



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 01:33 PM UTC

PDB ID : 9TDX / pdb_00009tdx
Title : Structure of D81A-Fructofuranosidase from *Purpureocillum lilacinum* in complex with sucrose
Authors : Ruiz-Nunez, M.; Sanz-Aparicio, J.
Deposited on : 2025-11-24
Resolution : 2.00 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

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:

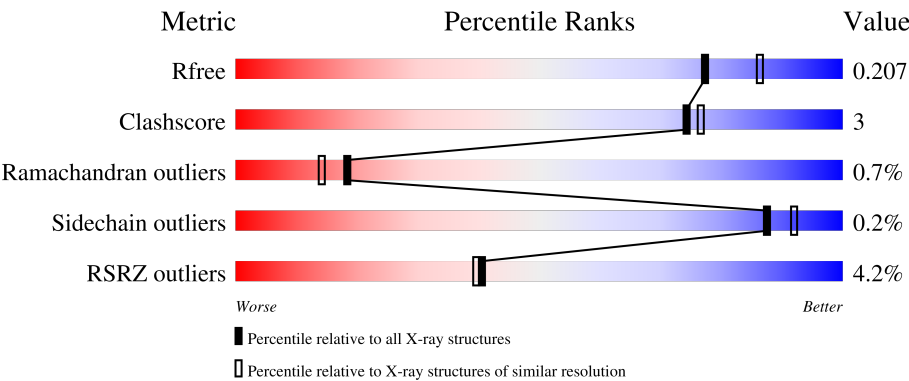
MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal 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 i

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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.



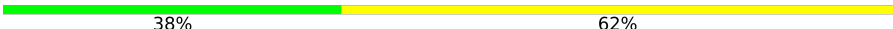
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R _{free}	180053	10052 (2.00-2.00)
Clashscore	190562	11152 (2.00-2.00)
Ramachandran outliers	187476	11031 (2.00-2.00)
Sidechain outliers	187428	11029 (2.00-2.00)
RSRZ outliers	180081	10067 (2.00-2.00)

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	576	3% 86% 7% 7%
1	B	576	4% 86% 6% 7%
1	C	576	4% 86% 7% 7%
1	D	576	4% 87% 5% 7%
2	E	8	25% 75%

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Mol	Chain	Length	Quality of chain
2	G	8	 38% 62%
2	H	8	 50% 50%
2	I	8	 38% 62%
3	F	2	 100%

2 Entry composition [i](#)

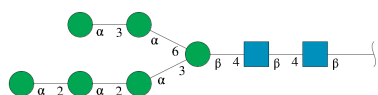
There are 10 unique types of molecules in this entry. The entry contains 18328 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Invertase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	535	Total	C	N	O	S	0	1	0
			4155	2644	707	790	14			
1	B	536	Total	C	N	O	S	0	0	0
			4151	2642	704	791	14			
1	C	536	Total	C	N	O	S	0	0	0
			4151	2642	704	791	14			
1	D	535	Total	C	N	O	S	0	0	0
			4144	2638	703	789	14			

- Molecule 2 is an oligosaccharide called alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-3)-alpha-D-mannopyranose-(1-6)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



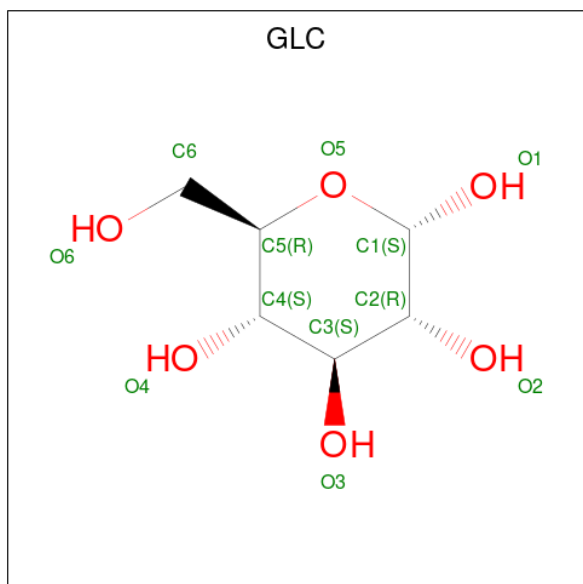
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	E	8	Total	C	N	O	0	0	0
			94	52	2	40			
2	G	8	Total	C	N	O	0	0	0
			94	52	2	40			
2	H	8	Total	C	N	O	0	0	0
			94	52	2	40			
2	I	8	Total	C	N	O	0	0	0
			94	52	2	40			

- Molecule 3 is an oligosaccharide called 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



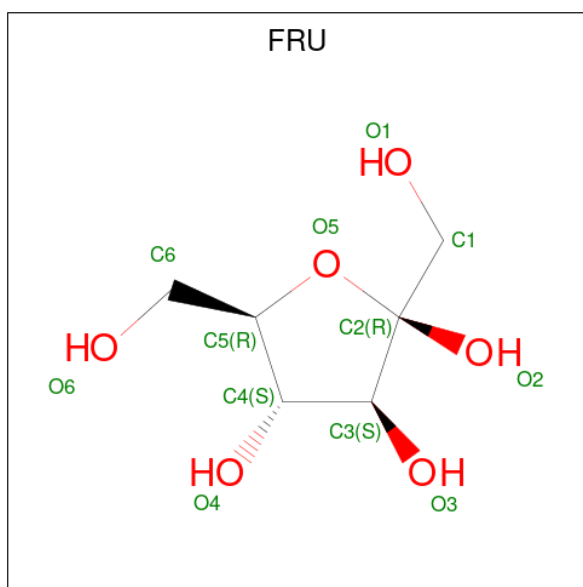
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace
3	F	2	Total	C	N	O	0	0
			28	16	2	10		

- Molecule 4 is alpha-D-glucopyranose (CCD ID: GLC) (formula: C₆H₁₂O₆) (labeled as "Ligand of Interest" by depositor).



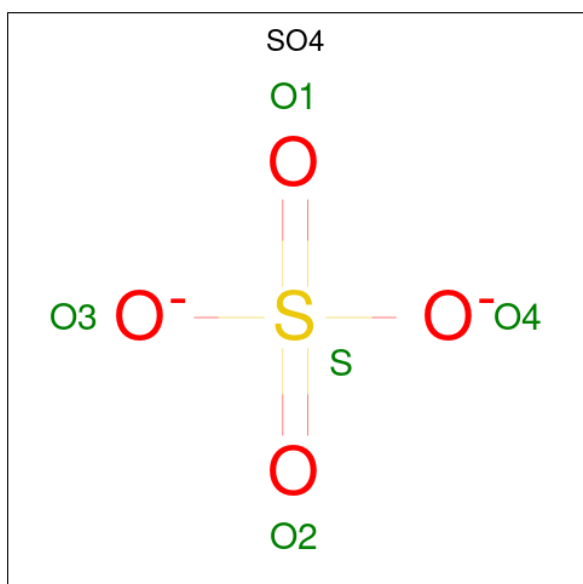
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			11	6	5		
4	B	1	Total	C	O	0	0
			11	6	5		
4	C	1	Total	C	O	0	0
			11	6	5		
4	D	1	Total	C	O	0	0
			11	6	5		

- Molecule 5 is beta-D-fructofuranose (CCD ID: FRU) (formula: C₆H₁₂O₆) (labeled as "Ligand of Interest" by depositor).



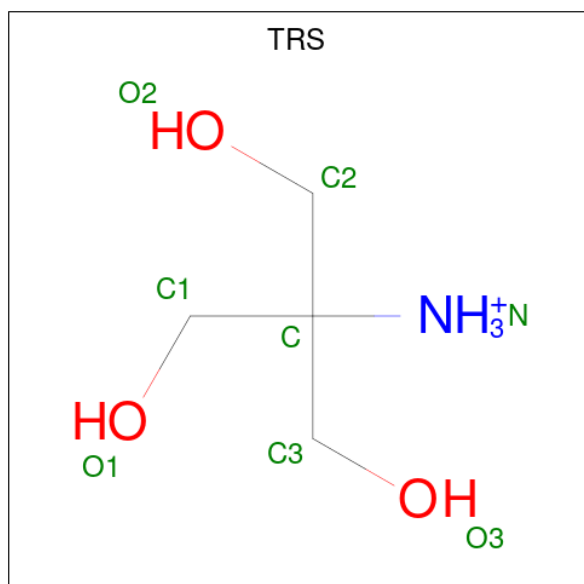
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			12	6	6		
5	B	1	Total	C	O	0	0
			12	6	6		
5	C	1	Total	C	O	0	0
			12	6	6		
5	D	1	Total	C	O	0	0
			12	6	6		

- Molecule 6 is SULFATE ION (CCD ID: SO4) (formula: O₄S) (labeled as "Ligand of Interest" by depositor).



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

- Molecule 7 is 2-AMINO-2-HYDROXYMETHYL-PROPANE-1,3-DIOL (CCD ID: TRS) (formula: $C_4H_{12}NO_3$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
7	A	1	Total	C	N	O	0	0
			8	4	1	3		
7	B	1	Total	C	N	O	0	0
			8	4	1	3		
7	C	1	Total	C	N	O	0	0
			8	4	1	3		
7	D	1	Total	C	N	O	0	0
			8	4	1	3		

- Molecule 8 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: $C_8H_{15}NO_6$) (labeled as "Ligand of Interest" by depositor).



