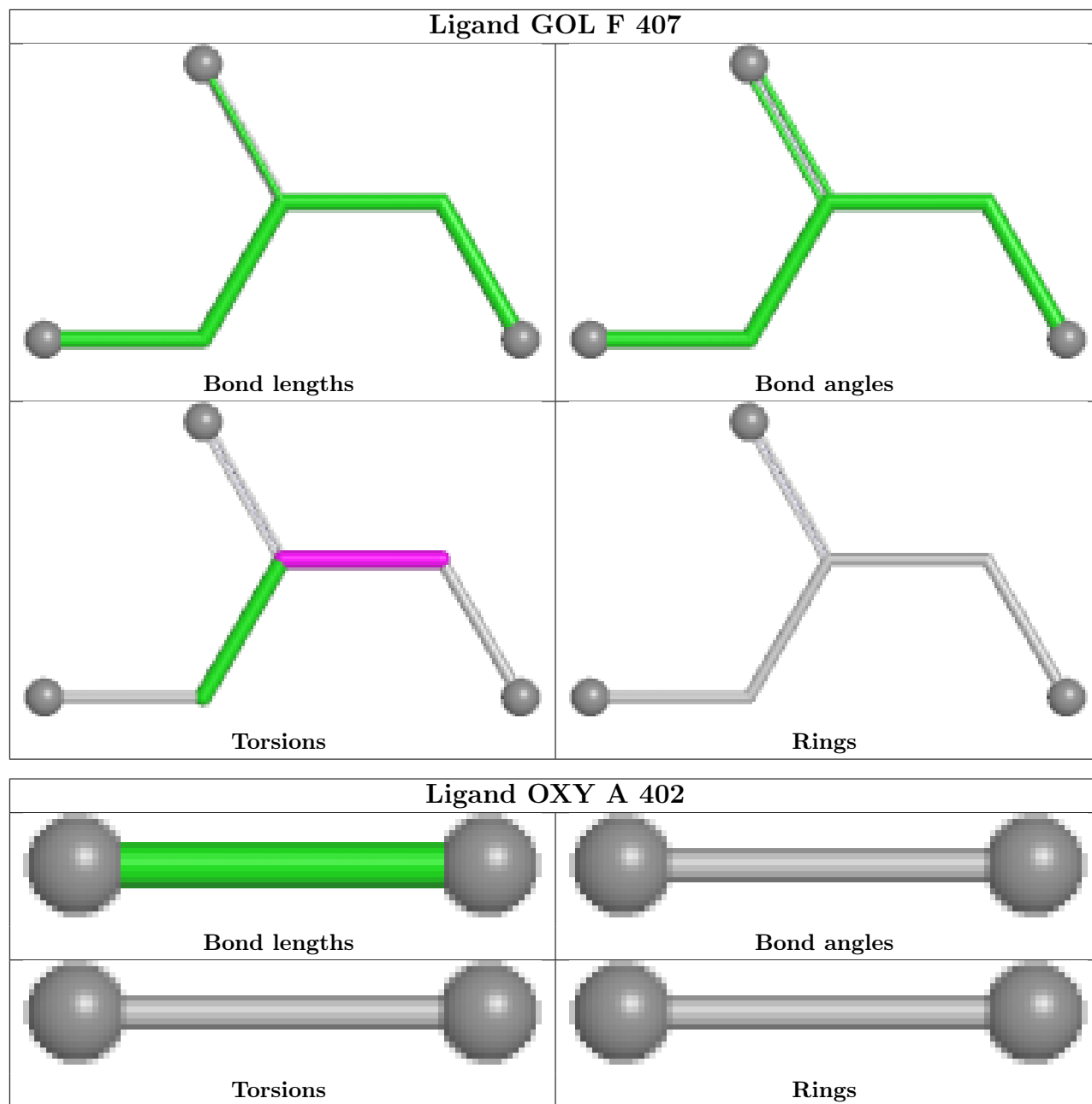


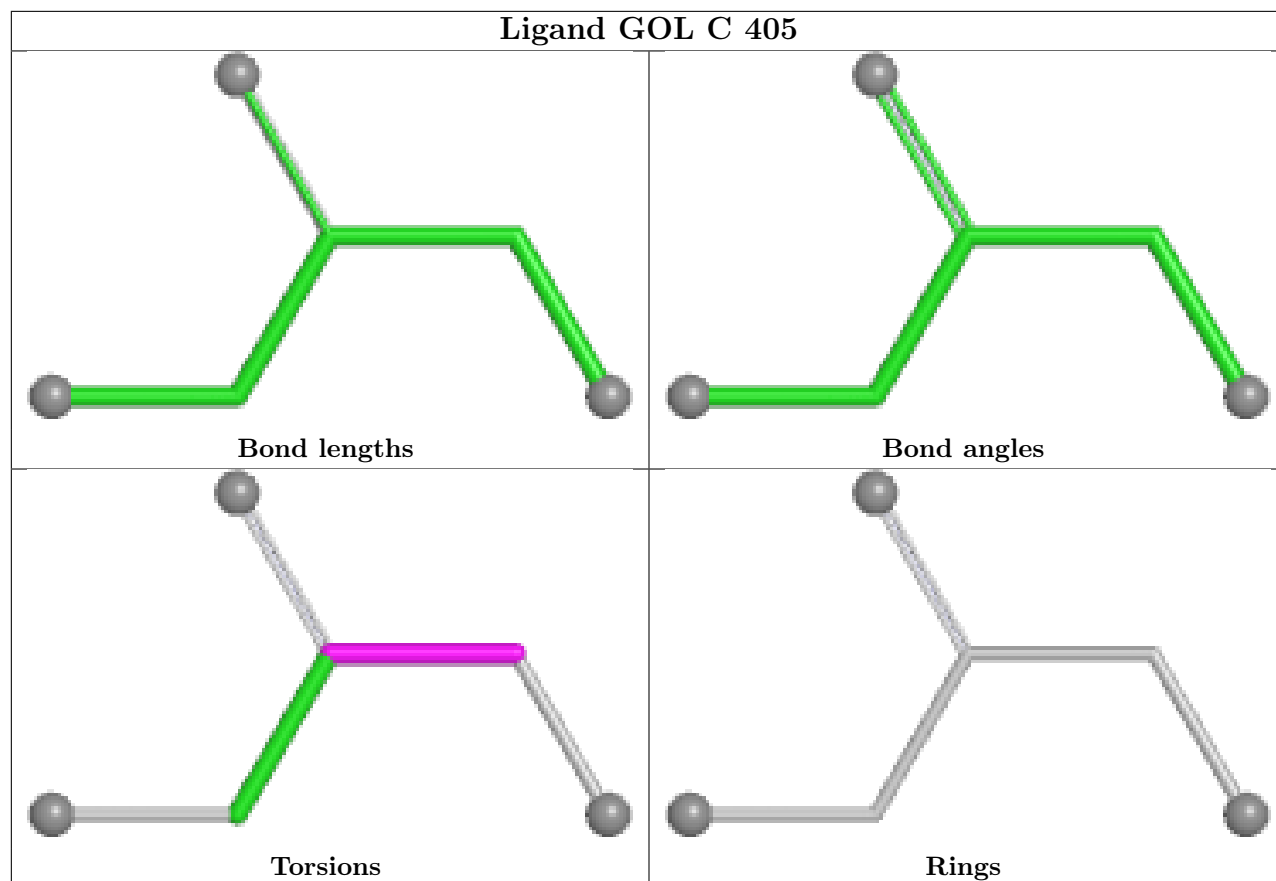


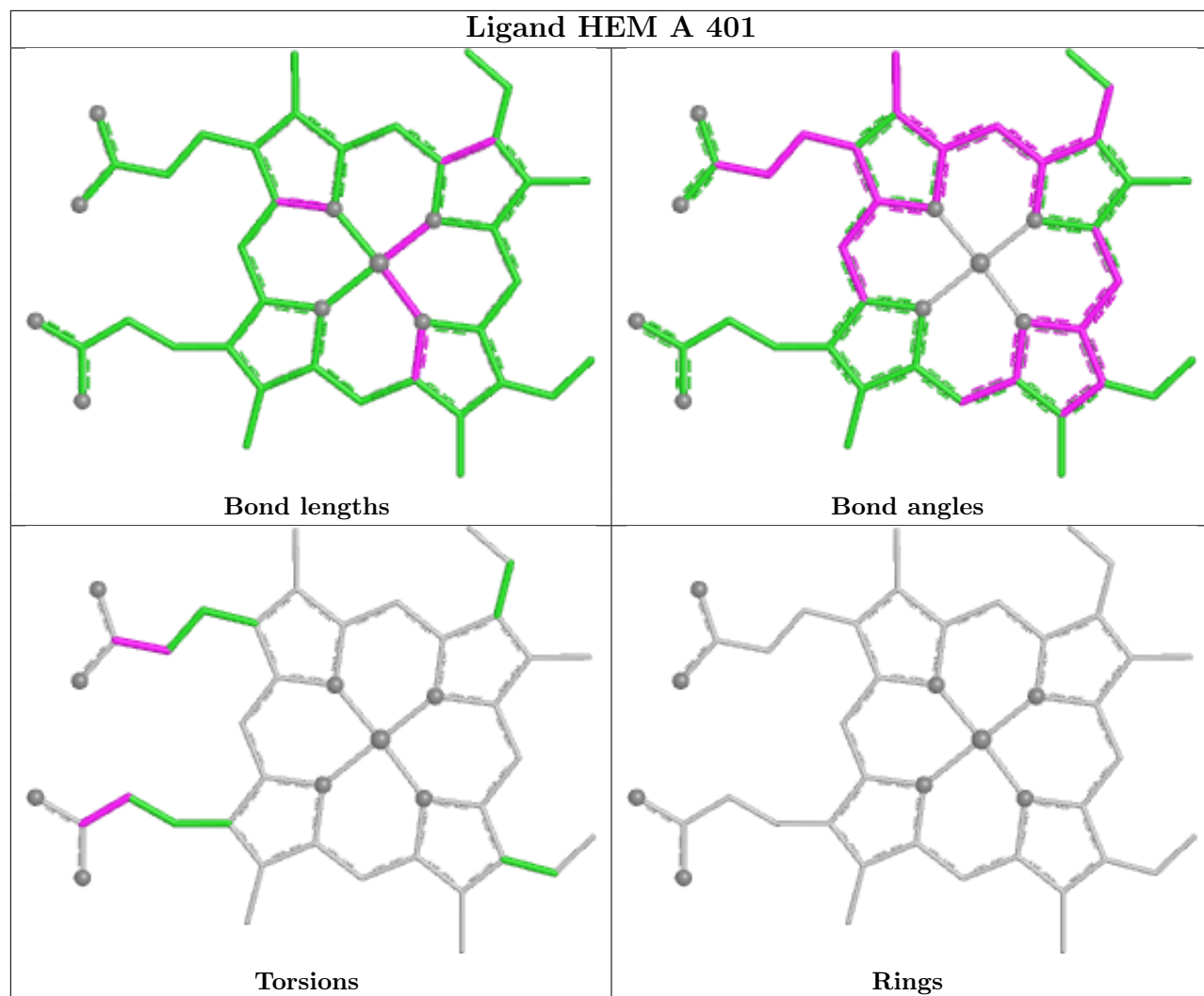


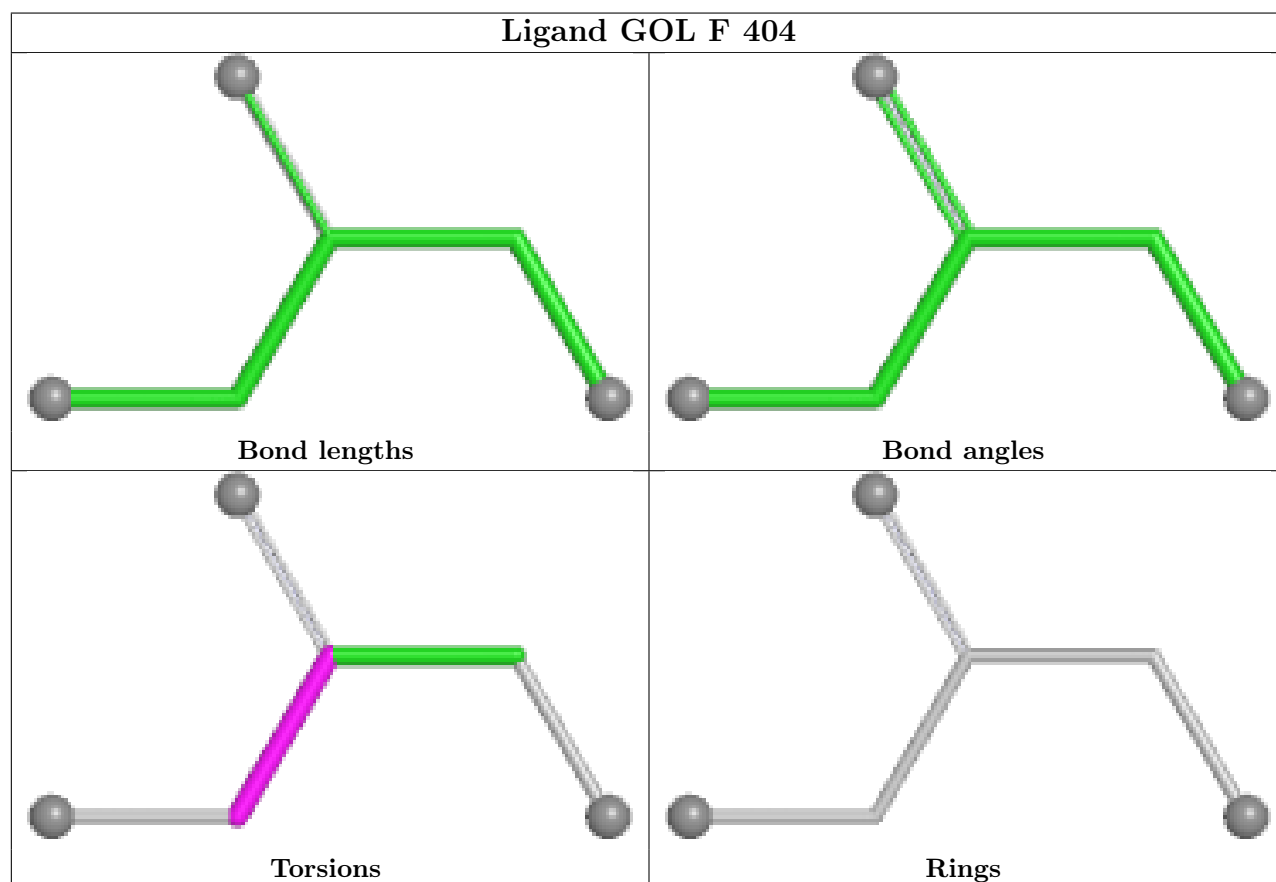
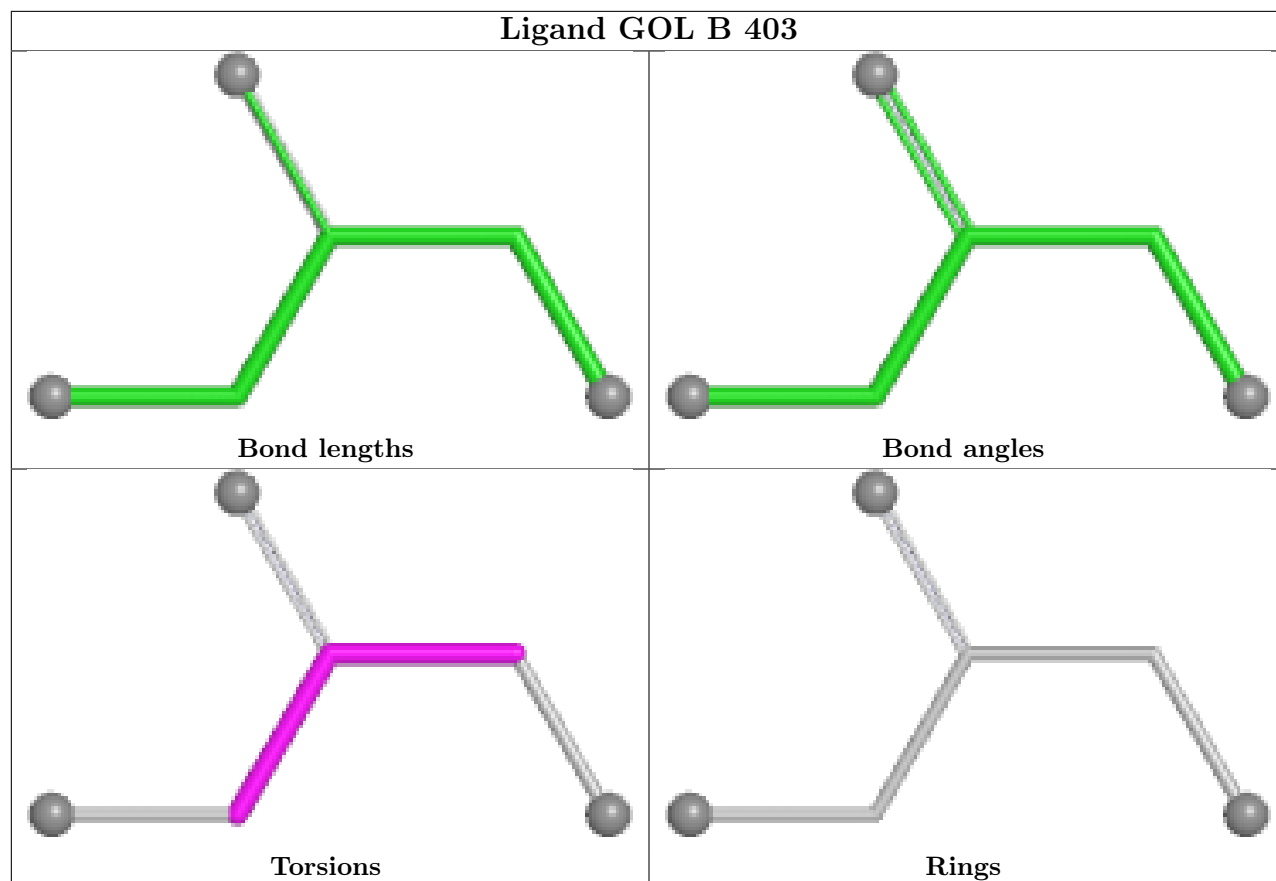