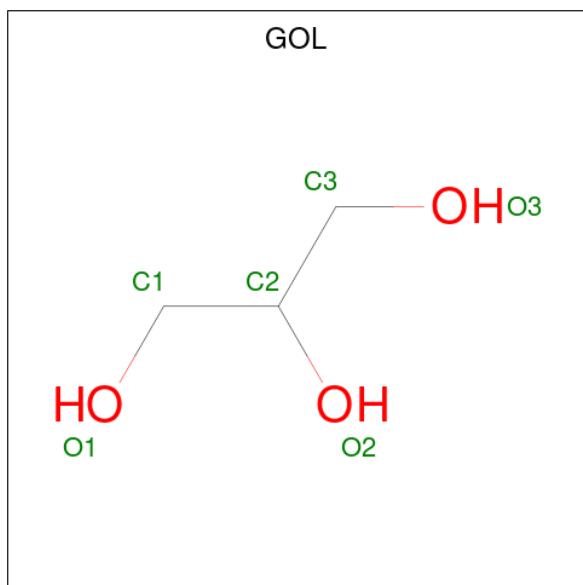
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
8	A	1	Total	C	N	O	0	0
			14	8	1	5		
8	A	1	Total	C	N	O	0	0
			14	8	1	5		
8	A	1	Total	C	N	O	0	0
			14	8	1	5		
8	A	1	Total	C	N	O	0	0
			14	8	1	5		
8	B	1	Total	C	N	O	0	0
			14	8	1	5		
8	B	1	Total	C	N	O	0	0
			14	8	1	5		
8	B	1	Total	C	N	O	0	0
			14	8	1	5		
8	C	1	Total	C	N	O	0	0
			14	8	1	5		
8	C	1	Total	C	N	O	0	0
			14	8	1	5		
8	C	1	Total	C	N	O	0	0
			14	8	1	5		
8	C	1	Total	C	N	O	0	0
			14	8	1	5		
8	D	1	Total	C	N	O	0	0
			14	8	1	5		
8	D	1	Total	C	N	O	0	0
			14	8	1	5		
8	D	1	Total	C	N	O	0	0
			14	8	1	5		

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
8	D	1	Total	C	N	O	0	0
			14	8	1	5		

- Molecule 9 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$) (labeled as "Ligand of Interest" by depositor).



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

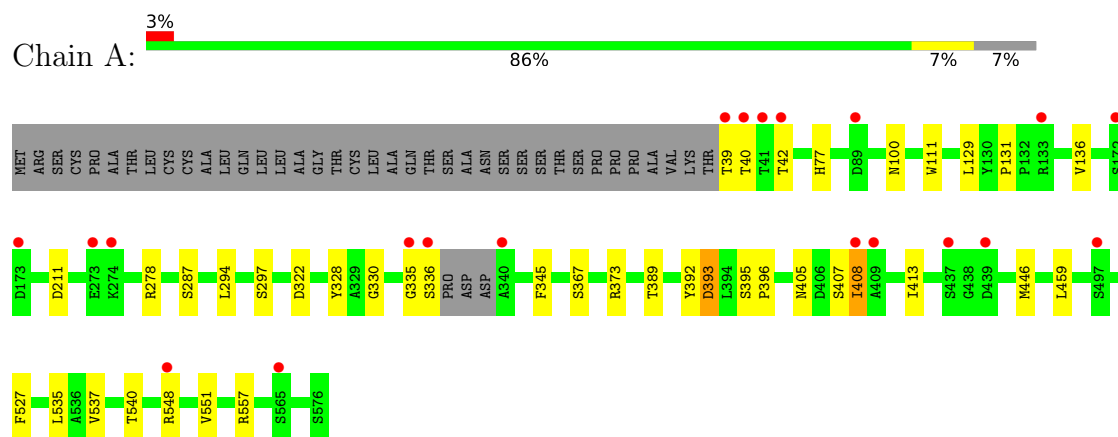
- Molecule 10 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
10	A	275	Total	O	0	0
			275	275		
10	B	232	Total	O	0	0
			232	232		
10	C	247	Total	O	0	0
			247	247		
10	D	199	Total	O	0	0
			199	199		

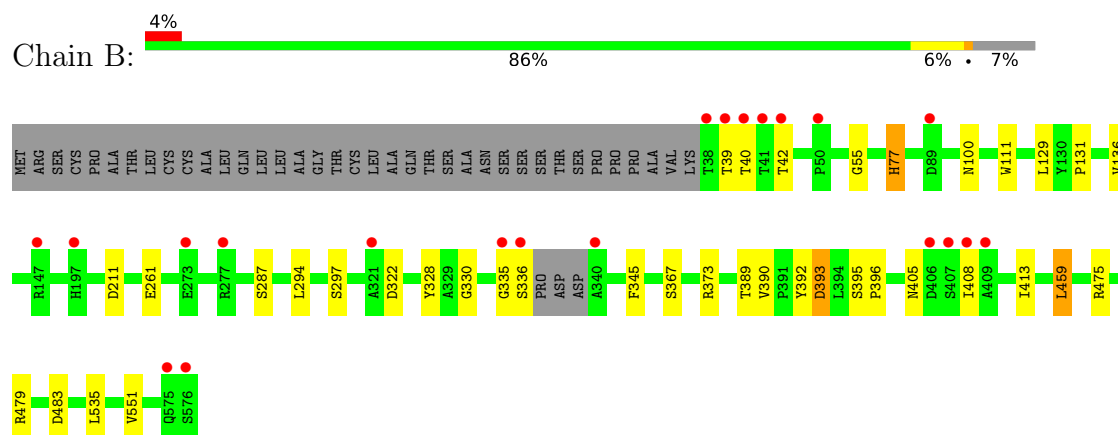
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

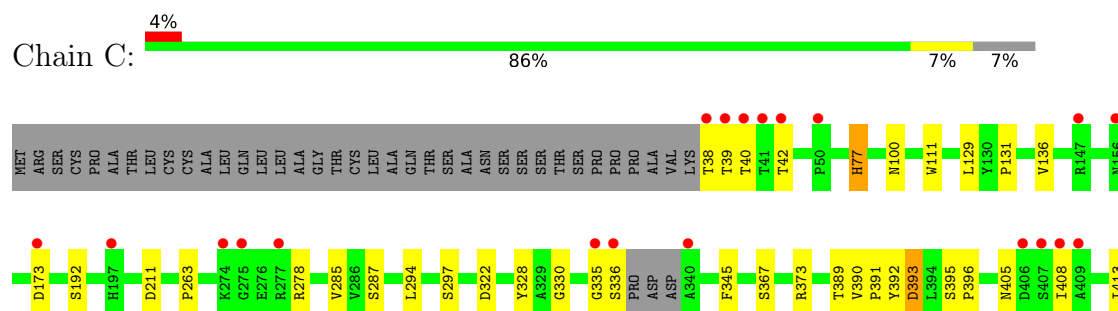
• Molecule 1: Invertase



• Molecule 1: Invertase



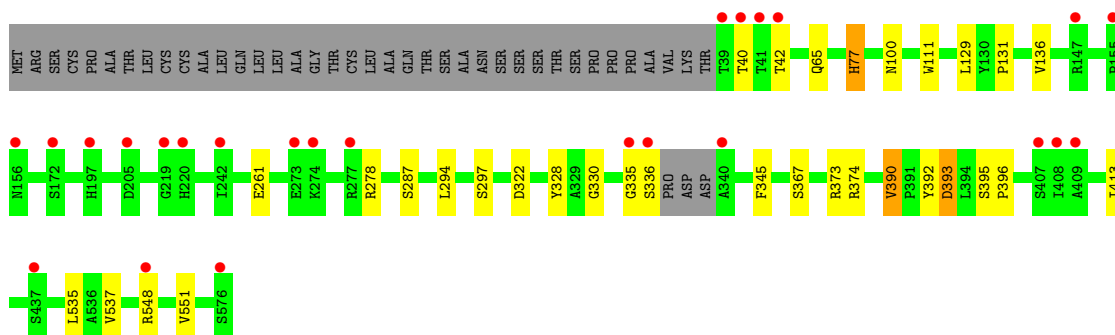
• Molecule 1: Invertase





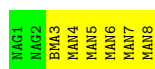
- Molecule 1: Invertase

Chain D: 4% 87% 5% 7%



- Molecule 2: alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-3)-alpha-D-mannopyranose-(1-6)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain E: 25% 75%



- Molecule 2: alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-3)-alpha-D-mannopyranose-(1-6)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain G: 38% 62%



- Molecule 2: alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-3)-alpha-D-mannopyranose-(1-6)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain H: 50% 50%



- Molecule 2: alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-3)-alpha-D-mannopyranose-(1-6)]beta-D-mannopyranose-(1

-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain I:  38% 62%



● Molecule 3: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain F:  100%



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	79.10Å 141.12Å 252.66Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.41 – 2.00 49.41 – 2.00	Depositor EDS
% Data completeness (in resolution range)	99.9 (49.41-2.00) 99.9 (49.41-2.00)	Depositor EDS
R_{merge}	0.13	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.45 (at 2.00Å)	Xtriage
Refinement program	REFMAC 5.8.0425	Depositor
R, R_{free}	0.179 , 0.198 0.187 , 0.207	Depositor DCC
R_{free} test set	9588 reflections (5.05%)	wwPDB-VP
Wilson B-factor (Å ²)	30.9	Xtriage
Anisotropy	0.039	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 32.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	18328	wwPDB-VP
Average B, all atoms (Å ²)	37.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.78% 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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: SO4, MAN, BMA, NAG, GLC, GOL, TRS, FRU

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	0.61	0/4295	0.94	1/5881 (0.0%)
1	B	0.59	0/4291	0.94	2/5877 (0.0%)
1	C	0.60	0/4291	0.94	3/5877 (0.1%)
1	D	0.59	0/4284	0.93	1/5867 (0.0%)
All	All	0.60	0/17161	0.94	7/23502 (0.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	2
1	B	0	1
1	C	0	1
1	D	0	2
All	All	0	6

There are no bond length outliers.

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	322	ASP	CA-CB-CG	6.94	119.54	112.60
1	C	322	ASP	CA-CB-CG	6.31	118.91	112.60
1	A	322	ASP	CA-CB-CG	6.17	118.77	112.60
1	B	322	ASP	CA-CB-CG	6.12	118.72	112.60
1	B	459	LEU	N-CA-CB	-5.40	101.60	110.68
1	C	38	THR	CA-CB-OG1	-5.04	102.03	109.60
1	C	479	ARG	NE-CZ-NH1	-5.04	116.46	121.50

There are no chirality outliers.

All (6) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	278	ARG	Sidechain
1	A	557	ARG	Sidechain
1	B	479	ARG	Sidechain
1	C	278	ARG	Sidechain
1	D	278	ARG	Sidechain
1	D	548	ARG	Sidechain

5.2 Too-close contacts

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	4155	0	3854	28	0
1	B	4151	0	3849	27	0
1	C	4151	0	3849	26	0
1	D	4144	0	3842	21	0
2	E	94	0	79	0	0
2	G	94	0	79	0	0
2	H	94	0	79	0	0
2	I	94	0	79	0	0
3	F	28	0	25	0	0
4	A	11	0	10	2	0
4	B	11	0	10	4	0
4	C	11	0	10	2	0
4	D	11	0	10	2	0
5	A	12	0	12	2	0
5	B	12	0	12	3	0
5	C	12	0	11	2	0
5	D	12	0	11	3	0
6	A	10	0	0	0	0
6	B	5	0	0	0	0
6	C	5	0	0	0	0
6	D	10	0	0	0	0
7	A	8	0	12	0	0
7	B	8	0	12	0	0
7	C	8	0	12	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
7	D	8	0	12	0	0
8	A	56	0	52	0	0
8	B	42	0	39	0	0
8	C	56	0	52	0	0
8	D	56	0	52	0	0
9	B	6	0	8	1	0
10	A	275	0	0	3	0
10	B	232	0	0	2	0
10	C	247	0	0	4	0
10	D	199	0	0	2	0
All	All	18328	0	16072	105	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 3.

All (105) 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)
4:C:607:GLC:C1	5:C:608:FRU:O2	1.65	1.43
4:D:601:GLC:C1	5:D:602:FRU:O2	1.65	1.41
4:A:601:GLC:C1	5:A:602:FRU:O2	1.69	1.40
4:B:601:GLC:C1	5:B:602:FRU:O2	1.71	1.38
1:B:535:LEU:HD21	1:C:535:LEU:HD21	1.40	1.04
1:A:535:LEU:HD21	1:D:535:LEU:HD21	1.52	0.90
1:C:39:THR:HG21	10:C:765:HOH:O	1.74	0.87
4:D:601:GLC:C1	5:D:602:FRU:C2	2.57	0.83
4:C:607:GLC:C1	5:C:608:FRU:C2	2.60	0.79
1:A:39:THR:HG21	10:A:846:HOH:O	1.84	0.77
4:B:601:GLC:C1	5:B:602:FRU:C2	2.63	0.76
1:A:294:LEU:HD23	1:B:294:LEU:HD23	1.69	0.74
4:A:601:GLC:C1	5:A:602:FRU:C2	2.67	0.72
1:A:40:THR:HG22	1:A:42:THR:H	1.57	0.69
1:B:40:THR:HG22	1:B:42:THR:H	1.59	0.67
1:D:40:THR:HG22	1:D:42:THR:H	1.61	0.66
1:D:131:PRO:HB3	1:D:136:VAL:O	1.97	0.65
1:C:131:PRO:HB3	1:C:136:VAL:O	1.97	0.65
1:A:131:PRO:HB3	1:A:136:VAL:O	1.97	0.65
1:C:294:LEU:HD23	1:D:294:LEU:HD23	1.79	0.63
1:C:40:THR:HG22	1:C:42:THR:H	1.63	0.63
1:B:287:SER:HB3	1:B:328:TYR:CE1	2.37	0.60
1:D:77:HIS:HE1	10:D:813:HOH:O	1.83	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:287:SER:HB3	1:C:328:TYR:CE1	2.36	0.59
1:A:287:SER:HB3	1:A:328:TYR:CE1	2.38	0.59
1:D:287:SER:HB3	1:D:328:TYR:CE1	2.39	0.58
1:B:131:PRO:HB3	1:B:136:VAL:O	2.05	0.56
1:C:77:HIS:HE1	10:C:807:HOH:O	1.92	0.53
1:D:65:GLN:NE2	10:D:704:HOH:O	2.42	0.53
1:C:297:SER:HB2	1:C:328:TYR:CD1	2.44	0.52
1:B:483:ASP:HA	9:B:607:GOL:H11	1.92	0.51
1:B:297:SER:HB2	1:B:328:TYR:CD1	2.46	0.51
1:A:407:SER:HB3	1:A:548:ARG:HH11	1.75	0.51
1:A:297:SER:HB2	1:A:328:TYR:CD1	2.46	0.50
1:D:297:SER:HB2	1:D:328:TYR:CD1	2.46	0.50
1:A:39:THR:HG22	1:A:40:THR:H	1.76	0.50
1:A:408:ILE:HD11	1:A:551:VAL:HG12	1.93	0.49
1:C:335:GLY:O	1:C:336:SER:C	2.55	0.49
1:B:335:GLY:O	1:B:336:SER:C	2.56	0.49
1:D:395:SER:N	1:D:396:PRO:CD	2.76	0.48
1:D:374:ARG:HB2	1:D:390:VAL:HG13	1.96	0.48
1:D:335:GLY:O	1:D:336:SER:C	2.57	0.48
1:B:395:SER:N	1:B:396:PRO:CD	2.77	0.47
1:C:395:SER:N	1:C:396:PRO:CD	2.77	0.47
1:A:413:ILE:HG13	1:A:551:VAL:HG11	1.96	0.47
1:D:413:ILE:HG13	1:D:551:VAL:HG11	1.96	0.47
1:A:395:SER:N	1:A:396:PRO:CD	2.78	0.47
1:A:335:GLY:O	1:A:336:SER:C	2.58	0.46
1:C:77:HIS:HD2	10:C:924:HOH:O	1.98	0.46
1:B:261:GLU:OE1	4:B:601:GLC:H3	2.15	0.46
1:C:413:ILE:HG13	1:C:551:VAL:HG11	1.98	0.45
1:B:77:HIS:HB3	1:B:100:ASN:O	2.16	0.45
1:B:408:ILE:HG22	1:B:408:ILE:O	2.16	0.45
1:B:459:LEU:HD12	1:B:475:ARG:HA	1.98	0.45
1:B:413:ILE:HG13	1:B:551:VAL:HG11	1.97	0.45
1:B:373:ARG:HD3	1:B:389:THR:CG2	2.47	0.45
1:A:39:THR:HG22	1:A:40:THR:N	2.31	0.45
1:C:408:ILE:O	1:C:408:ILE:HG22	2.16	0.45
1:B:392:TYR:O	1:B:393:ASP:C	2.59	0.44
1:C:392:TYR:O	1:C:393:ASP:C	2.60	0.44
1:B:77:HIS:HD2	10:B:912:HOH:O	2.00	0.44
1:A:77:HIS:HD2	10:A:957:HOH:O	2.01	0.44
1:A:537:VAL:HG21	1:D:535:LEU:HD22	2.00	0.44
1:A:405:ASN:HB3	1:A:408:ILE:HD11	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:392:TYR:O	1:A:393:ASP:C	2.61	0.43
1:A:77:HIS:HB3	1:A:100:ASN:O	2.19	0.43
1:B:77:HIS:HE1	10:B:722:HOH:O	1.99	0.43
1:D:392:TYR:O	1:D:393:ASP:C	2.60	0.43
1:A:77:HIS:HE1	10:A:942:HOH:O	2.01	0.43
1:C:330:GLY:HA3	1:C:345:PHE:CZ	2.54	0.43
1:D:373:ARG:HH21	1:D:373:ARG:HG3	1.84	0.43
1:A:373:ARG:CG	1:A:373:ARG:HH21	2.32	0.43
1:B:373:ARG:CD	1:B:389:THR:CG2	2.97	0.43
1:C:373:ARG:HH21	1:C:373:ARG:HG3	1.83	0.43
1:C:373:ARG:HH21	1:C:373:ARG:CG	2.32	0.42
1:B:373:ARG:HH21	1:B:373:ARG:HG3	1.83	0.42
1:C:111:TRP:CD2	1:C:129:LEU:HD12	2.54	0.42
1:C:405:ASN:HB3	1:C:408:ILE:HD11	2.02	0.42
1:D:77:HIS:HB3	1:D:100:ASN:O	2.19	0.42
1:A:535:LEU:HD22	1:D:537:VAL:HG21	2.01	0.42
1:B:373:ARG:HB3	1:B:389:THR:HG23	2.02	0.42
1:C:390:VAL:HG22	1:C:391:PRO:HD2	2.01	0.42
1:B:373:ARG:HH21	1:B:373:ARG:CG	2.33	0.42
1:A:373:ARG:HH21	1:A:373:ARG:HG3	1.83	0.42
1:A:111:TRP:CD2	1:A:129:LEU:HD12	2.55	0.42
1:C:459:LEU:HD12	1:C:475:ARG:HA	2.02	0.41
1:A:330:GLY:HA3	1:A:345:PHE:CZ	2.55	0.41
1:B:330:GLY:HA3	1:B:345:PHE:CZ	2.55	0.41
1:C:39:THR:HG23	1:C:192:SER:HB3	2.02	0.41
1:C:419:ASP:HB2	10:C:751:HOH:O	2.19	0.41
1:D:111:TRP:CD2	1:D:129:LEU:HD12	2.56	0.41
1:D:330:GLY:HA3	1:D:345:PHE:CZ	2.55	0.41
1:D:373:ARG:HH21	1:D:373:ARG:CG	2.33	0.41
1:A:373:ARG:HD3	1:A:389:THR:CG2	2.50	0.41
1:C:263:PRO:HA	1:C:285:VAL:O	2.20	0.41
1:B:111:TRP:CD2	1:B:129:LEU:HD12	2.56	0.41
1:C:373:ARG:HD3	1:C:389:THR:CG2	2.51	0.41
1:A:446:MET:HG3	1:A:540:THR:HG22	2.02	0.41
1:B:39:THR:HB	1:B:55:GLY:HA2	2.03	0.41
1:B:408:ILE:HD12	1:B:413:ILE:CG1	2.51	0.41
4:B:601:GLC:C2	5:B:602:FRU:O2	2.58	0.41
1:D:261:GLU:OE2	5:D:602:FRU:H12	2.21	0.41
1:C:77:HIS:HB3	1:C:100:ASN:O	2.20	0.40
1:A:459:LEU:HD11	1:A:527:PHE:CD2	2.56	0.40
1:B:405:ASN:HB3	1:B:408:ILE:HD11	2.03	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	532/576 (92%)	509 (96%)	20 (4%)	3 (1%)	21	17
1	B	532/576 (92%)	510 (96%)	18 (3%)	4 (1%)	16	11
1	C	532/576 (92%)	509 (96%)	19 (4%)	4 (1%)	16	11
1	D	531/576 (92%)	508 (96%)	20 (4%)	3 (1%)	21	17
All	All	2127/2304 (92%)	2036 (96%)	77 (4%)	14 (1%)	18	14

All (14) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	367	SER
1	A	393	ASP
1	B	367	SER
1	B	393	ASP
1	C	367	SER
1	C	393	ASP
1	D	367	SER
1	B	77	HIS
1	D	393	ASP
1	C	77	HIS
1	D	77	HIS
1	A	211	ASP
1	B	211	ASP
1	C	211	ASP

5.3.2 Protein sidechains [i](#)

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

resolution.