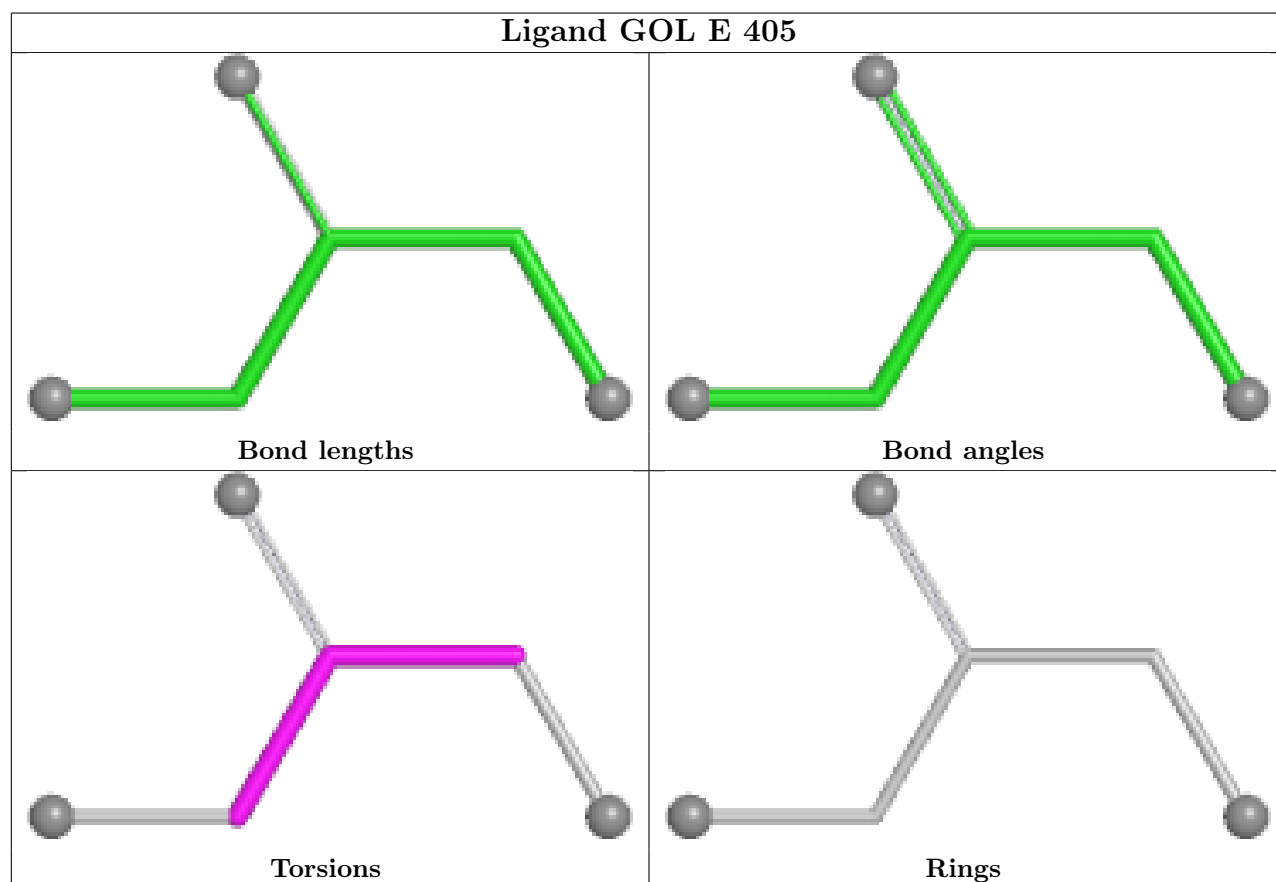
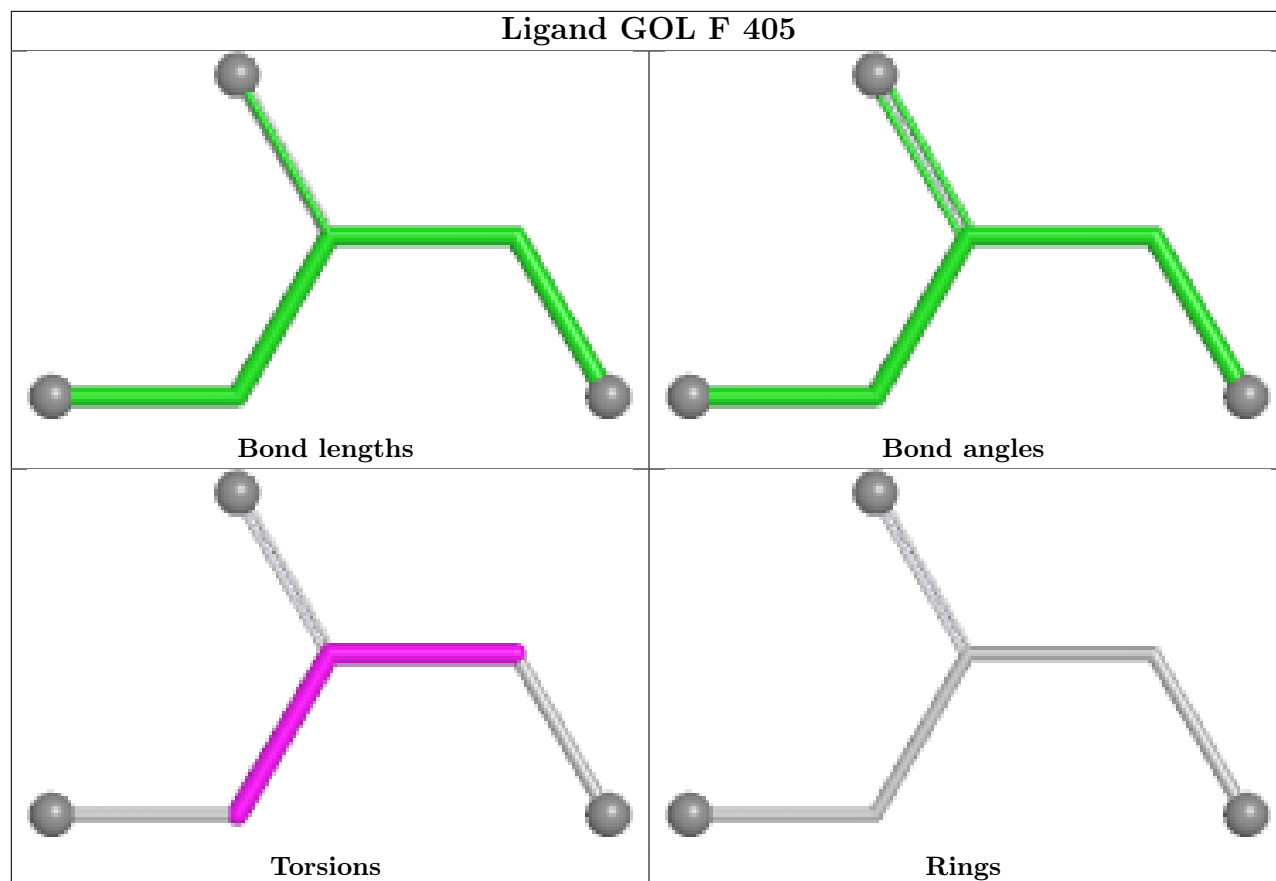


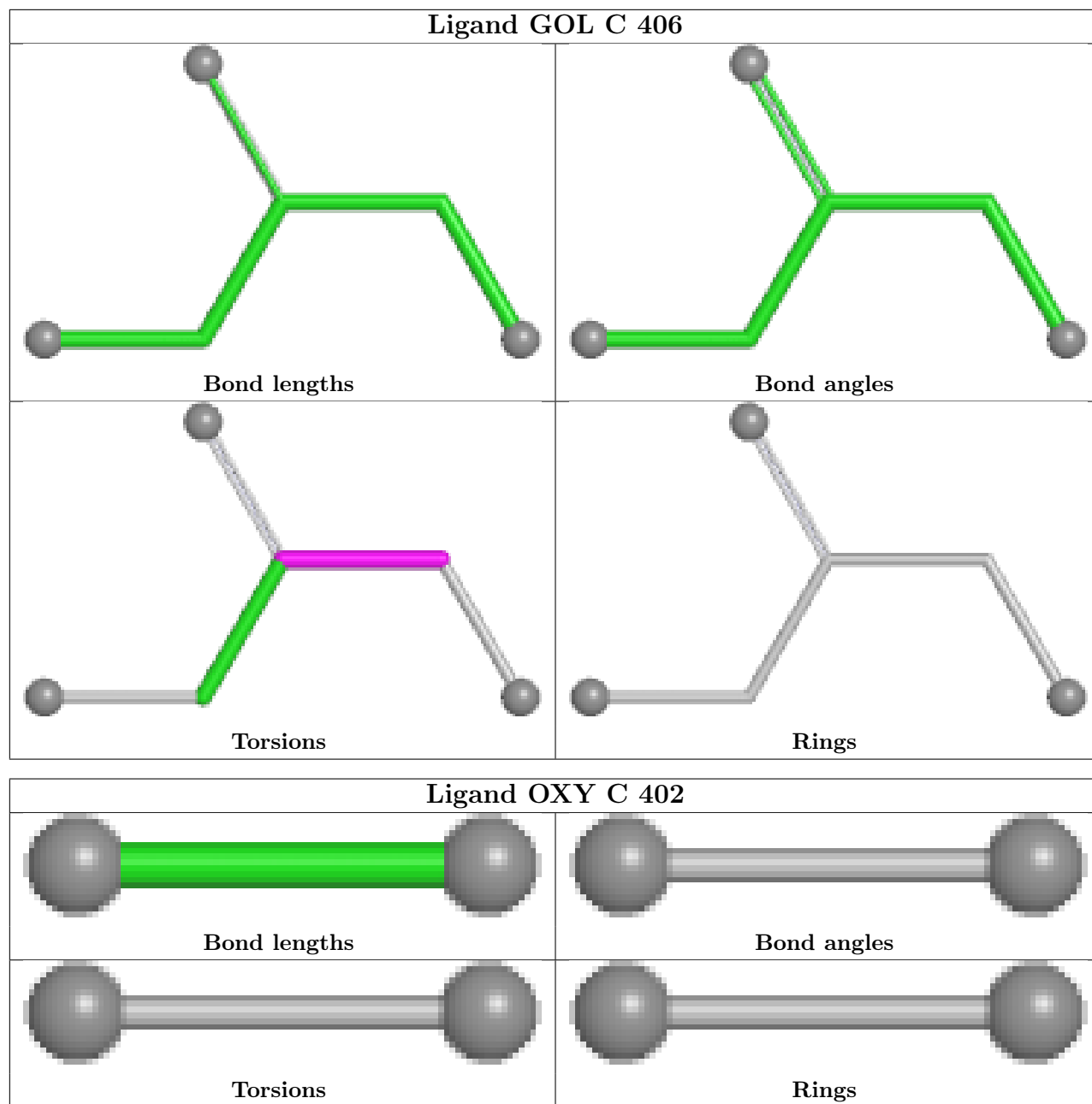












5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

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

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

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th 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 ²)	Q<0.9
1	A	363/363 (100%)	-0.10	5 (1%) 73 75	27, 42, 75, 139	0
1	B	361/363 (99%)	-0.11	4 (1%) 78 80	27, 43, 73, 110	0
1	C	363/363 (100%)	-0.09	7 (1%) 66 69	28, 42, 74, 160	0
1	D	361/363 (99%)	-0.14	5 (1%) 73 75	20, 43, 71, 122	1 (0%)
1	E	362/363 (99%)	0.05	3 (0%) 82 84	23, 48, 74, 138	2 (0%)
1	F	360/363 (99%)	0.22	9 (2%) 58 61	23, 49, 80, 167	2 (0%)
All	All	2170/2178 (99%)	-0.03	33 (1%) 72 74	20, 44, 75, 167	5 (0%)

The worst 5 of 33 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	1	MET	7.8
1	A	61	ALA	4.7
1	E	90	PHE	4.1
1	C	90	PHE	4.1
1	A	2	GLU	3.4

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

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, 95th 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.