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	441/475 (93%)	440 (100%)	1 (0%)	87	92
1	B	441/475 (93%)	440 (100%)	1 (0%)	87	92
1	C	441/475 (93%)	440 (100%)	1 (0%)	87	92
1	D	440/475 (93%)	439 (100%)	1 (0%)	87	92
All	All	1763/1900 (93%)	1759 (100%)	4 (0%)	87	92

All (4) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	408	ILE
1	B	390	VAL
1	C	173	ASP
1	D	390	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (27) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	51	ASN
1	A	65	GLN
1	A	77	HIS
1	A	98	GLN
1	A	197	HIS
1	B	51	ASN
1	B	65	GLN
1	B	69	GLN
1	B	77	HIS
1	B	80	ASN
1	B	98	GLN
1	B	197	HIS
1	B	289	ASN
1	B	423	ASN
1	B	531	GLN
1	B	554	ASN
1	C	65	GLN
1	C	77	HIS
1	C	98	GLN

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Mol	Chain	Res	Type
1	C	423	ASN
1	C	502	ASN
1	C	554	ASN
1	D	51	ASN
1	D	65	GLN
1	D	77	HIS
1	D	98	GLN
1	D	423	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

34 monosaccharides are modelled in this entry.

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	$\# Z > 2$
2	NAG	E	1	1,2	14,14,15	0.40	0	17,19,21	0.71	0
2	NAG	E	2	2	14,14,15	0.35	0	17,19,21	0.59	0
2	BMA	E	3	2	11,11,12	1.11	1 (9%)	15,15,17	0.58	0
2	MAN	E	4	2	11,11,12	0.92	0	15,15,17	0.92	1 (6%)
2	MAN	E	5	2	11,11,12	1.20	2 (18%)	15,15,17	0.63	0
2	MAN	E	6	2	11,11,12	1.16	2 (18%)	15,15,17	0.62	0
2	MAN	E	7	2	11,11,12	0.83	1 (9%)	15,15,17	0.82	1 (6%)
2	MAN	E	8	2	11,11,12	0.54	0	15,15,17	0.87	1 (6%)
3	NAG	F	1	1,3	14,14,15	0.42	0	17,19,21	0.70	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	NAG	F	2	3	14,14,15	0.38	0	17,19,21	0.52	0
2	NAG	G	1	1,2	14,14,15	0.42	0	17,19,21	0.72	0
2	NAG	G	2	2	14,14,15	0.42	0	17,19,21	0.67	1 (5%)
2	BMA	G	3	2	11,11,12	0.91	1 (9%)	15,15,17	0.66	0
2	MAN	G	4	2	11,11,12	0.71	1 (9%)	15,15,17	0.60	0
2	MAN	G	5	2	11,11,12	0.59	0	15,15,17	0.61	0
2	MAN	G	6	2	11,11,12	0.55	0	15,15,17	0.56	0
2	MAN	G	7	2	11,11,12	1.31	1 (9%)	15,15,17	0.45	0
2	MAN	G	8	2	11,11,12	0.50	0	15,15,17	0.88	1 (6%)
2	NAG	H	1	1,2	14,14,15	0.38	0	17,19,21	0.56	0
2	NAG	H	2	2	14,14,15	0.51	0	17,19,21	0.72	0
2	BMA	H	3	2	11,11,12	0.82	1 (9%)	15,15,17	0.64	0
2	MAN	H	4	2	11,11,12	0.77	0	15,15,17	1.02	1 (6%)
2	MAN	H	5	2	11,11,12	0.89	0	15,15,17	0.71	0
2	MAN	H	6	2	11,11,12	0.65	0	15,15,17	0.69	0
2	MAN	H	7	2	11,11,12	1.11	1 (9%)	15,15,17	0.54	0
2	MAN	H	8	2	11,11,12	0.60	0	15,15,17	0.83	1 (6%)
2	NAG	I	1	1,2	14,14,15	0.37	0	17,19,21	0.71	0
2	NAG	I	2	2	14,14,15	0.40	0	17,19,21	0.76	0
2	BMA	I	3	2	11,11,12	0.97	1 (9%)	15,15,17	0.68	0
2	MAN	I	4	2	11,11,12	0.65	0	15,15,17	0.99	2 (13%)
2	MAN	I	5	2	11,11,12	0.53	0	15,15,17	0.68	0
2	MAN	I	6	2	11,11,12	0.92	1 (9%)	15,15,17	0.64	0
2	MAN	I	7	2	11,11,12	0.88	1 (9%)	15,15,17	0.72	0
2	MAN	I	8	2	11,11,12	0.88	1 (9%)	15,15,17	0.67	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
2	NAG	E	1	1,2	-	0/6/23/26	0/1/1/1
2	NAG	E	2	2	-	0/6/23/26	0/1/1/1
2	BMA	E	3	2	-	0/2/19/22	0/1/1/1
2	MAN	E	4	2	-	0/2/19/22	0/1/1/1
2	MAN	E	5	2	-	0/2/19/22	0/1/1/1
2	MAN	E	6	2	-	0/2/19/22	0/1/1/1
2	MAN	E	7	2	-	0/2/19/22	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	MAN	E	8	2	-	0/2/19/22	0/1/1/1
3	NAG	F	1	1,3	-	0/6/23/26	0/1/1/1
3	NAG	F	2	3	-	0/6/23/26	0/1/1/1
2	NAG	G	1	1,2	-	0/6/23/26	0/1/1/1
2	NAG	G	2	2	-	0/6/23/26	0/1/1/1
2	BMA	G	3	2	-	0/2/19/22	0/1/1/1
2	MAN	G	4	2	-	0/2/19/22	0/1/1/1
2	MAN	G	5	2	-	0/2/19/22	0/1/1/1
2	MAN	G	6	2	-	0/2/19/22	0/1/1/1
2	MAN	G	7	2	-	0/2/19/22	0/1/1/1
2	MAN	G	8	2	-	0/2/19/22	0/1/1/1
2	NAG	H	1	1,2	-	0/6/23/26	0/1/1/1
2	NAG	H	2	2	-	0/6/23/26	0/1/1/1
2	BMA	H	3	2	-	0/2/19/22	0/1/1/1
2	MAN	H	4	2	-	0/2/19/22	0/1/1/1
2	MAN	H	5	2	-	0/2/19/22	0/1/1/1
2	MAN	H	6	2	-	0/2/19/22	0/1/1/1
2	MAN	H	7	2	-	0/2/19/22	0/1/1/1
2	MAN	H	8	2	-	0/2/19/22	0/1/1/1
2	NAG	I	1	1,2	-	0/6/23/26	0/1/1/1
2	NAG	I	2	2	-	0/6/23/26	0/1/1/1
2	BMA	I	3	2	-	0/2/19/22	0/1/1/1
2	MAN	I	4	2	-	0/2/19/22	0/1/1/1
2	MAN	I	5	2	-	0/2/19/22	0/1/1/1
2	MAN	I	6	2	-	0/2/19/22	0/1/1/1
2	MAN	I	7	2	-	0/2/19/22	0/1/1/1
2	MAN	I	8	2	-	0/2/19/22	0/1/1/1

All (15) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	G	7	MAN	O5-C5	4.08	1.51	1.43
2	E	3	BMA	O5-C5	3.17	1.49	1.43
2	H	7	MAN	O5-C5	3.15	1.49	1.43
2	E	6	MAN	C2-C3	-2.92	1.48	1.52
2	E	5	MAN	O5-C5	2.80	1.48	1.43
2	I	3	BMA	O5-C5	2.43	1.48	1.43
2	E	7	MAN	O5-C5	2.34	1.48	1.43
2	I	6	MAN	O5-C5	2.31	1.47	1.43
2	H	3	BMA	O5-C5	2.25	1.47	1.43
2	I	7	MAN	O5-C5	2.22	1.47	1.43
2	G	3	BMA	O5-C5	2.16	1.47	1.43

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	E	6	MAN	O5-C5	2.09	1.47	1.43
2	I	8	MAN	O4-C4	2.06	1.48	1.43
2	E	5	MAN	C2-C3	-2.06	1.49	1.52
2	G	4	MAN	O5-C5	2.01	1.47	1.43

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	G	8	MAN	C1-O5-C5	2.72	115.83	112.19
2	E	8	MAN	C1-O5-C5	2.64	115.73	112.19
2	H	4	MAN	C1-C2-C3	2.34	113.05	109.64
2	H	8	MAN	C1-O5-C5	2.33	115.30	112.19
2	E	7	MAN	C1-O5-C5	2.32	115.30	112.19
2	I	4	MAN	C1-C2-C3	2.16	112.80	109.64
2	I	4	MAN	C1-O5-C5	2.10	115.00	112.19
2	G	2	NAG	C2-N2-C7	2.05	125.65	122.90
2	E	4	MAN	O3-C3-C2	-2.01	105.96	110.05

There are no chirality outliers.

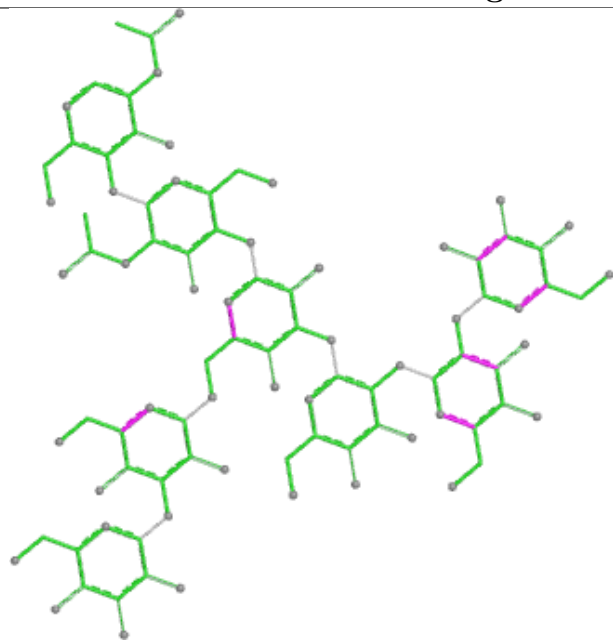
There are no torsion outliers.

There are no ring outliers.

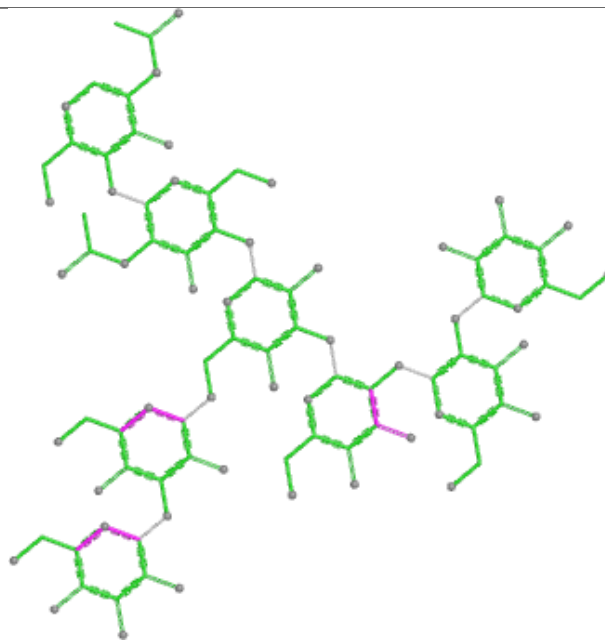
No monomer is involved in short contacts.

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.