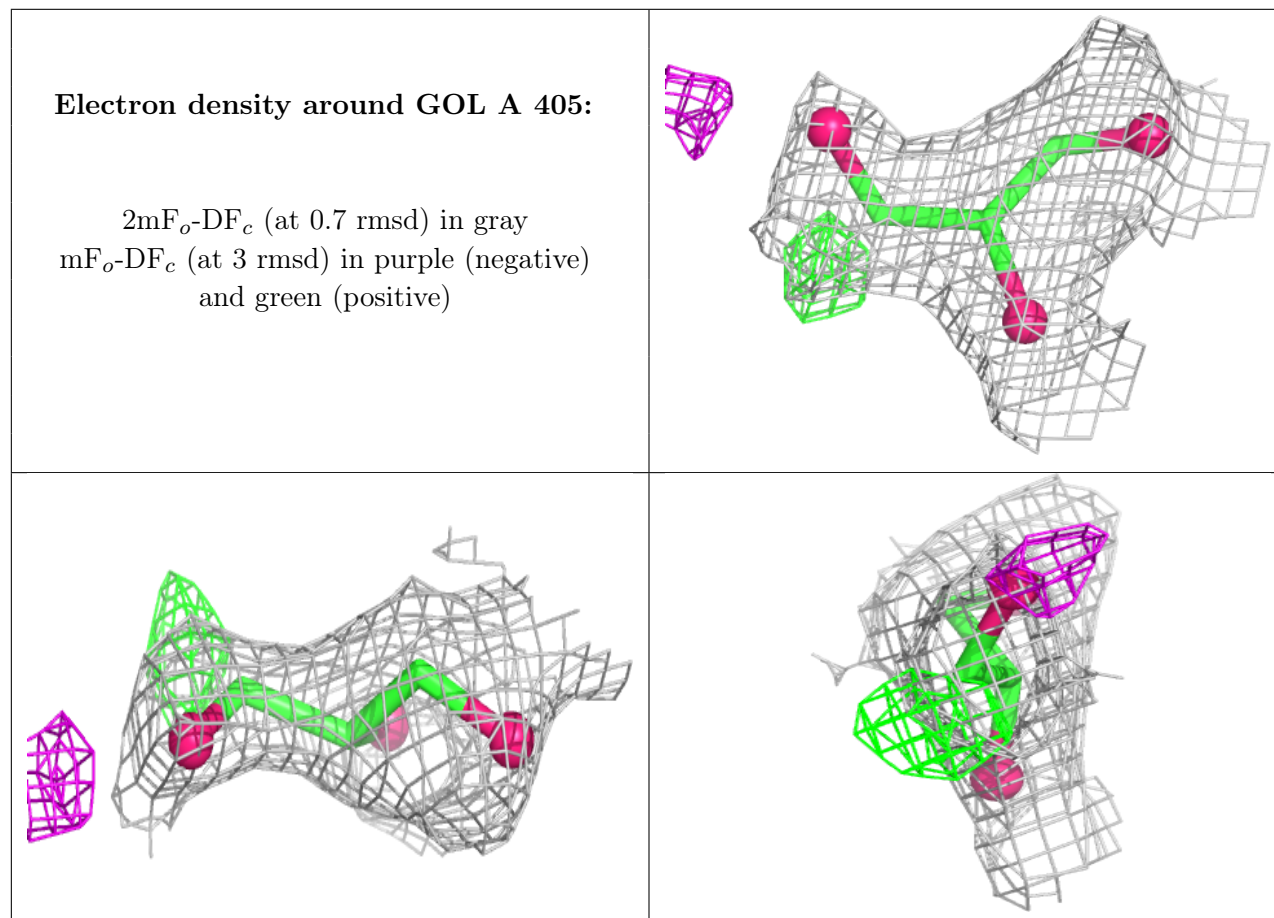
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
5	GOL	A	405	6/6	0.71	0.16	63,76,85,87	0
8	VOH	E	403	12/12	0.82	0.20	30,43,53,55	12
5	GOL	C	406	6/6	0.83	0.14	65,69,74,75	0
7	PO4	A	408	5/5	0.84	0.15	88,98,110,111	0
5	GOL	E	405	6/6	0.84	0.15	67,73,83,85	0
5	GOL	D	403	6/6	0.85	0.16	54,71,79,81	0
5	GOL	F	406	6/6	0.86	0.09	87,98,101,103	0
9	EPE	F	403	15/15	0.86	0.13	81,93,113,117	0
5	GOL	F	405	6/6	0.87	0.12	81,86,94,96	0
8	VOH	C	403	12/12	0.88	0.18	45,59,71,72	12
4	MPD	C	404	8/8	0.89	0.17	49,67,79,83	0
5	GOL	F	404	6/6	0.90	0.09	73,76,84,85	0
5	GOL	B	403	6/6	0.92	0.16	66,76,97,102	0
5	GOL	E	404	6/6	0.92	0.14	70,77,81,84	6
5	GOL	A	406	6/6	0.92	0.13	60,71,75,82	0
5	GOL	F	407	6/6	0.92	0.15	45,48,66,67	0
6	CL	A	407	1/1	0.93	0.12	75,75,75,75	0
6	CL	B	404	1/1	0.93	0.10	77,77,77,77	0
3	OXY	F	402	2/2	0.93	0.13	58,58,58,61	0
3	OXY	C	402	2/2	0.93	0.11	44,44,44,49	0
5	GOL	A	404	6/6	0.93	0.10	69,88,94,96	0
5	GOL	C	405	6/6	0.93	0.12	25,36,41,43	6
3	OXY	B	402	2/2	0.94	0.10	52,52,52,65	0
4	MPD	A	403	8/8	0.94	0.15	67,94,101,107	0
2	HEM	F	401	43/43	0.95	0.08	35,44,59,65	0
6	CL	F	408	1/1	0.95	0.13	89,89,89,89	0
3	OXY	D	402	2/2	0.96	0.10	45,45,45,58	0
2	HEM	E	401	43/43	0.96	0.07	34,41,48,51	1
2	HEM	B	401	43/43	0.96	0.07	26,36,40,41	1
6	CL	C	407	1/1	0.97	0.07	70,70,70,70	0
6	CL	E	406	1/1	0.97	0.11	65,65,65,65	0
2	HEM	C	401	43/43	0.97	0.06	24,32,37,39	1
6	CL	F	409	1/1	0.97	0.08	64,64,64,64	0
3	OXY	E	402	2/2	0.97	0.07	55,55,55,66	0
3	OXY	A	402	2/2	0.97	0.10	57,57,57,65	0
2	HEM	D	401	43/43	0.97	0.06	28,33,43,45	1
2	HEM	A	401	43/43	0.97	0.06	29,34,39,42	1

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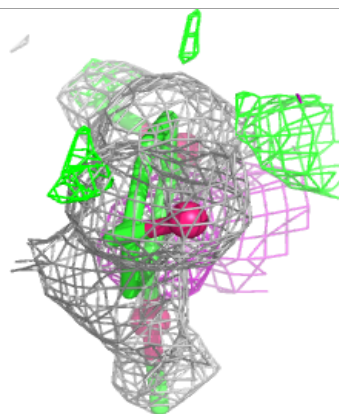
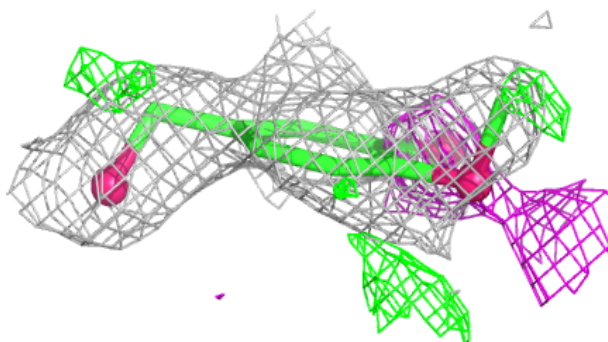
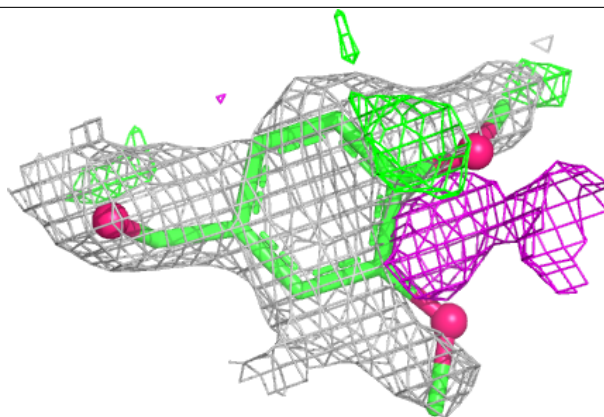
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
6	CL	D	405	1/1	0.98	0.15	78,78,78,78	0
6	CL	D	404	1/1	0.98	0.07	61,61,61,61	0

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.

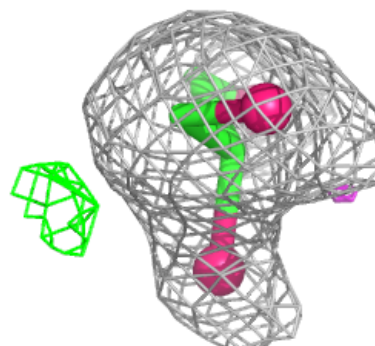
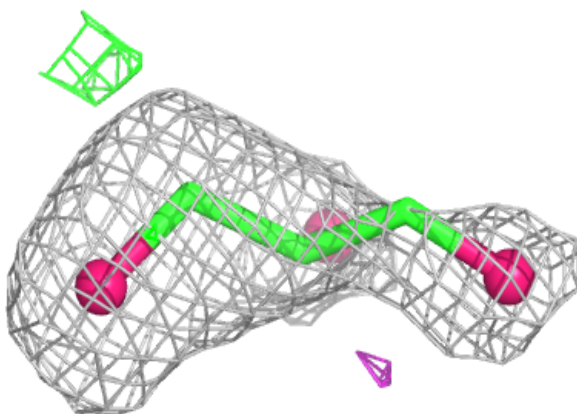
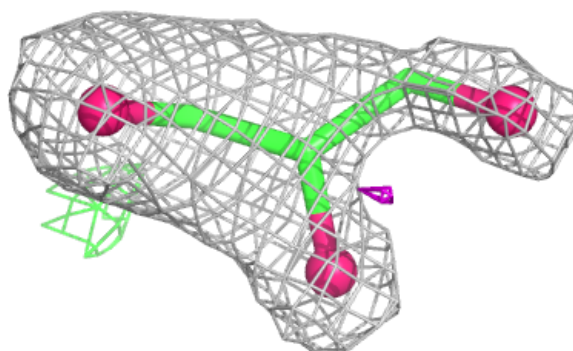


Electron density around VOH E 403:

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

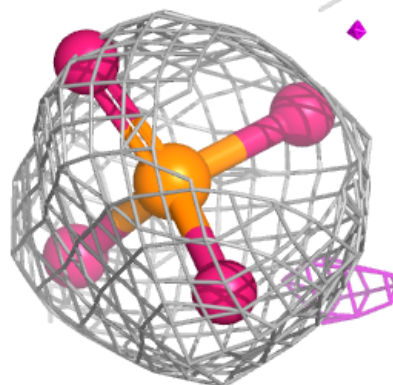
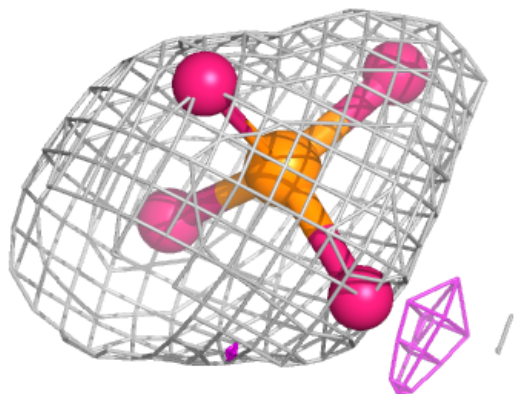
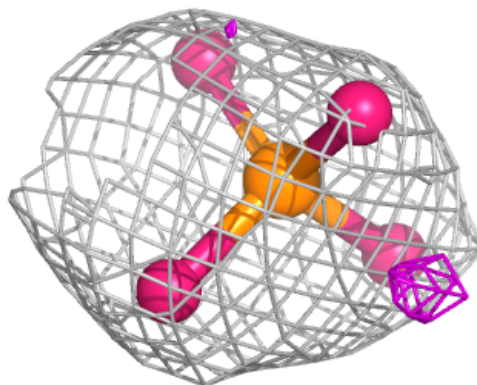
**Electron density around GOL C 406:**

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



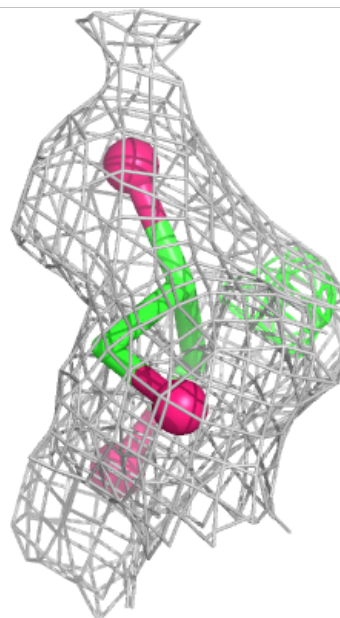
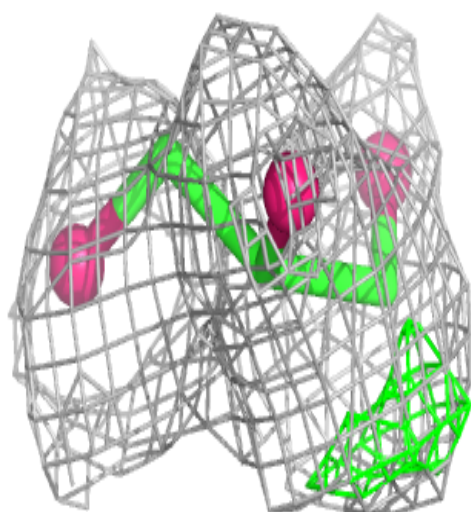
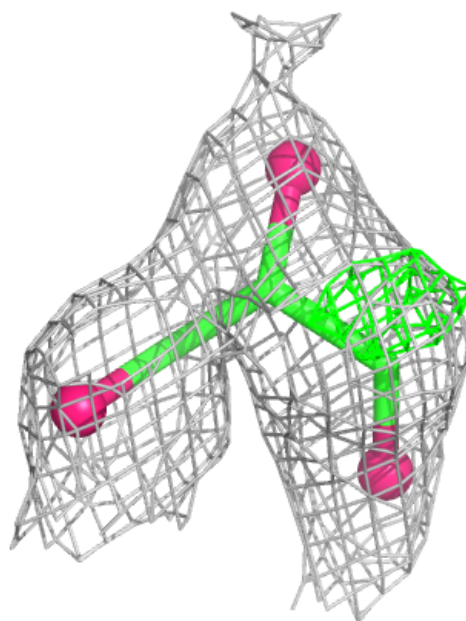
Electron density around PO4 A 408:

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



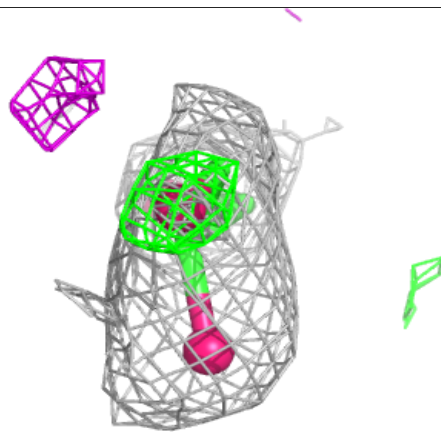
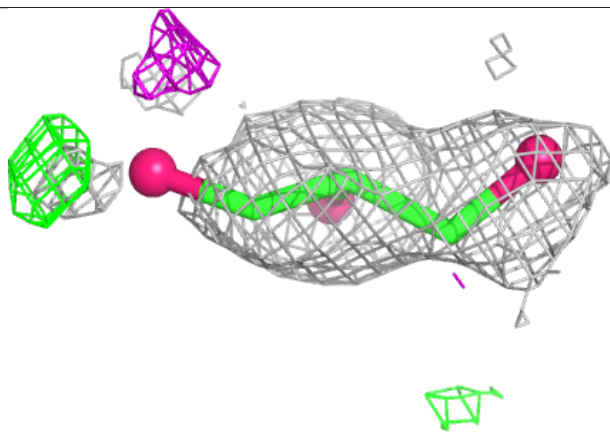
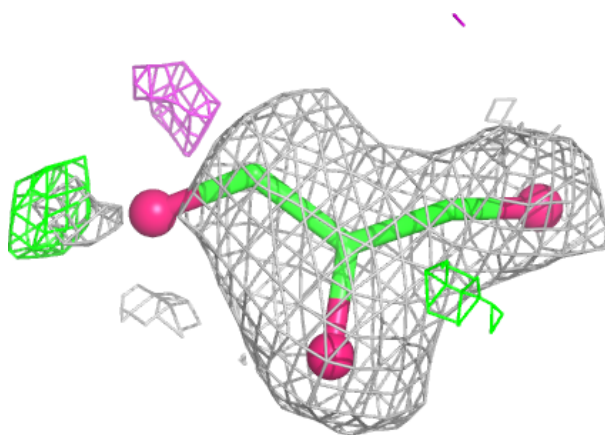
Electron density around GOL E 405:

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



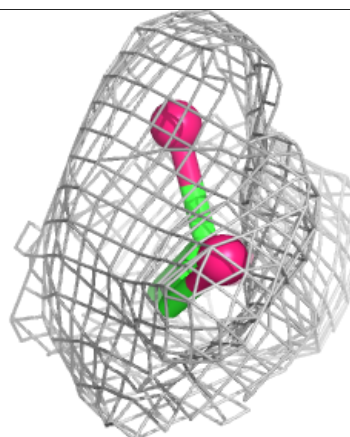
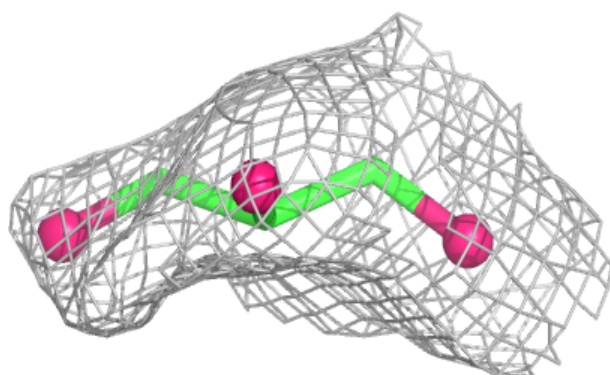
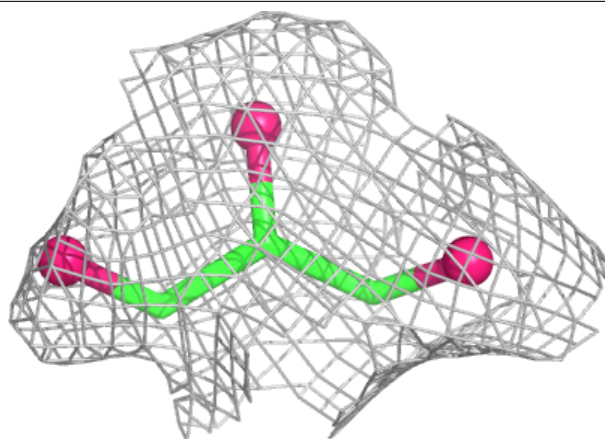
Electron density around GOL D 403:

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

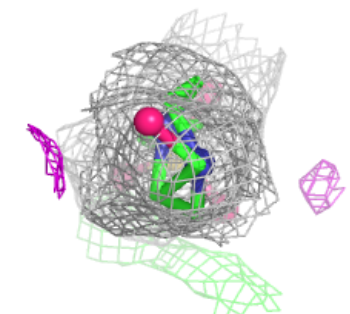
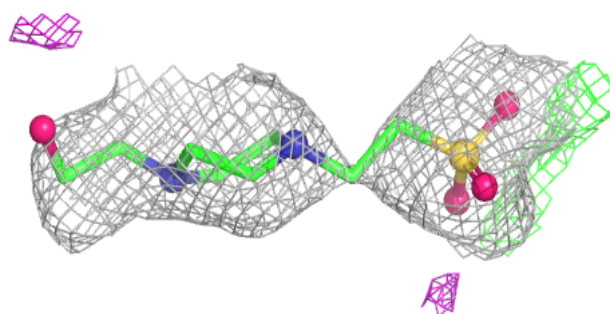
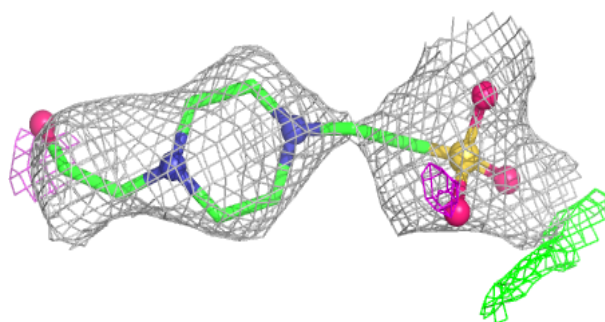


Electron density around GOL F 406:

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

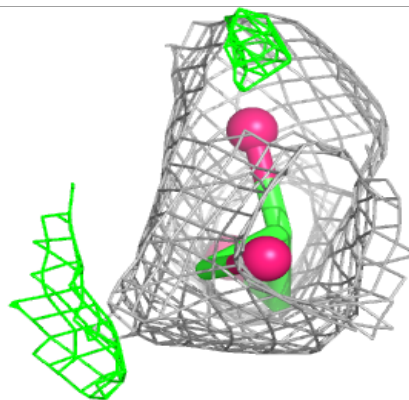
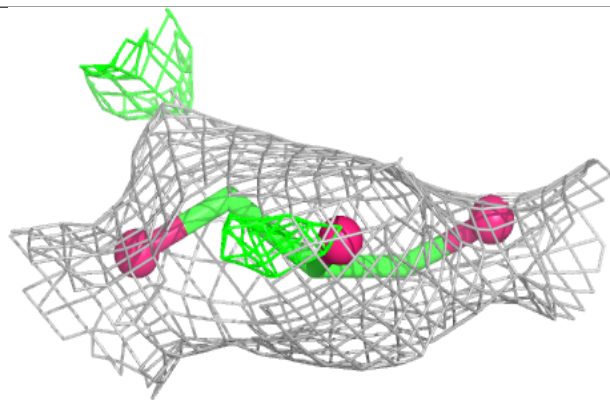
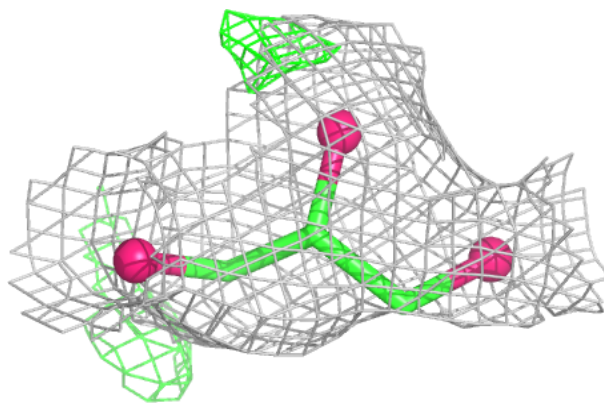
**Electron density around EPE F 403:**

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



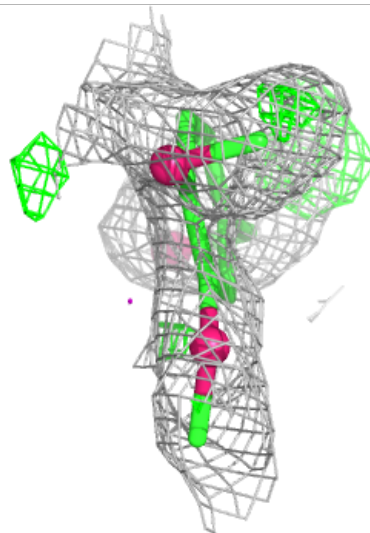
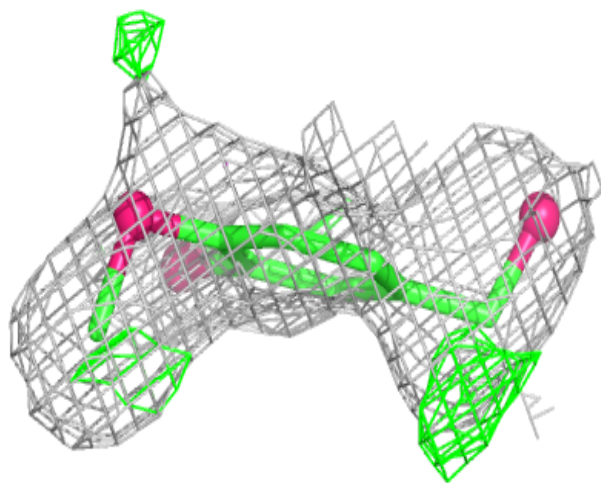
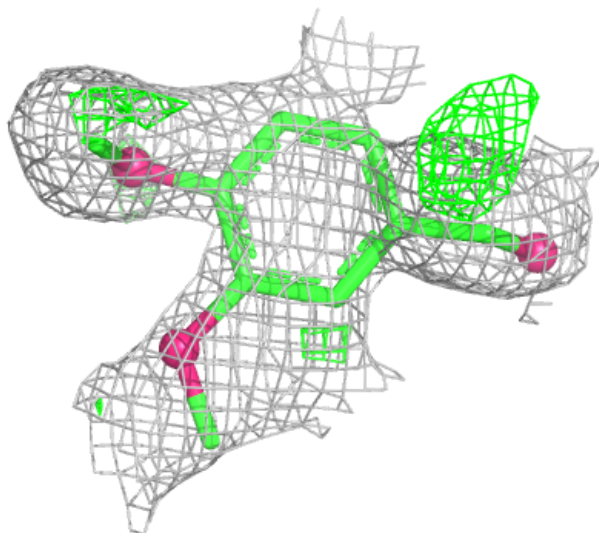
Electron density around GOL F 405:

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



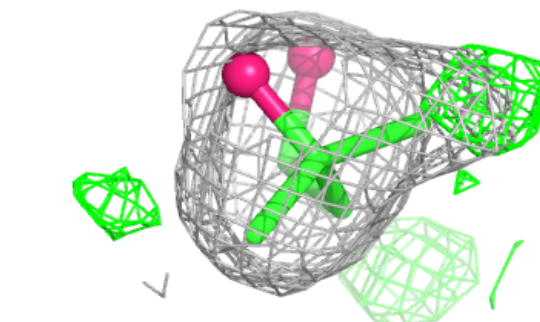
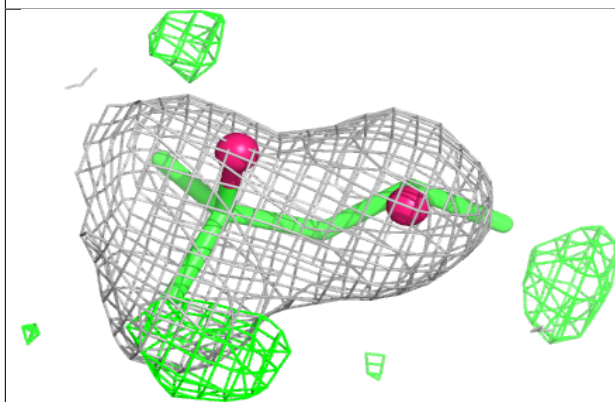
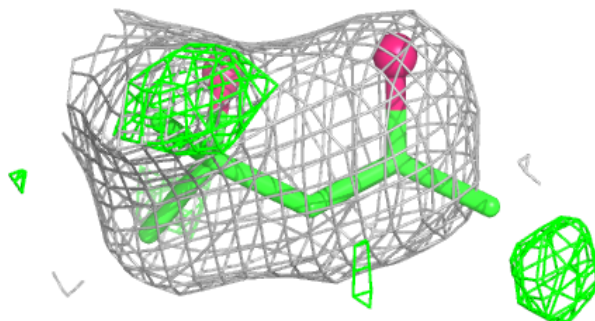
Electron density around VOH C 403:

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

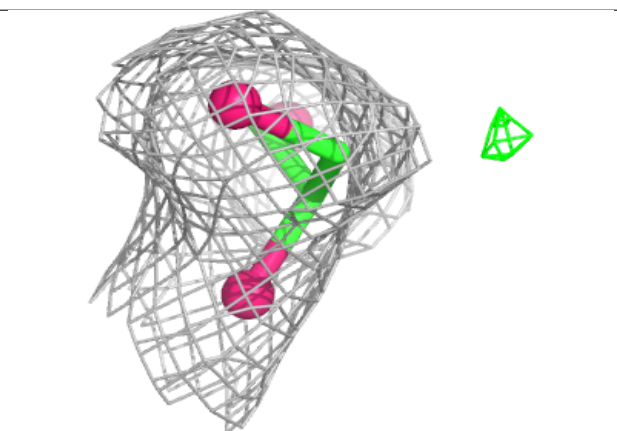
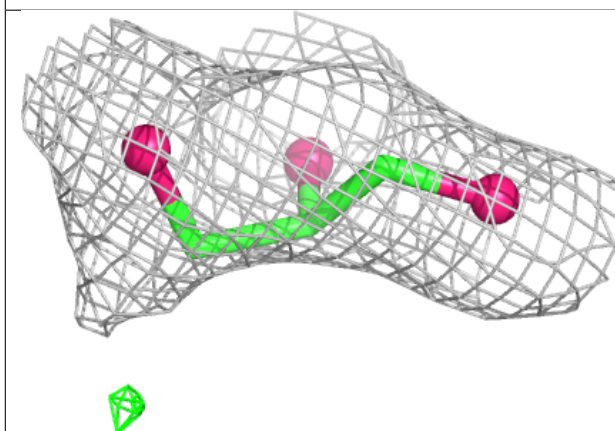
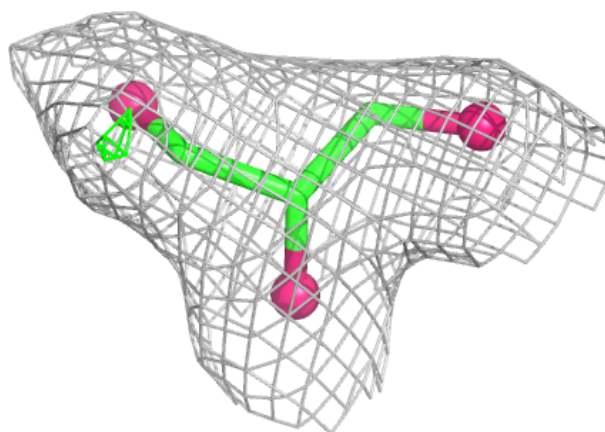


Electron density around MPD C 404:

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

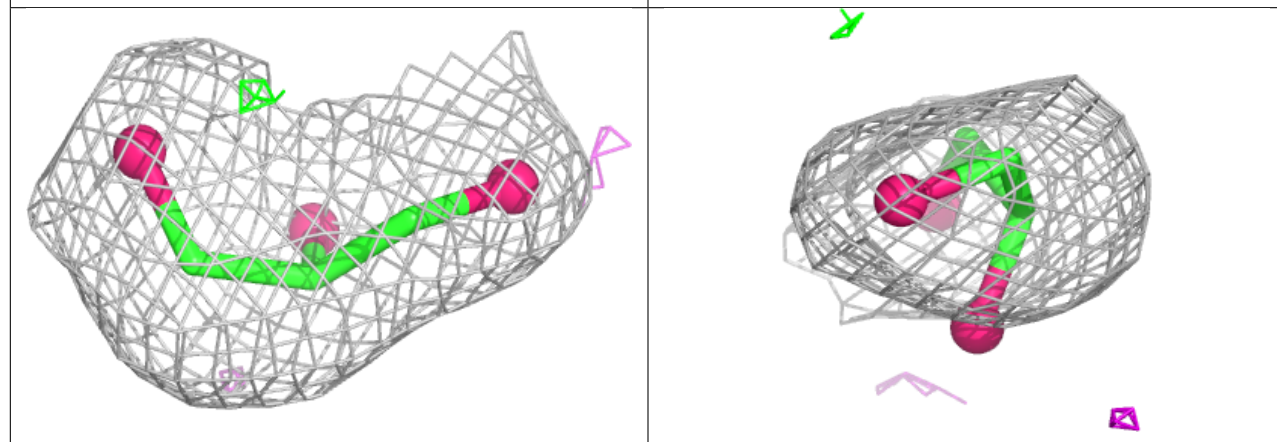
**Electron density around GOL F 404:**

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



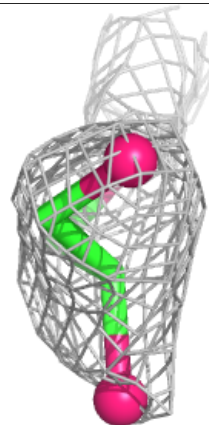
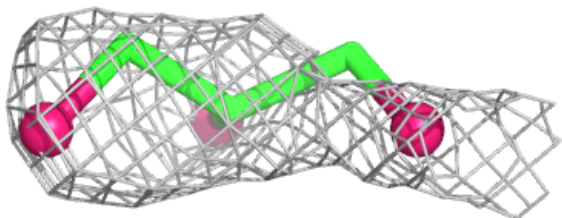
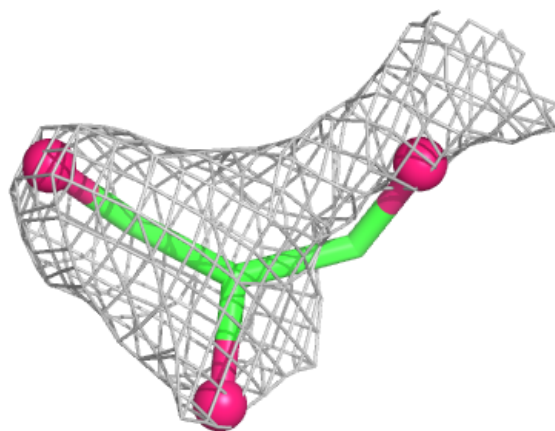
Electron density around GOL B 403:

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



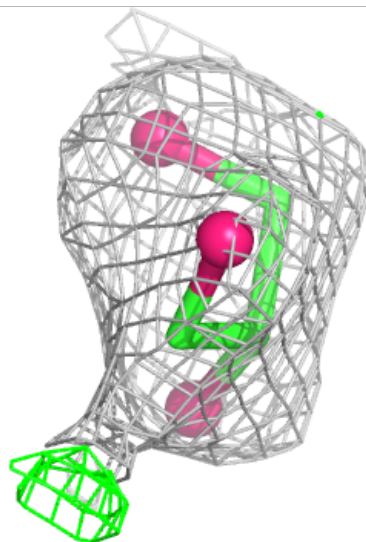
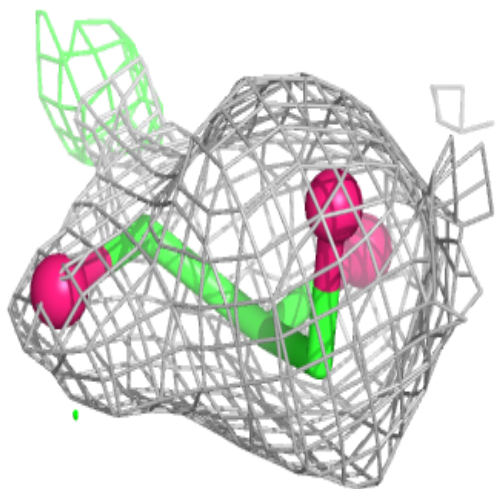
Electron density around GOL E 404:

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



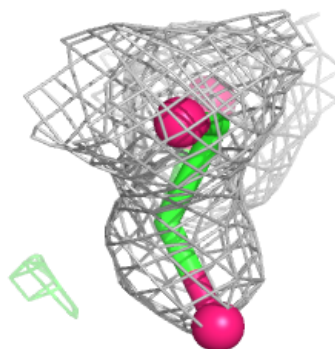
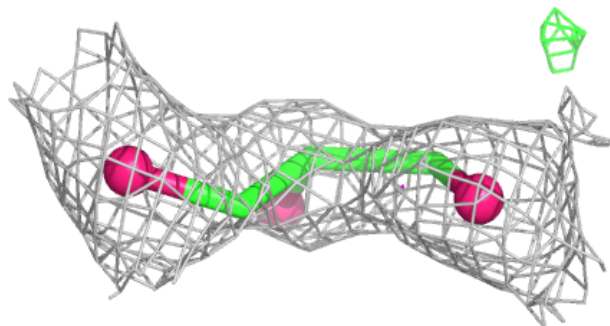
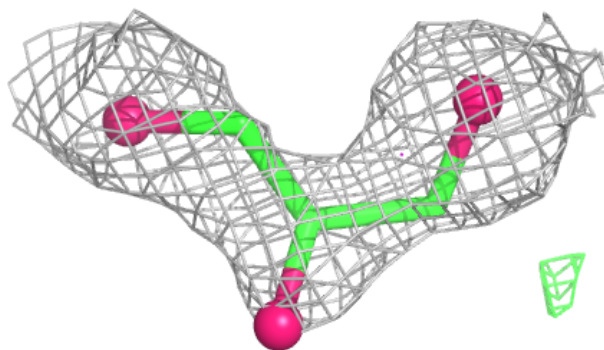
Electron density around GOL A 406:

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



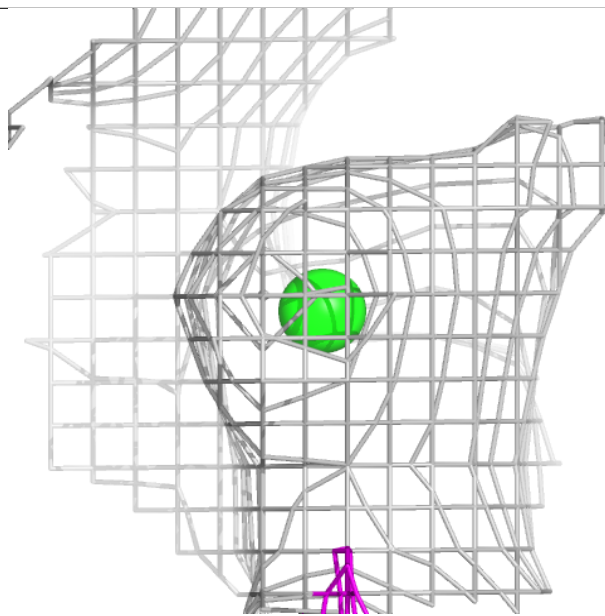
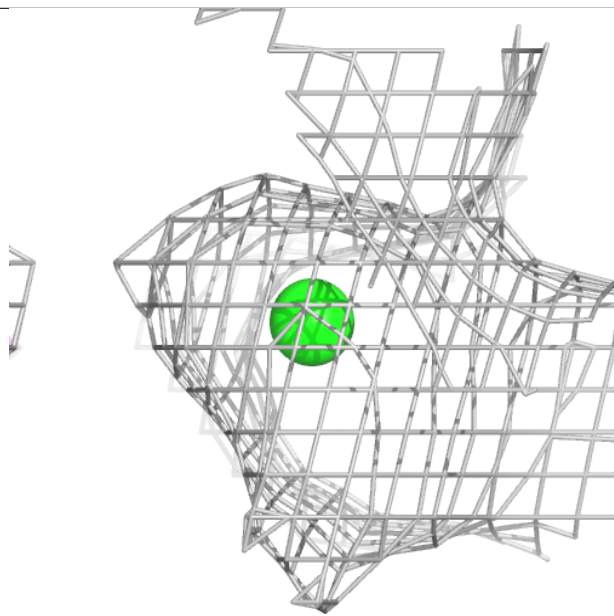
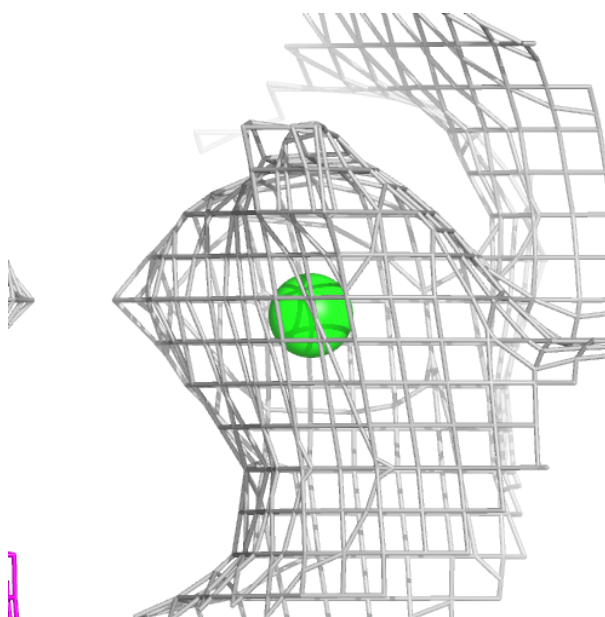
Electron density around GOL F 407:

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



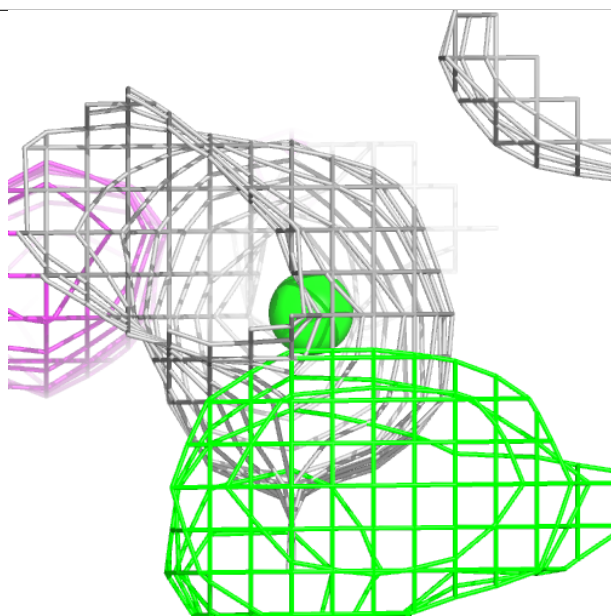
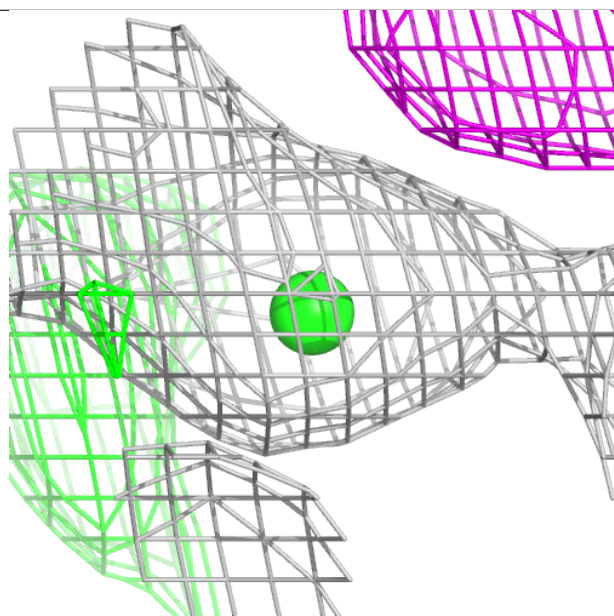
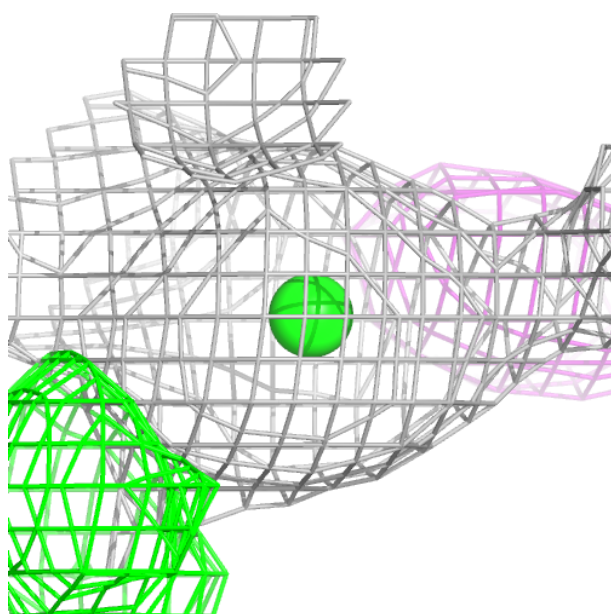
Electron density around CL A 407:

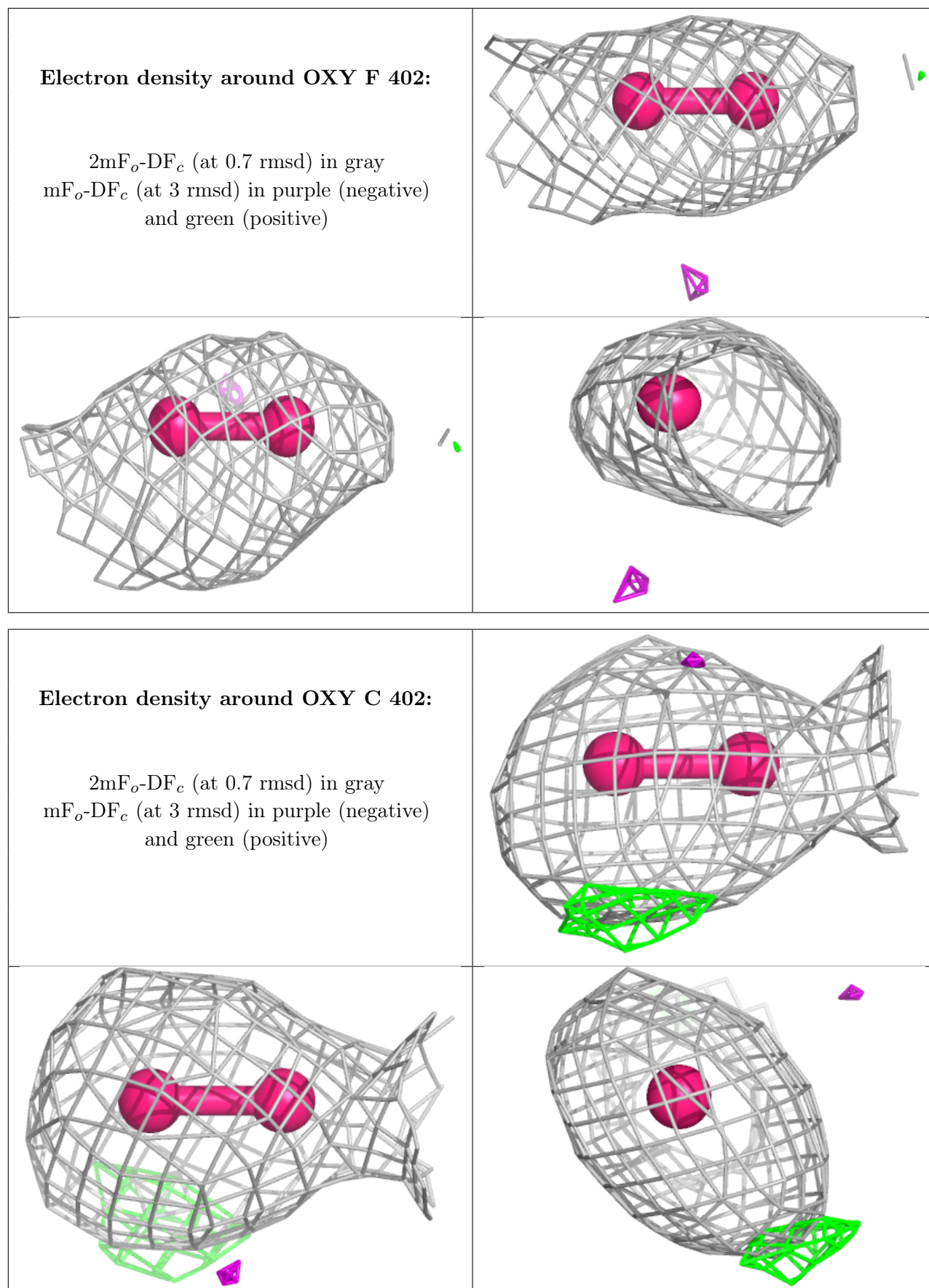
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around CL B 404:

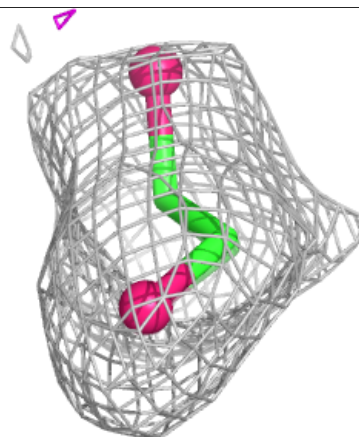
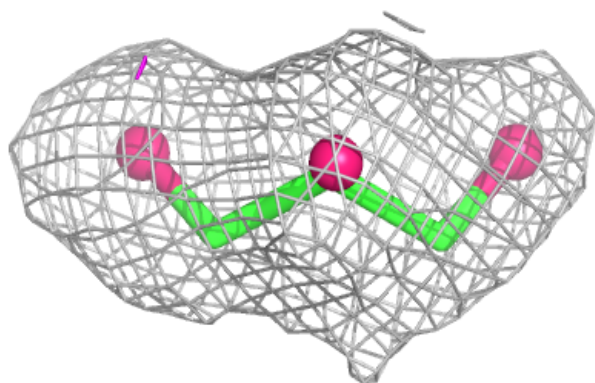
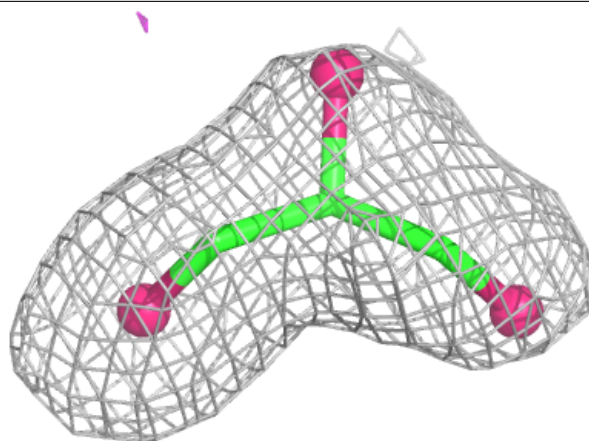
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)





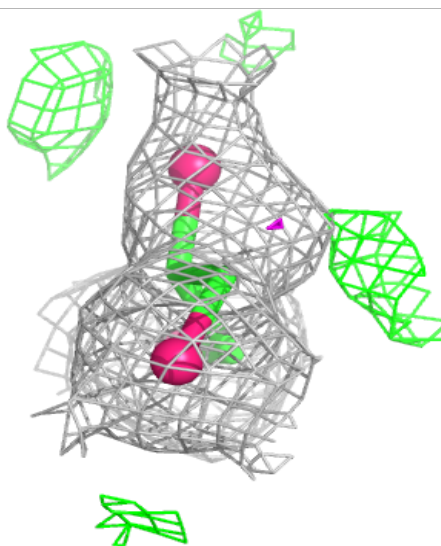
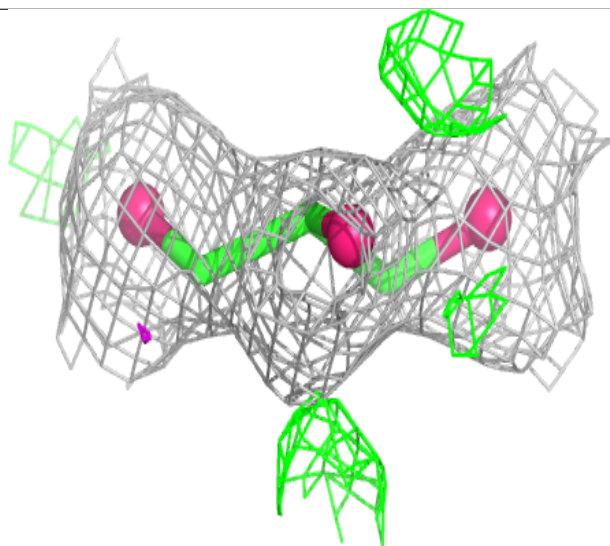
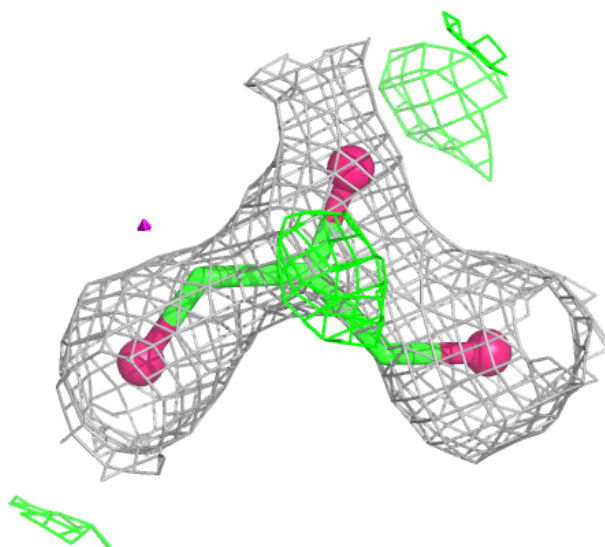
Electron density around GOL A 404:

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



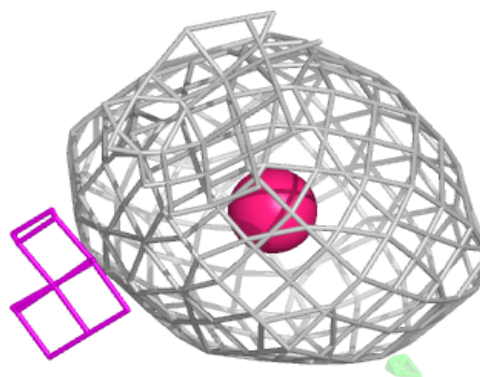
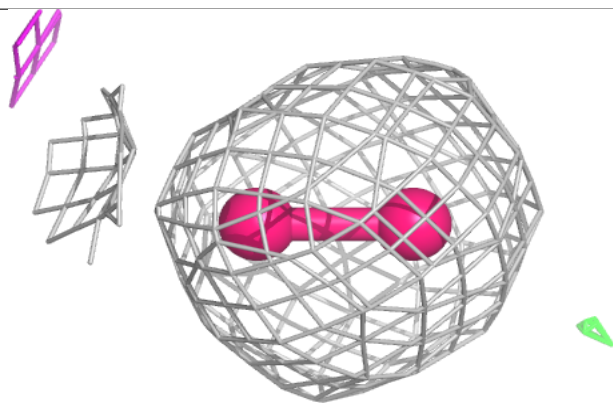
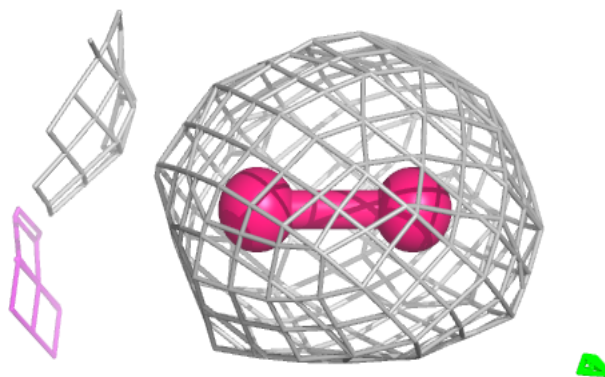
Electron density around GOL C 405:

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



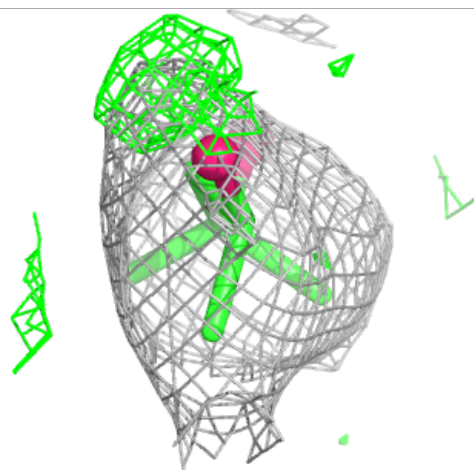
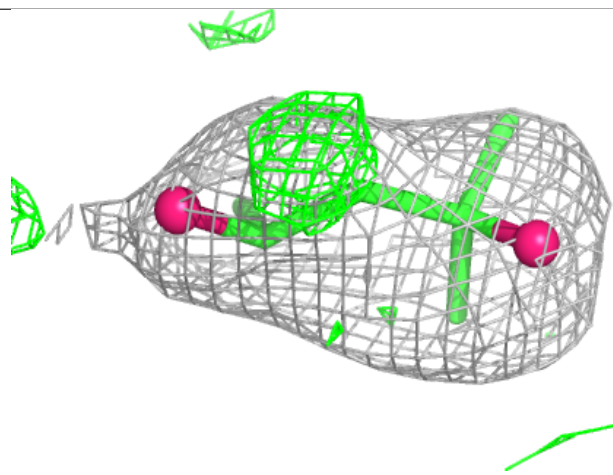
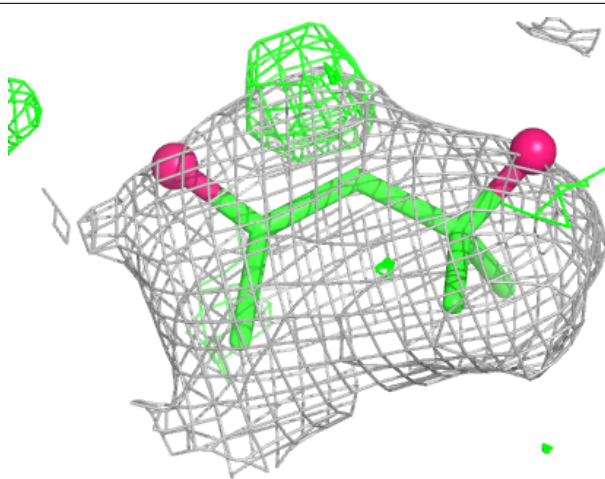
Electron density around OXY B 402:

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



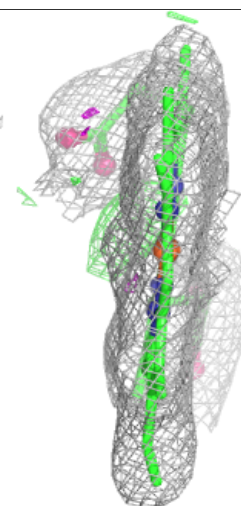
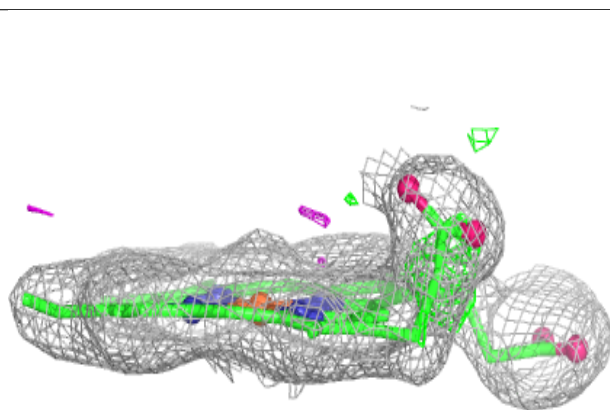
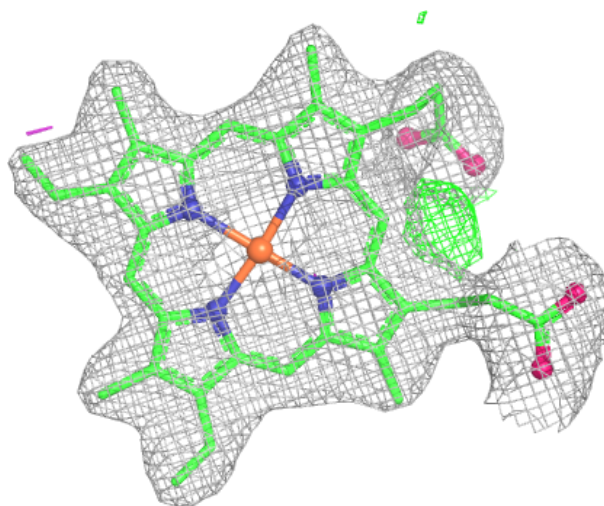
Electron density around MPD A 403:

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



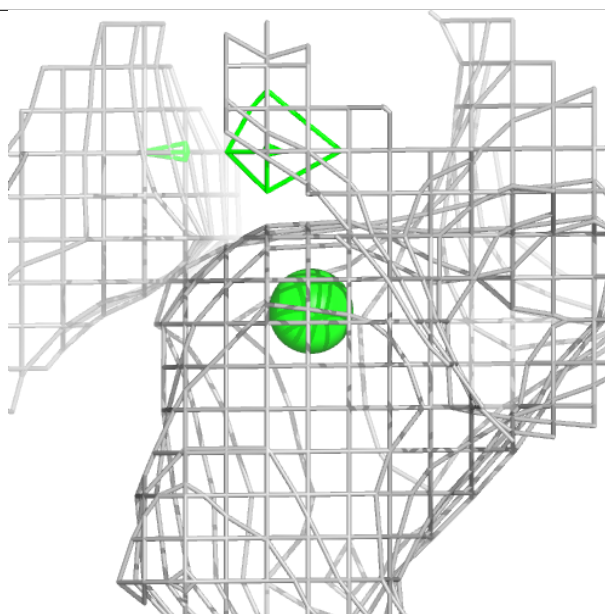
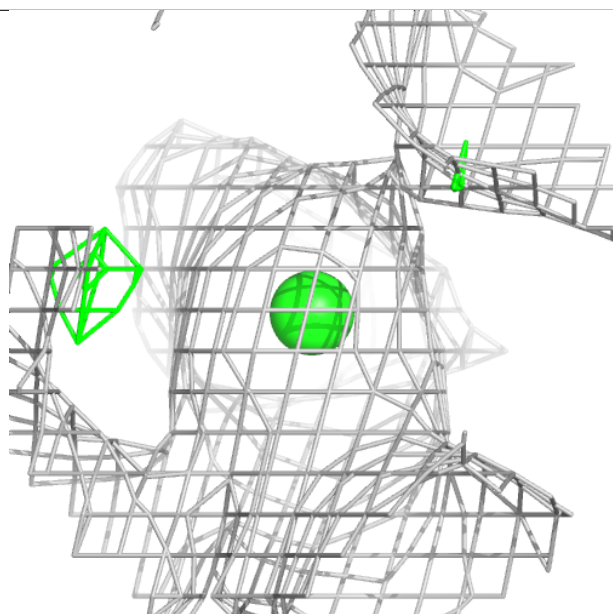
Electron density around HEM F 401:

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



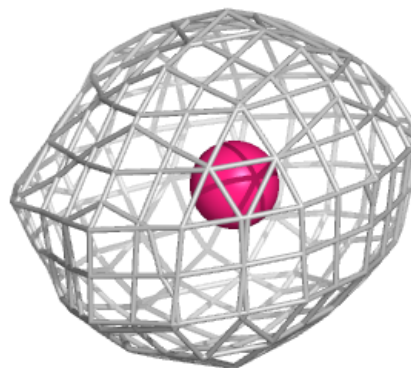
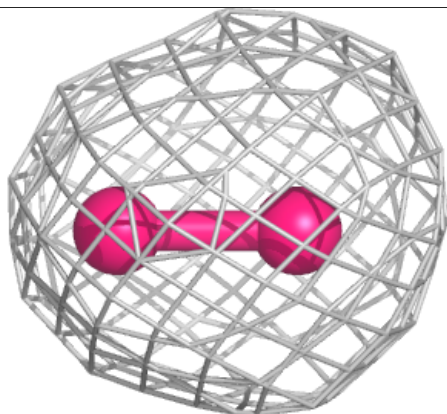
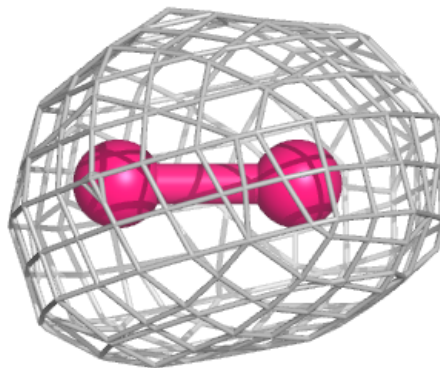
Electron density around CL F 408:

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



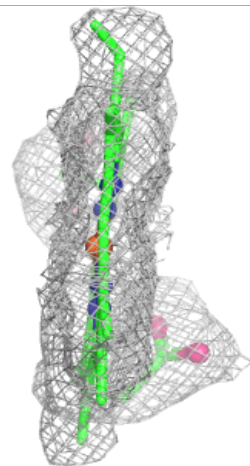
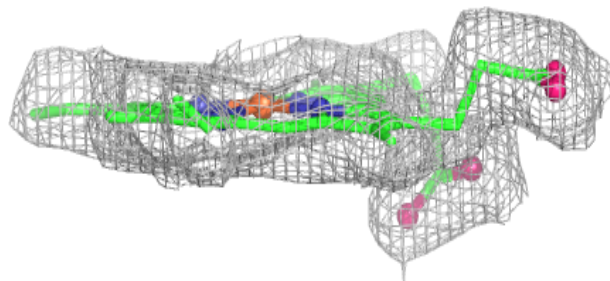
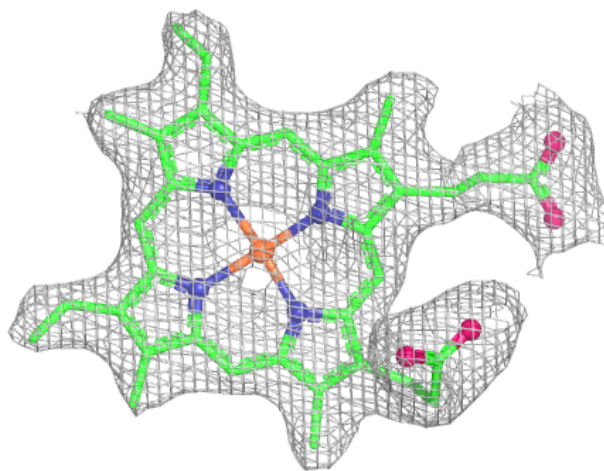
Electron density around OXY D 402:

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



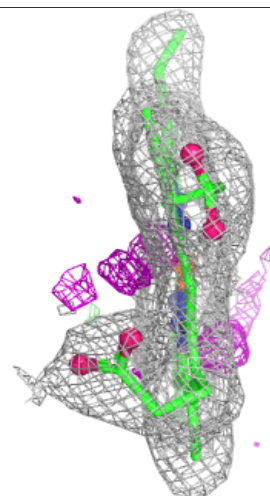
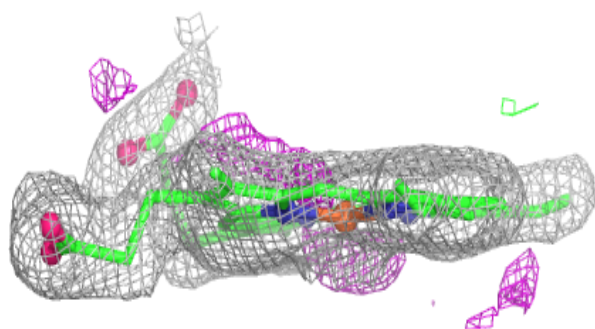
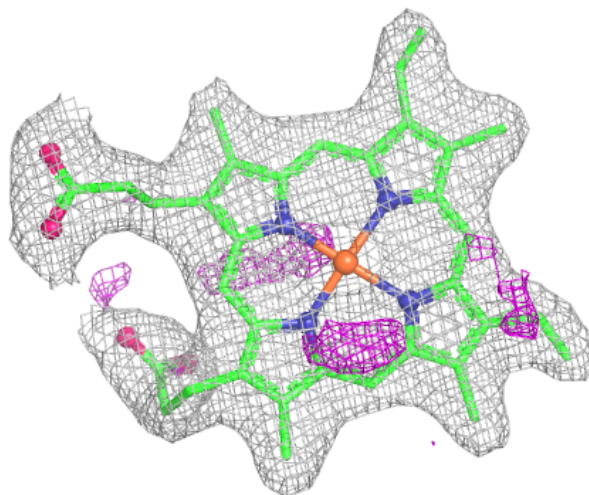
Electron density around HEM E 401:

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



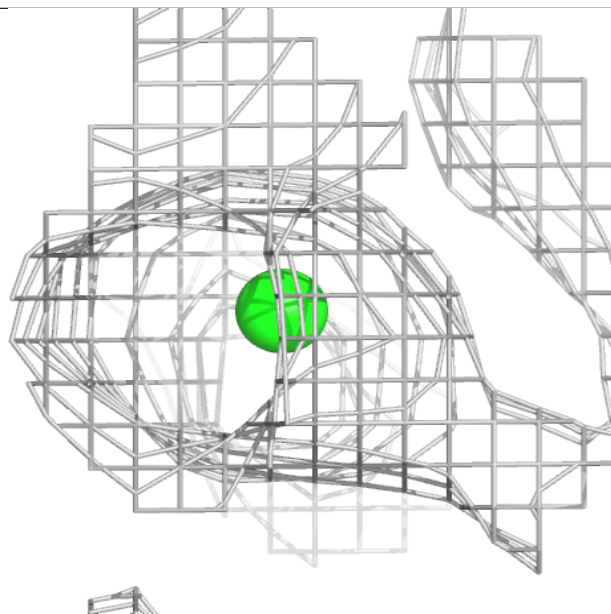
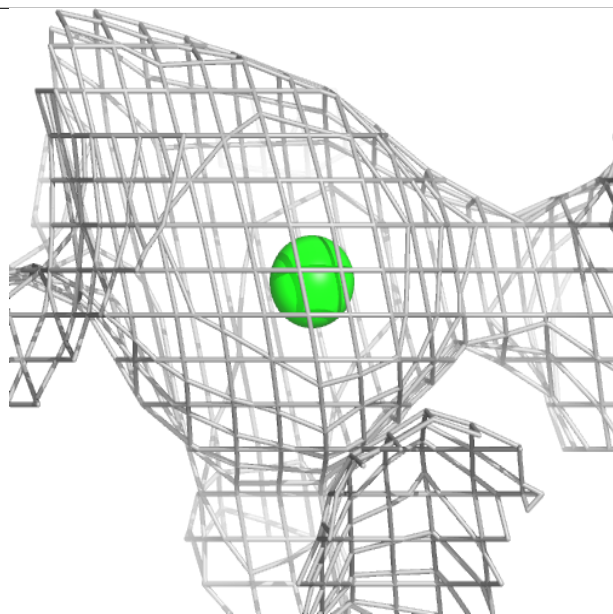
Electron density around HEM B 401:

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



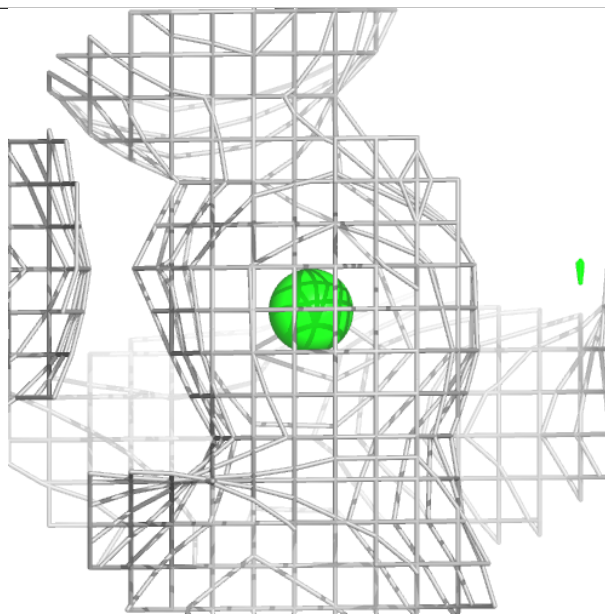
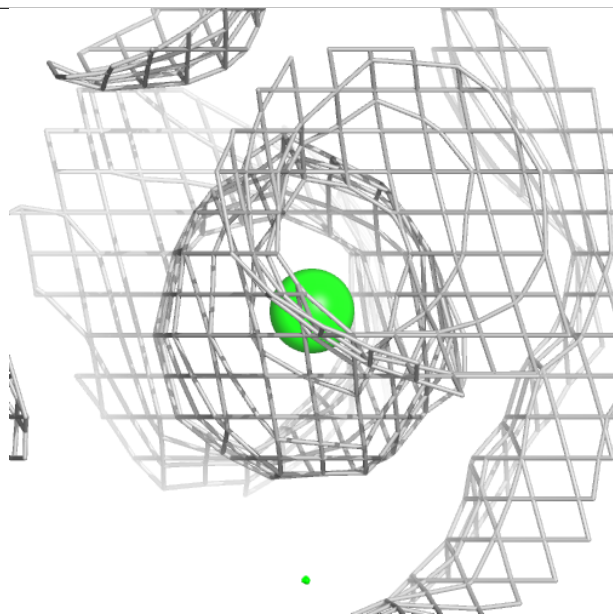
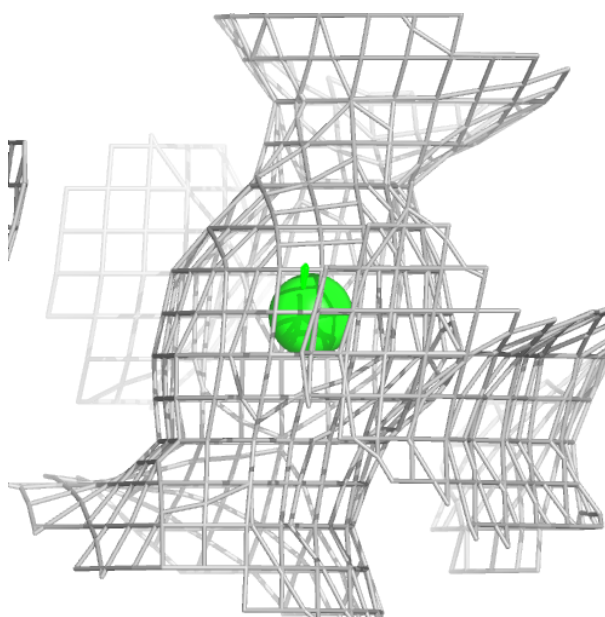
Electron density around CL C 407:

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



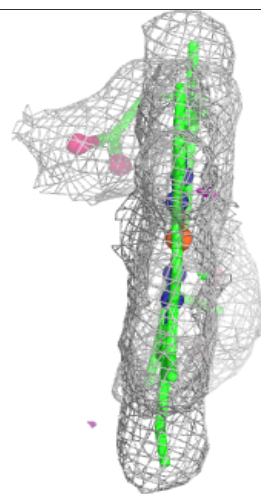
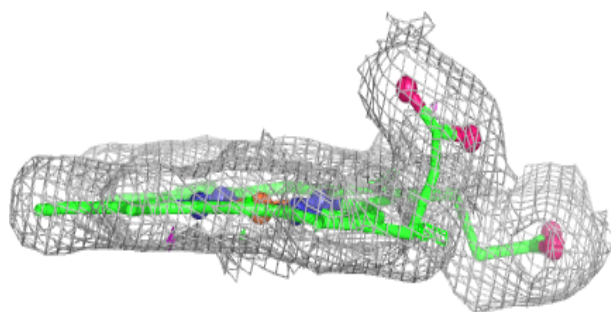
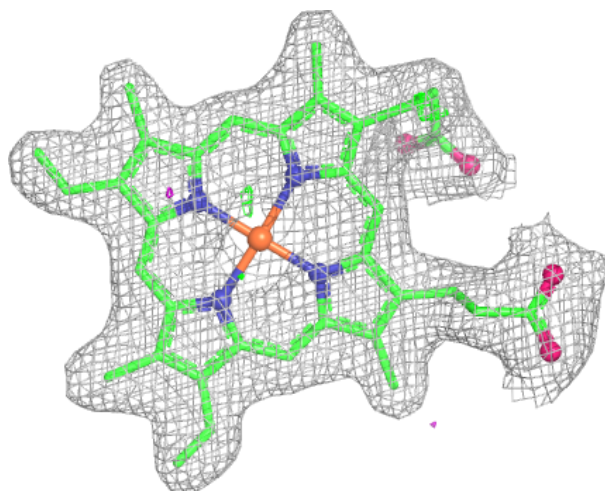
Electron density around CL E 406:

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



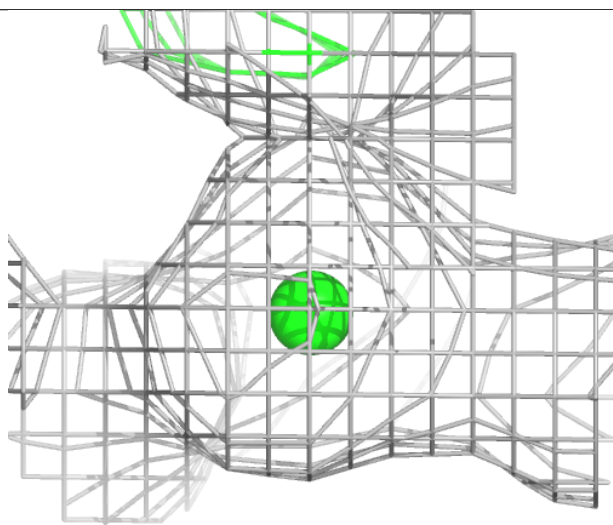
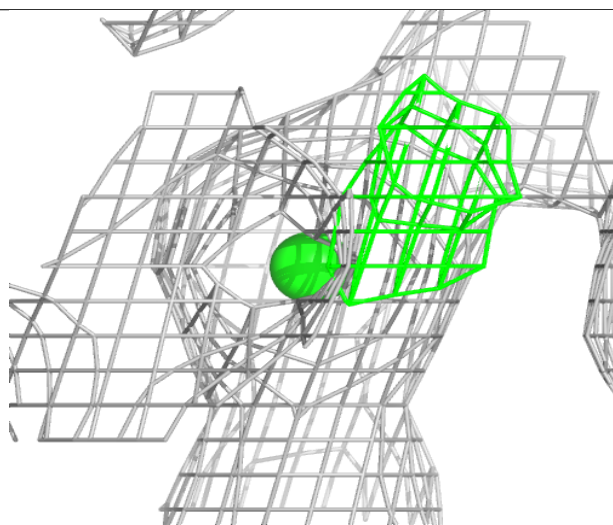
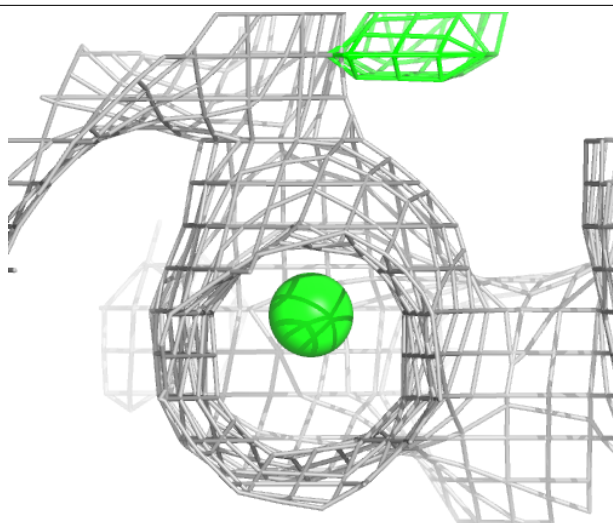
Electron density around HEM C 401:

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



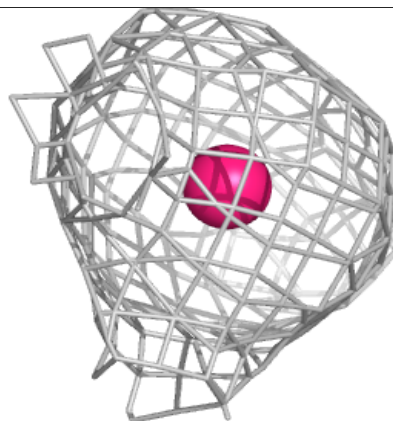
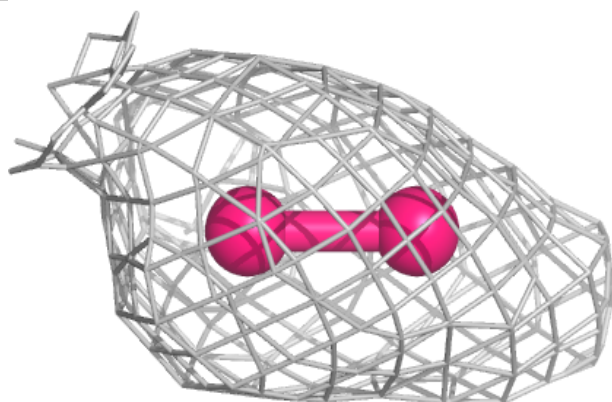
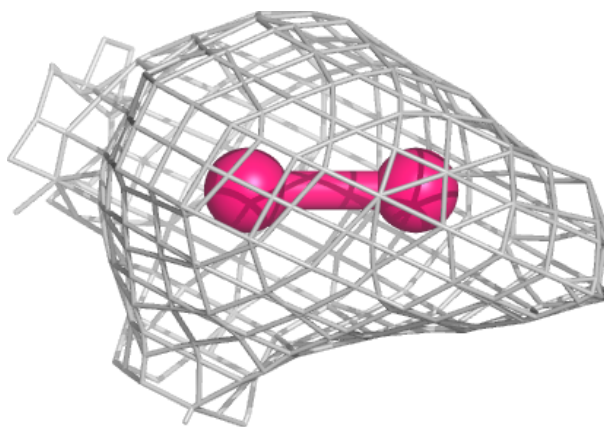
Electron density around CL F 409:

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



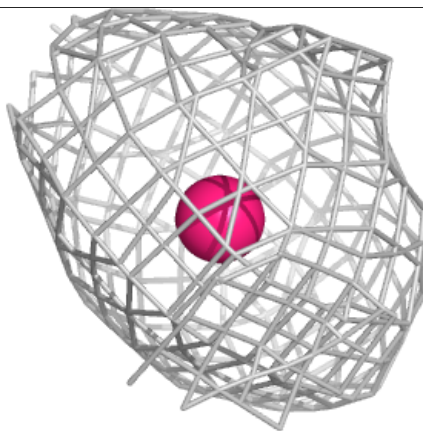
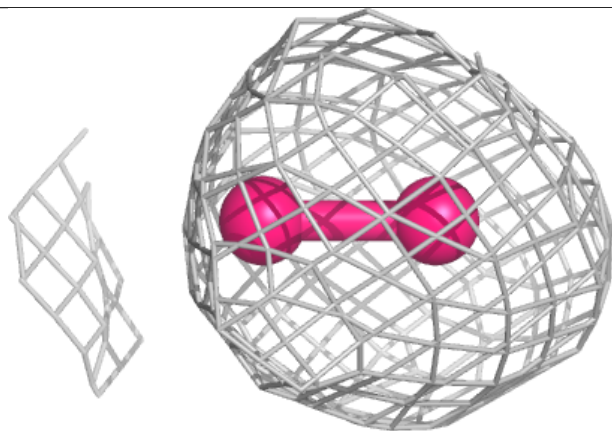
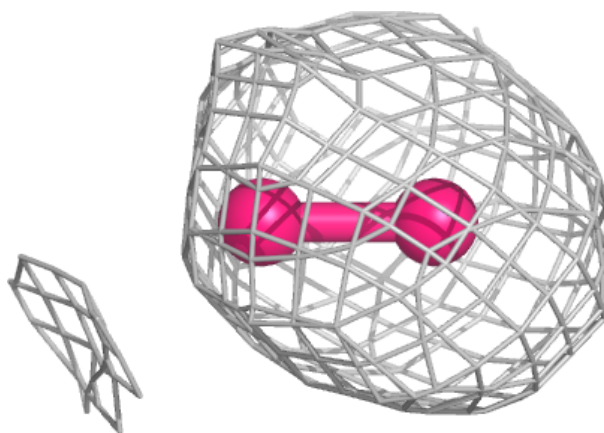
Electron density around OXY E 402:

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



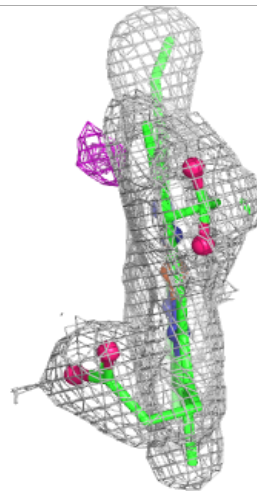
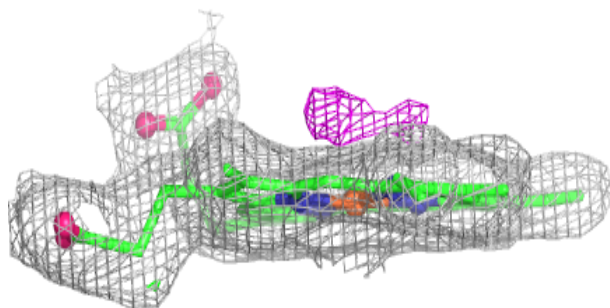
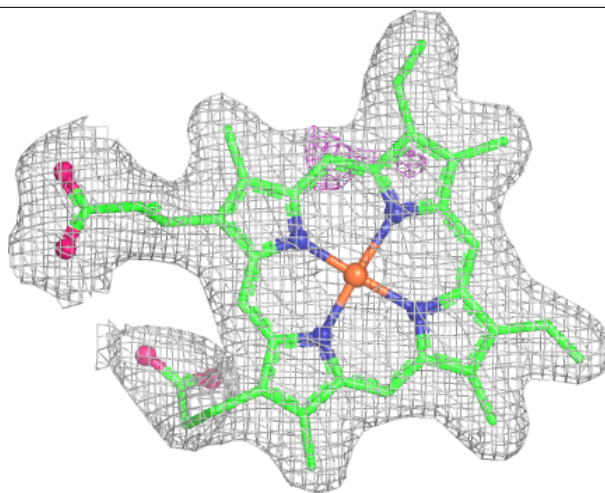
Electron density around OXY A 402:

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



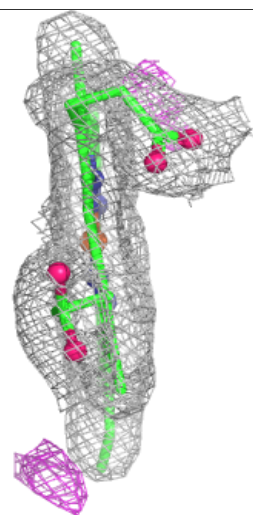
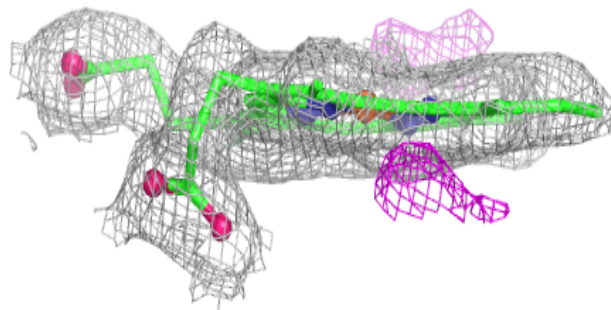
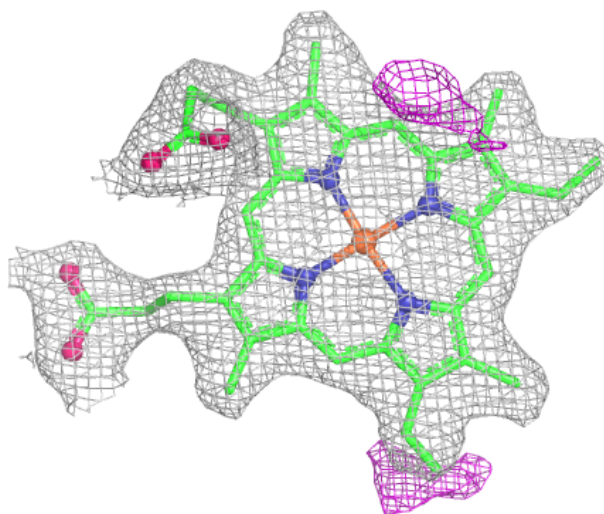
Electron density around HEM D 401:

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



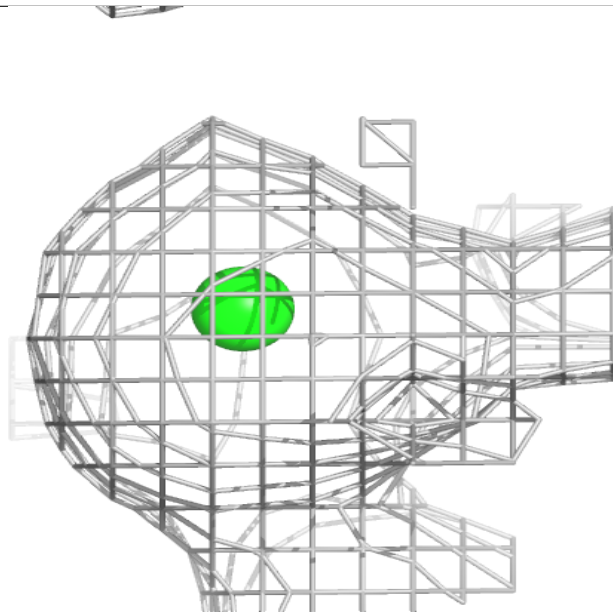
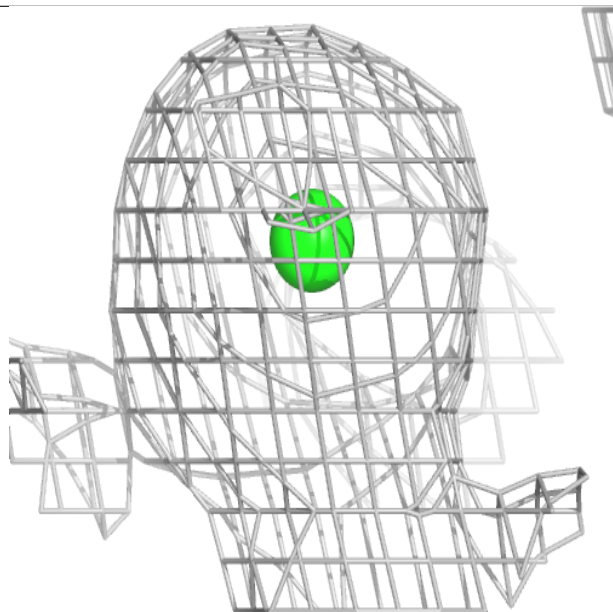
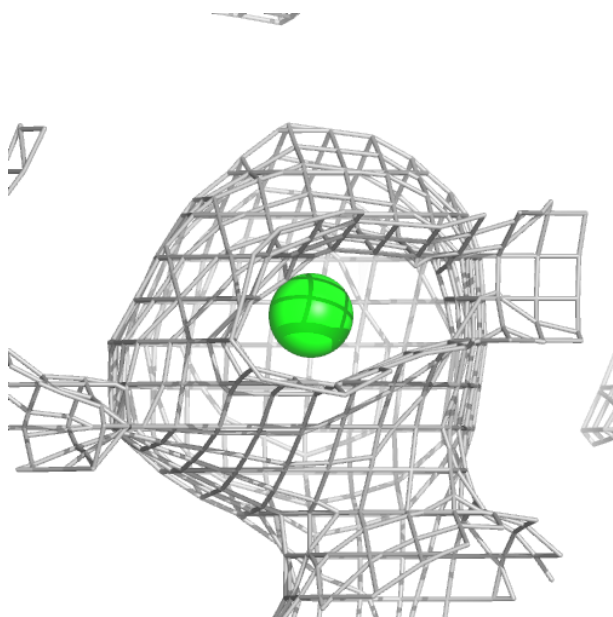
Electron density around HEM A 401:

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



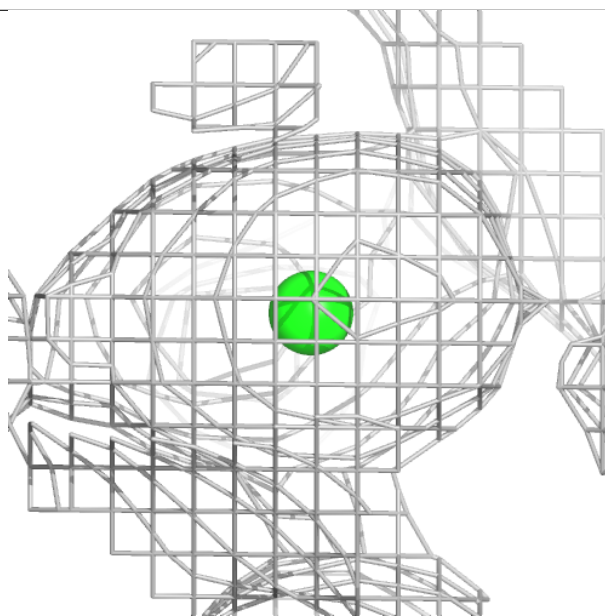
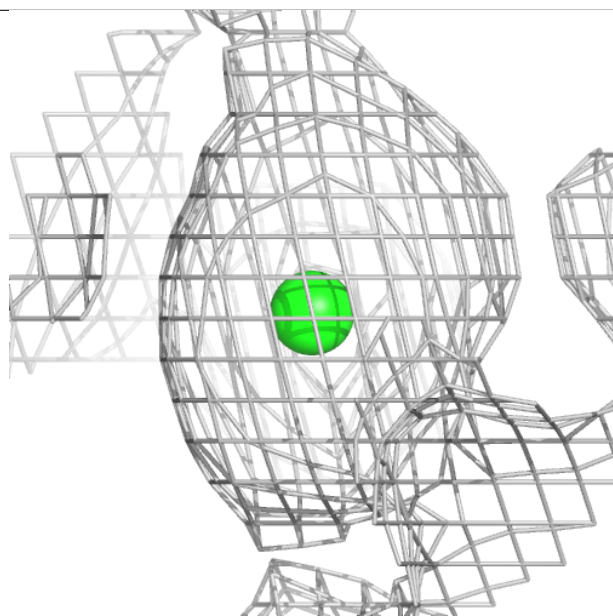
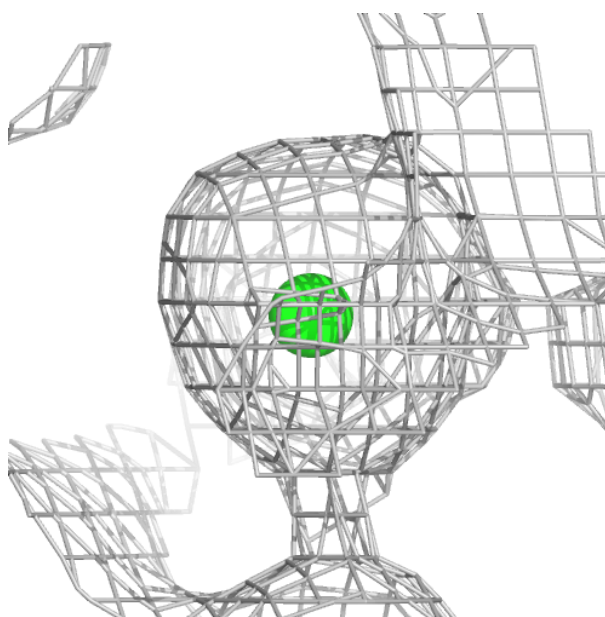
Electron density around CL D 405:

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



Electron density around CL D 404:

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



6.5 Other polymers ⓘ

There are no such residues in this entry.