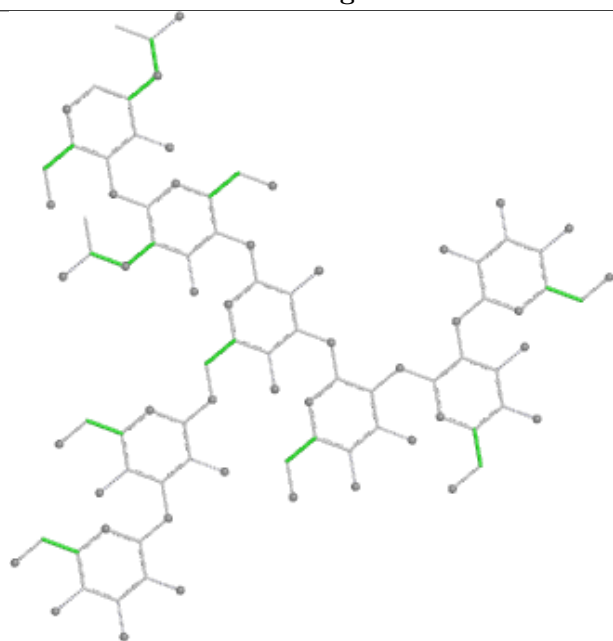
Oligosaccharide Chain E



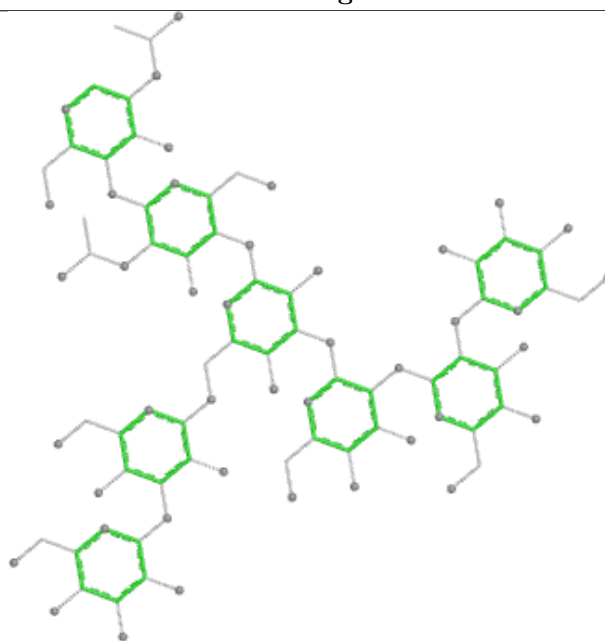
Bond lengths



Bond angles

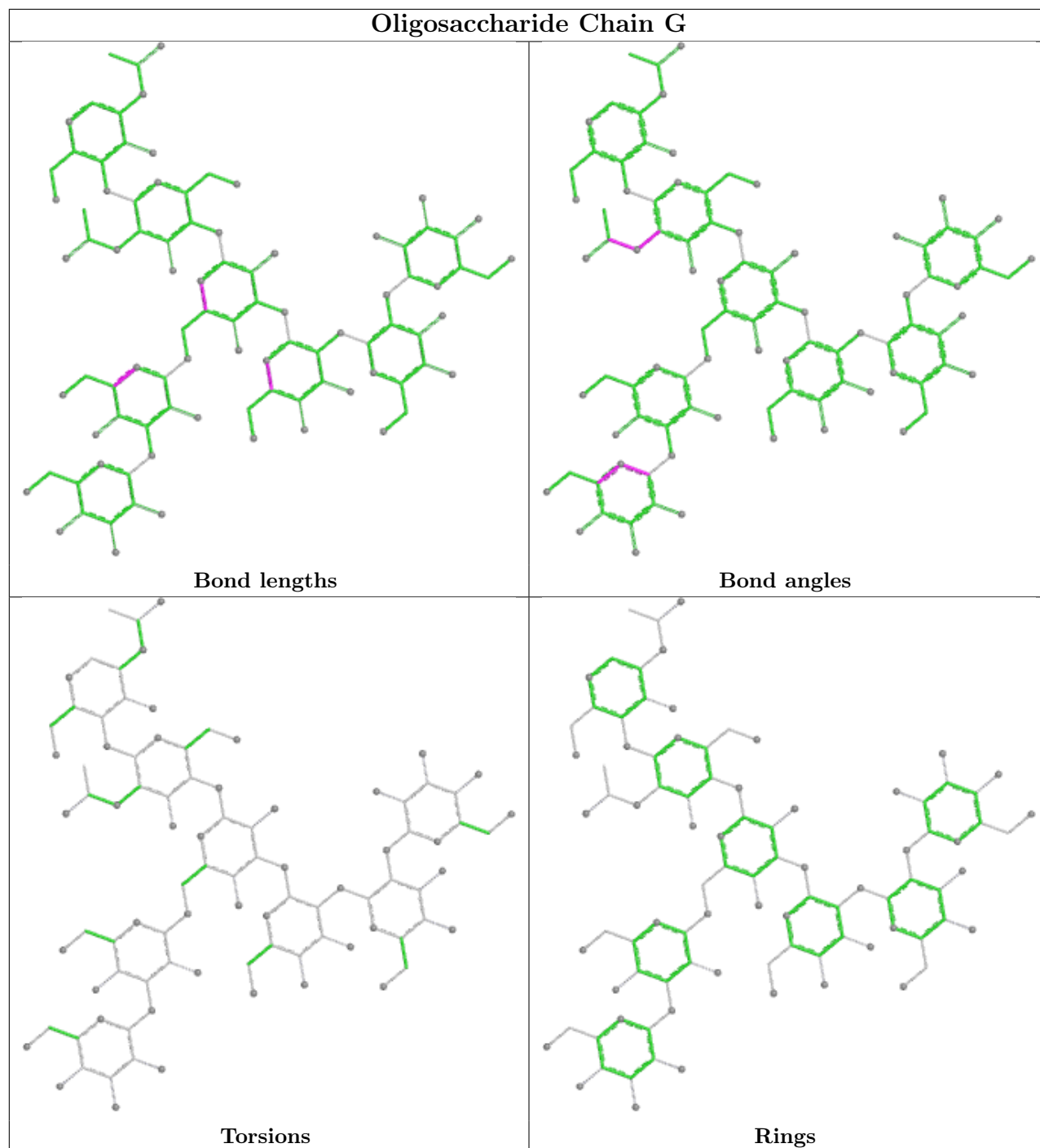


Torsions

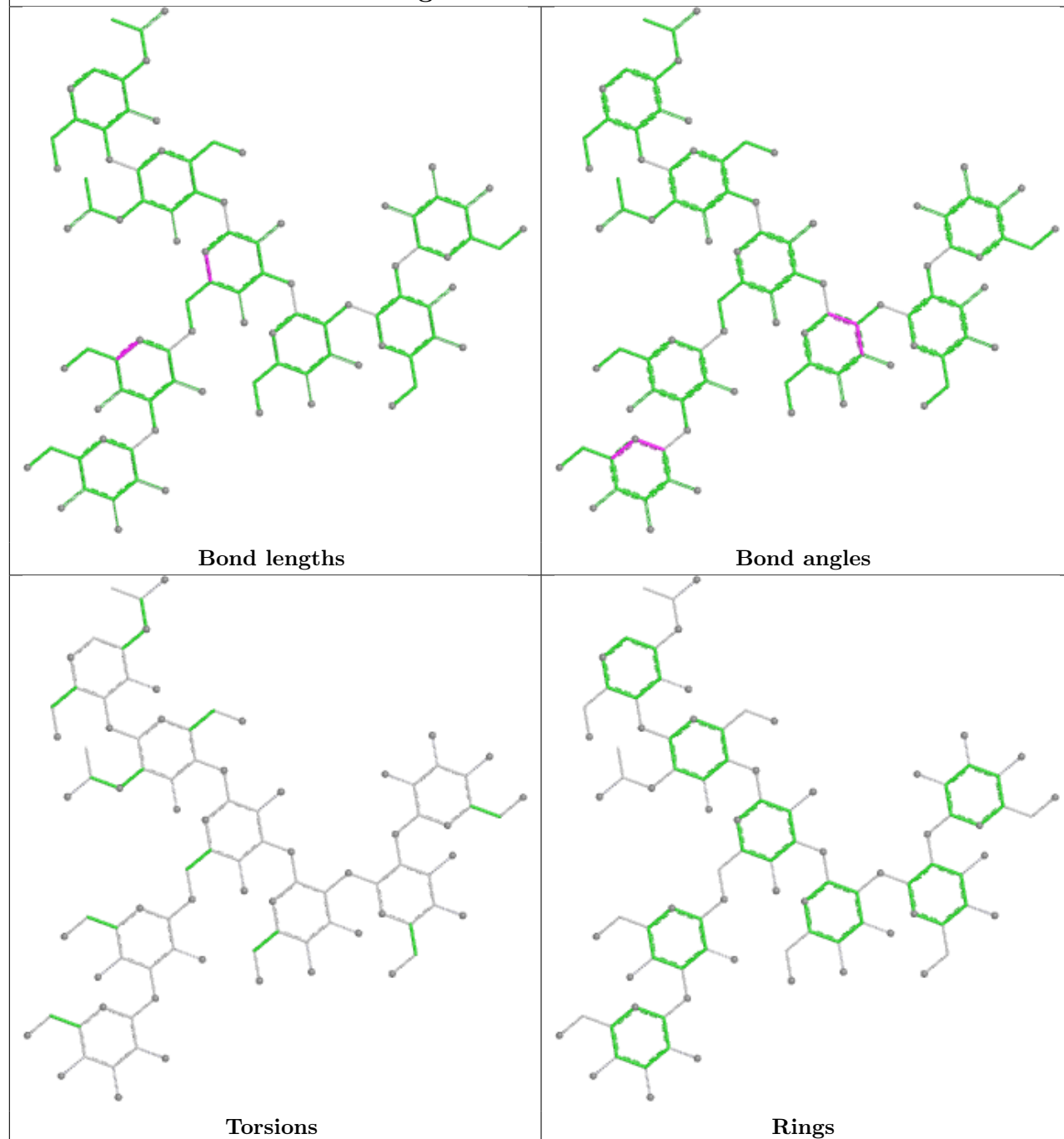


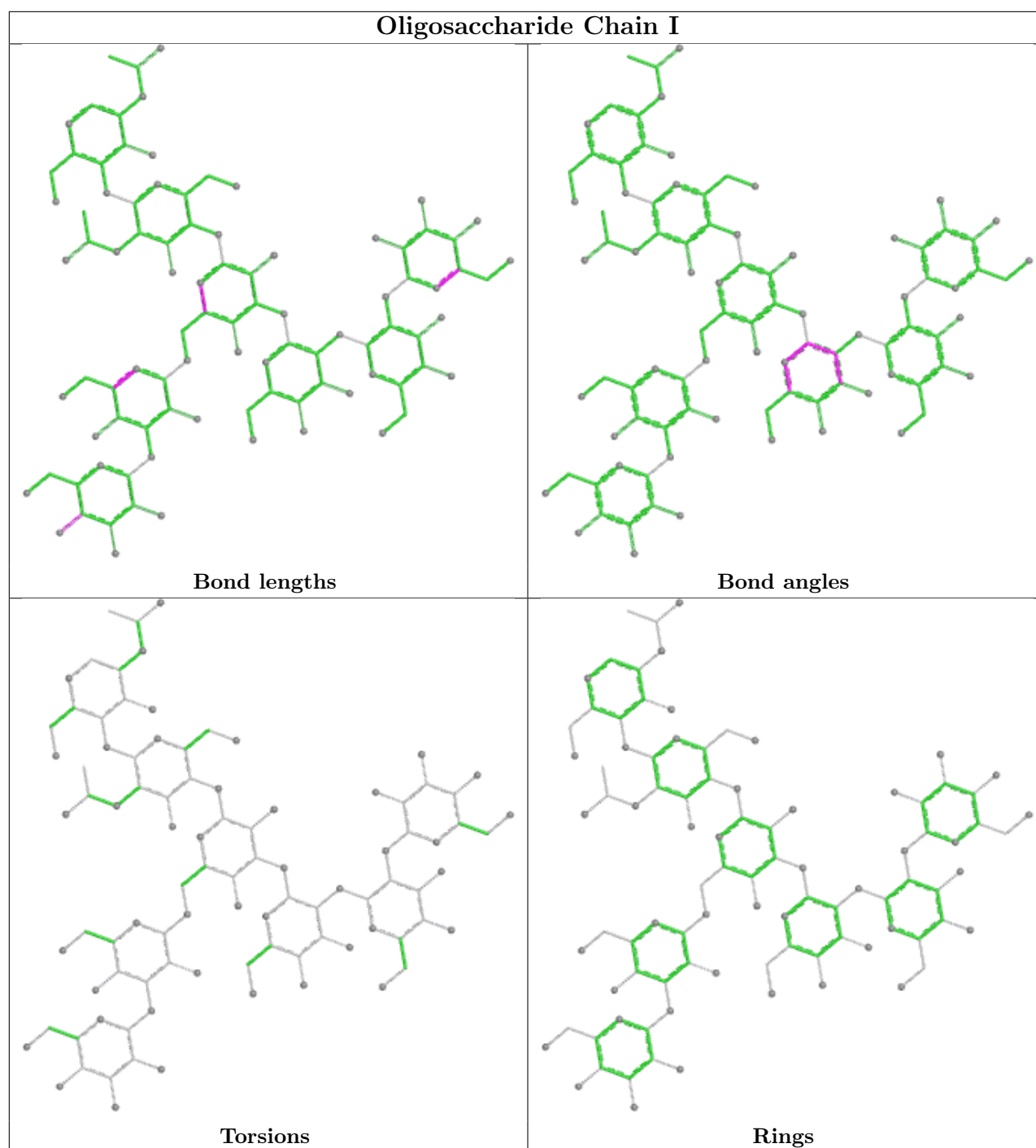
Rings

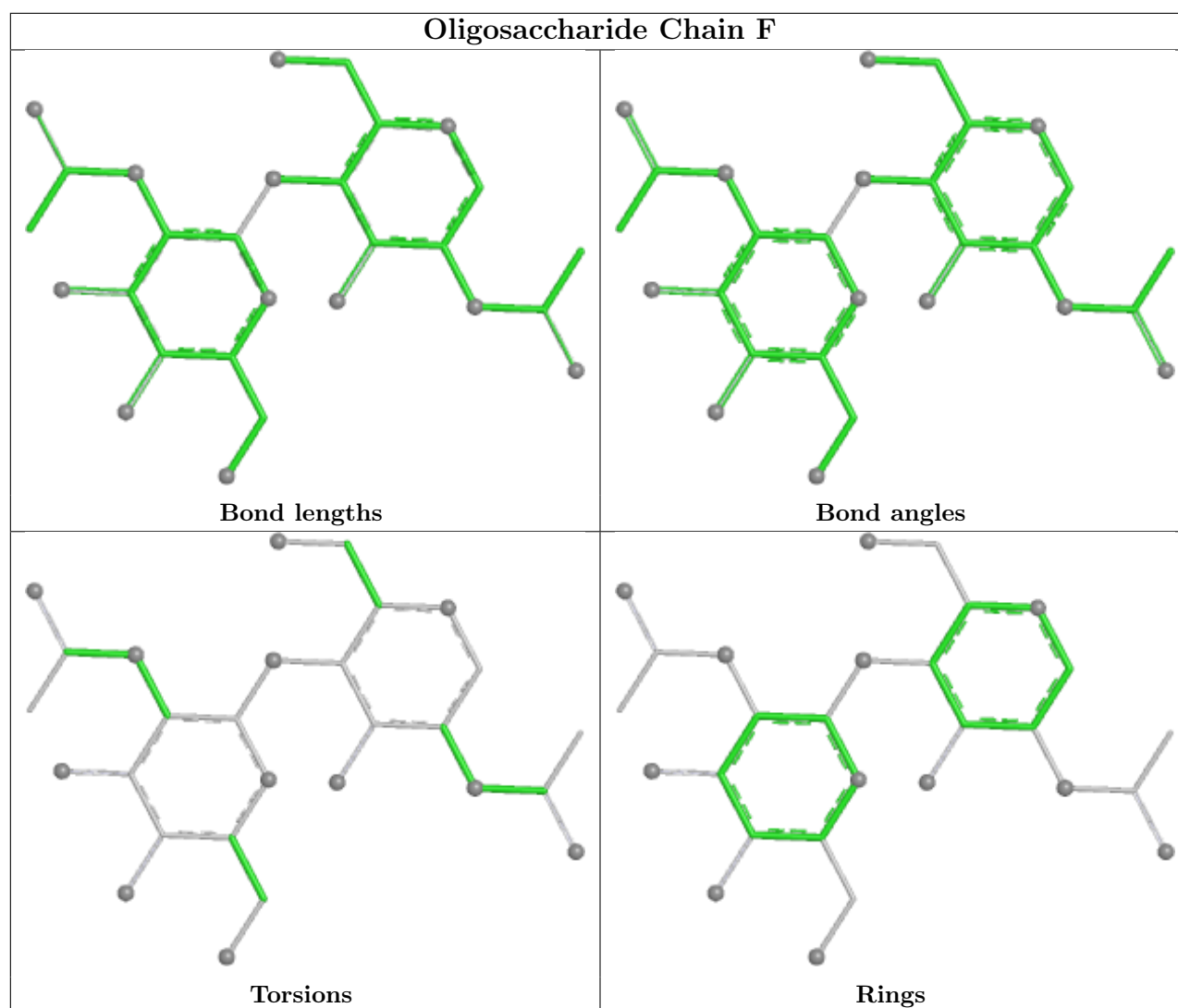
Oligosaccharide Chain G



Oligosaccharide Chain H







5.6 Ligand geometry [i](#)

34 ligands are modelled in this entry.

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	$\# Z > 2$
6	SO4	A	609	-	4,4,4	0.20	0	6,6,6	0.28	0
4	GLC	D	601	-	11,11,12	0.34	0	15,15,17	0.71	0
5	FRU	C	608	-	11,12,12	0.57	0	10,18,18	0.61	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
8	NAG	B	605	1	14,14,15	0.44	0	17,19,21	0.87	0
8	NAG	C	604	1	14,14,15	0.45	0	17,19,21	0.76	0
6	SO4	B	603	-	4,4,4	0.52	0	6,6,6	0.26	0
9	GOL	B	607	-	5,5,5	0.16	0	5,5,5	0.45	0
8	NAG	C	606	1	14,14,15	0.37	0	17,19,21	0.72	0
8	NAG	A	608	1	14,14,15	0.39	0	17,19,21	0.69	0
8	NAG	D	606	1	14,14,15	0.26	0	17,19,21	0.65	0
4	GLC	C	607	-	11,11,12	0.47	0	15,15,17	0.68	0
8	NAG	A	606	1	14,14,15	0.47	0	17,19,21	0.63	0
4	GLC	B	601	-	11,11,12	1.24	2 (18%)	15,15,17	1.38	2 (13%)
5	FRU	A	602	-	11,12,12	0.32	0	10,18,18	0.71	0
6	SO4	C	601	-	4,4,4	0.41	0	6,6,6	0.24	0
8	NAG	A	607	1	14,14,15	0.56	0	17,19,21	0.79	1 (5%)
8	NAG	D	609	1	14,14,15	0.38	0	17,19,21	0.55	0
7	TRS	B	608	-	7,7,7	0.16	0	9,9,9	0.44	0
4	GLC	A	601	-	11,11,12	0.52	0	15,15,17	0.72	0
6	SO4	D	604	-	4,4,4	0.39	0	6,6,6	0.25	0
7	TRS	A	604	-	7,7,7	0.25	0	9,9,9	0.44	0
5	FRU	B	602	-	11,12,12	0.50	0	10,18,18	1.07	1 (10%)
6	SO4	D	603	-	4,4,4	0.32	0	6,6,6	0.23	0
8	NAG	D	607	1	14,14,15	0.37	0	17,19,21	0.67	0
6	SO4	A	603	-	4,4,4	0.44	0	6,6,6	0.48	0
7	TRS	D	605	-	7,7,7	0.22	0	9,9,9	0.28	0
8	NAG	C	605	1	14,14,15	0.45	0	17,19,21	0.77	1 (5%)
5	FRU	D	602	-	11,12,12	0.70	0	10,18,18	0.66	0
8	NAG	B	604	1	14,14,15	0.45	0	17,19,21	0.59	0
8	NAG	A	605	1	14,14,15	0.38	0	17,19,21	0.77	1 (5%)
8	NAG	B	606	1	14,14,15	0.42	0	17,19,21	0.71	0
8	NAG	D	608	1	14,14,15	0.43	0	17,19,21	0.93	1 (5%)
8	NAG	C	603	1	14,14,15	0.40	0	17,19,21	0.47	0
7	TRS	C	602	-	7,7,7	0.39	0	9,9,9	0.47	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
8	NAG	B	605	1	-	0/6/23/26	0/1/1/1
4	GLC	D	601	-	-	2/2/19/22	0/1/1/1
5	FRU	C	608	-	-	1/5/24/24	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
8	NAG	C	604	1	-	0/6/23/26	0/1/1/1
9	GOL	B	607	-	-	0/4/4/4	-
8	NAG	C	606	1	-	0/6/23/26	0/1/1/1
8	NAG	A	608	1	-	0/6/23/26	0/1/1/1
8	NAG	D	606	1	-	0/6/23/26	0/1/1/1
4	GLC	C	607	-	-	0/2/19/22	0/1/1/1
8	NAG	A	606	1	-	0/6/23/26	0/1/1/1
4	GLC	B	601	-	-	2/2/19/22	0/1/1/1
5	FRU	A	602	-	-	1/5/24/24	0/1/1/1
8	NAG	A	607	1	-	0/6/23/26	0/1/1/1
8	NAG	D	609	1	-	0/6/23/26	0/1/1/1
7	TRS	B	608	-	-	6/9/9/9	-
4	GLC	A	601	-	-	0/2/19/22	0/1/1/1
7	TRS	A	604	-	-	3/9/9/9	-
5	FRU	B	602	-	-	1/5/24/24	0/1/1/1
8	NAG	D	607	1	-	0/6/23/26	0/1/1/1
7	TRS	D	605	-	-	7/9/9/9	-
8	NAG	C	605	1	-	0/6/23/26	0/1/1/1
5	FRU	D	602	-	-	3/5/24/24	0/1/1/1
8	NAG	B	604	1	-	0/6/23/26	0/1/1/1
8	NAG	A	605	1	-	0/6/23/26	0/1/1/1
8	NAG	B	606	1	-	0/6/23/26	0/1/1/1
8	NAG	D	608	1	-	0/6/23/26	0/1/1/1
8	NAG	C	603	1	-	0/6/23/26	0/1/1/1
7	TRS	C	602	-	-	5/9/9/9	-

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	601	GLC	C2-C3	-2.62	1.48	1.52
4	B	601	GLC	O5-C5	2.13	1.47	1.43

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	601	GLC	C1-C2-C3	3.70	115.03	109.64
4	B	601	GLC	O2-C2-C3	-2.95	104.03	110.15
5	B	602	FRU	O1-C1-C2	-2.54	106.06	111.67
8	C	605	NAG	C1-O5-C5	2.28	115.24	112.19
8	D	608	NAG	C2-N2-C7	2.26	125.92	122.90
8	A	607	NAG	C1-O5-C5	2.18	115.10	112.19

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
8	A	605	NAG	C1-C2-N2	2.15	113.82	110.43

There are no chirality outliers.

All (31) torsion outliers are listed below:

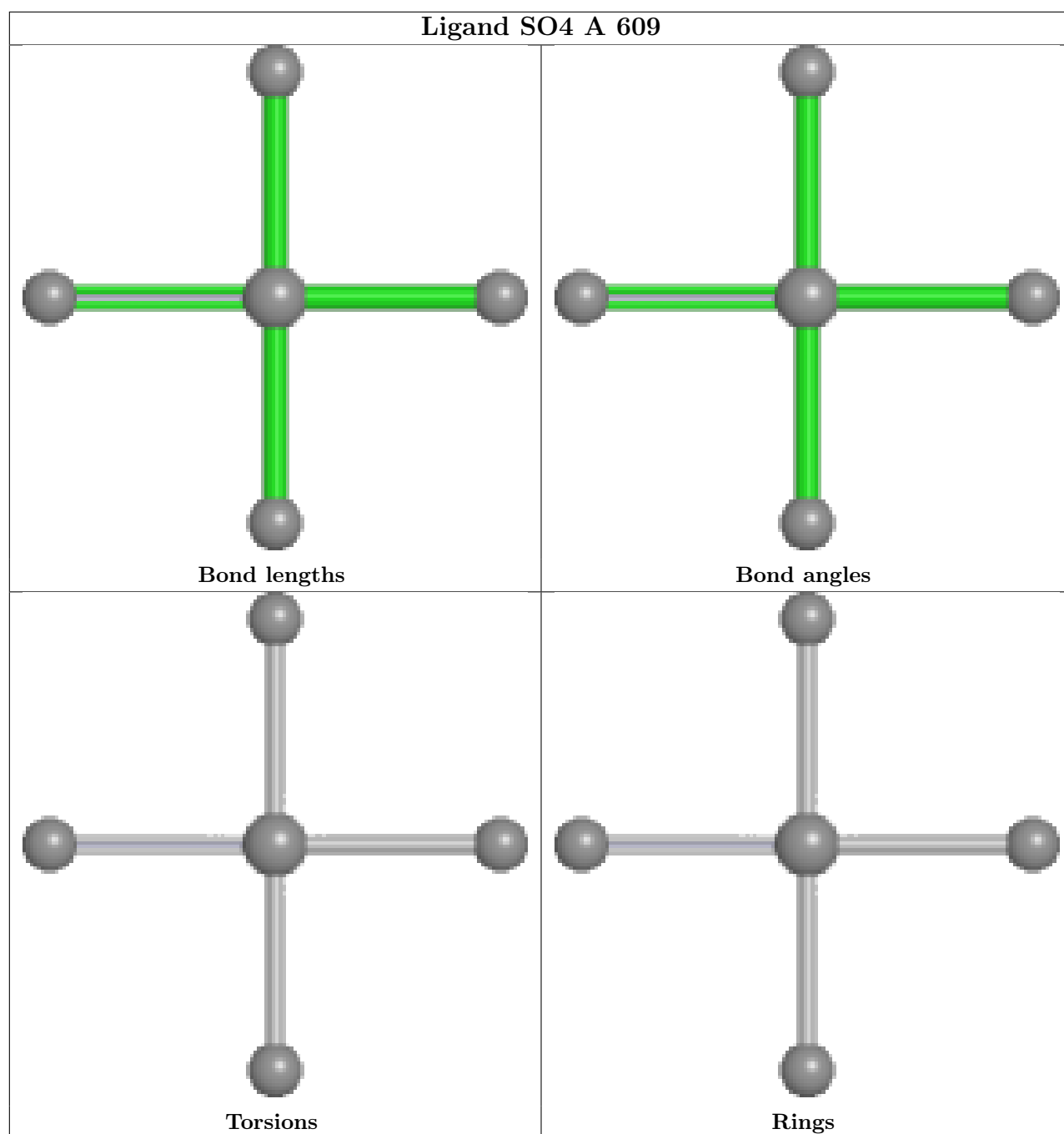
Mol	Chain	Res	Type	Atoms
5	D	602	FRU	O1-C1-C2-O2
7	A	604	TRS	C2-C-C1-O1
7	A	604	TRS	C3-C-C1-O1
7	A	604	TRS	N-C-C1-O1
7	B	608	TRS	N-C-C1-O1
7	B	608	TRS	C3-C-C2-O2
7	B	608	TRS	N-C-C2-O2
7	D	605	TRS	C2-C-C1-O1
7	D	605	TRS	C3-C-C1-O1
7	D	605	TRS	N-C-C1-O1
7	D	605	TRS	C2-C-C3-O3
7	D	605	TRS	N-C-C3-O3
7	B	608	TRS	C2-C-C1-O1
7	C	602	TRS	C3-C-C2-O2
5	D	602	FRU	O1-C1-C2-C3
7	B	608	TRS	C1-C-C2-O2
7	C	602	TRS	N-C-C2-O2
7	D	605	TRS	C1-C-C3-O3
4	B	601	GLC	O5-C5-C6-O6
4	D	601	GLC	C4-C5-C6-O6
4	B	601	GLC	C4-C5-C6-O6
7	B	608	TRS	C3-C-C1-O1
7	C	602	TRS	C1-C-C2-O2
5	D	602	FRU	O1-C1-C2-O5
7	C	602	TRS	C3-C-C1-O1
7	D	605	TRS	C1-C-C2-O2
5	A	602	FRU	O1-C1-C2-C3
7	C	602	TRS	N-C-C1-O1
5	B	602	FRU	O1-C1-C2-C3
5	C	608	FRU	O1-C1-C2-C3
4	D	601	GLC	O5-C5-C6-O6

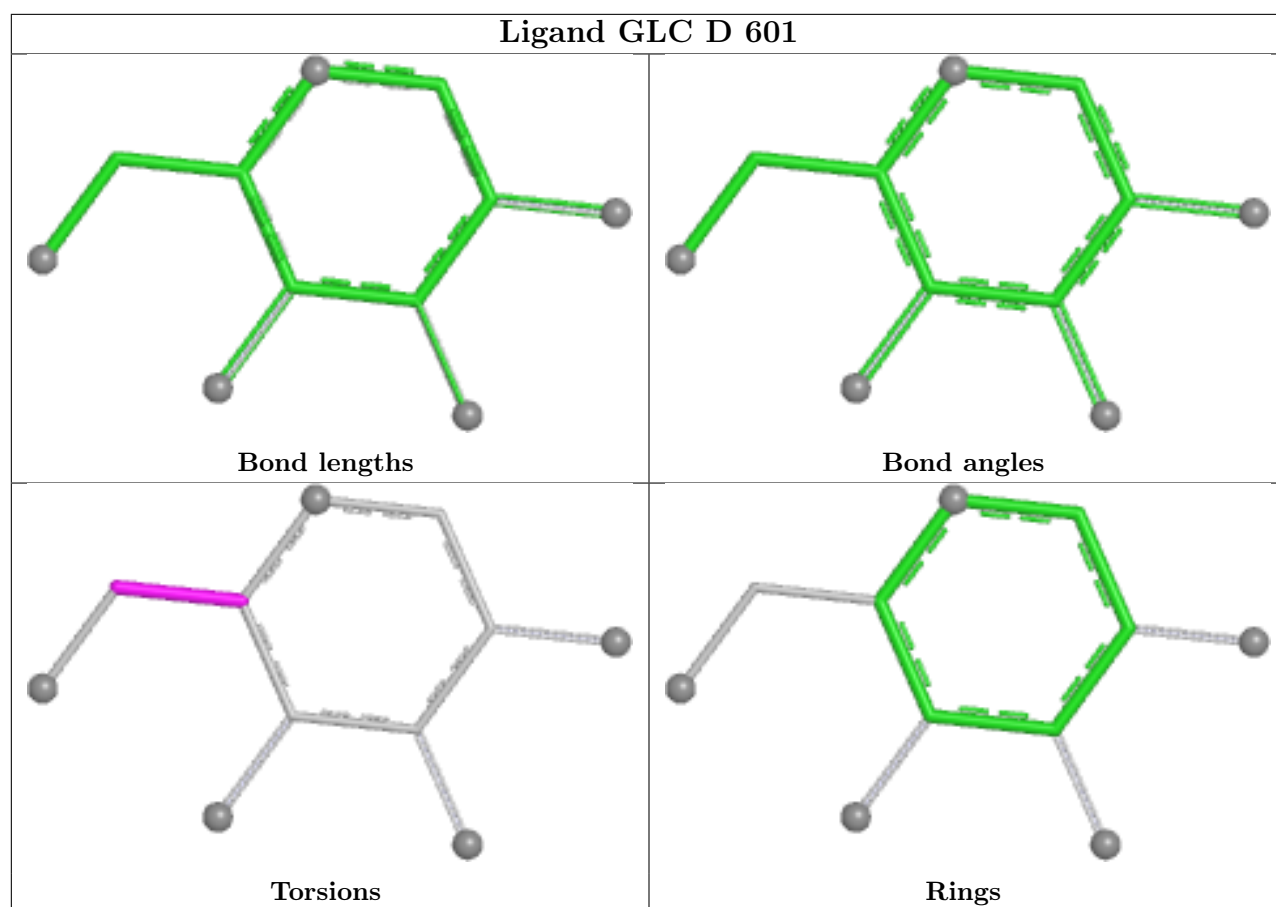
There are no ring outliers.

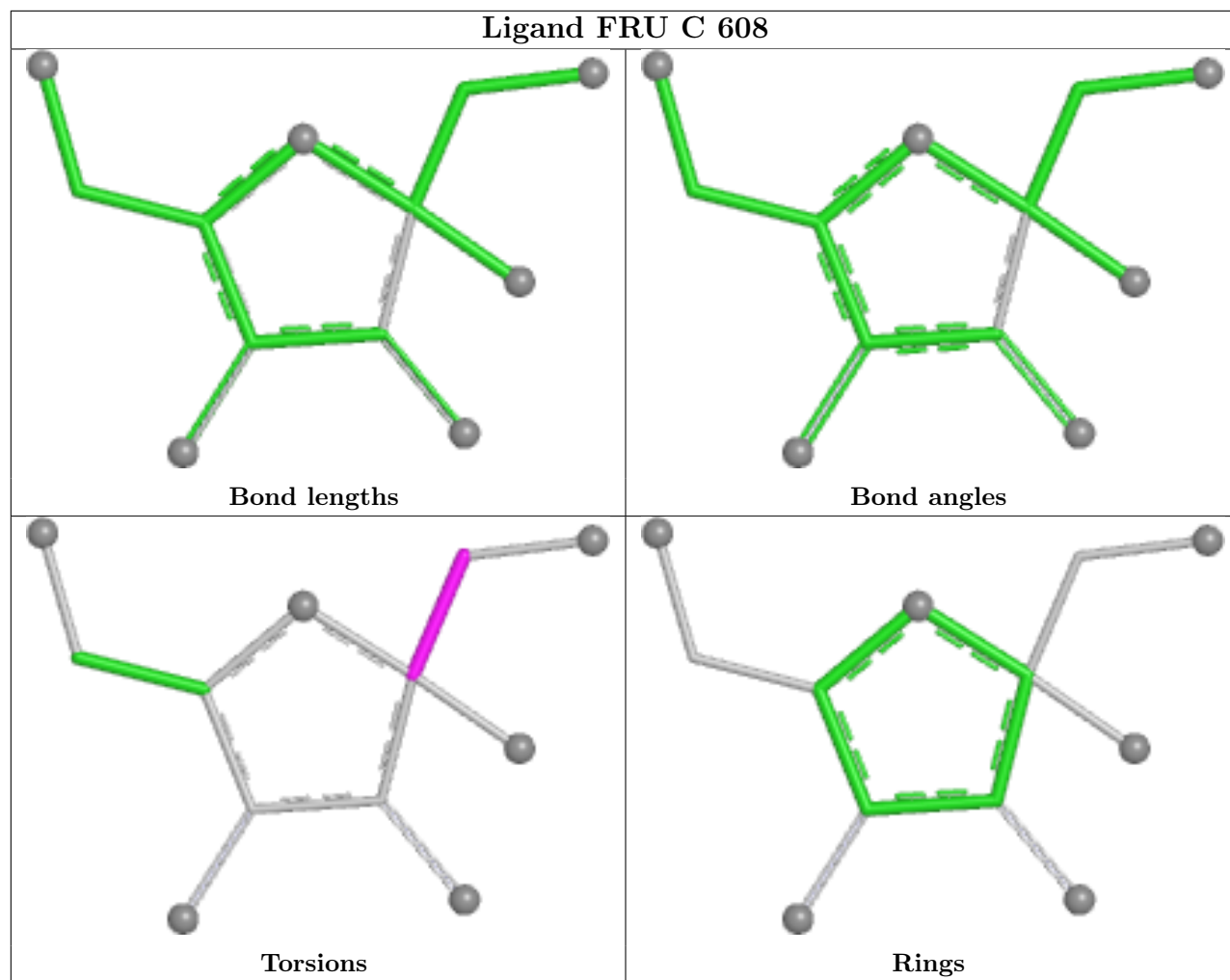
9 monomers are involved in 12 short contacts:

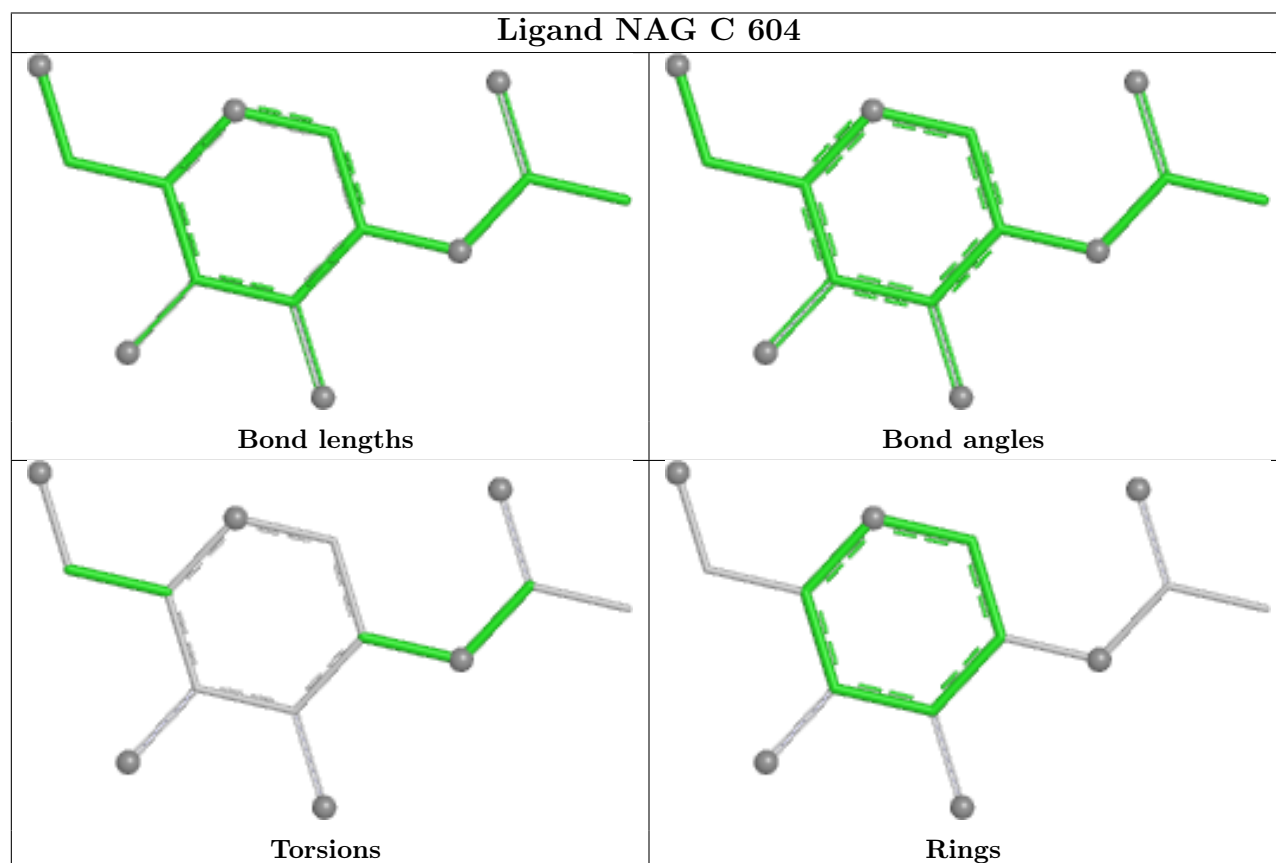
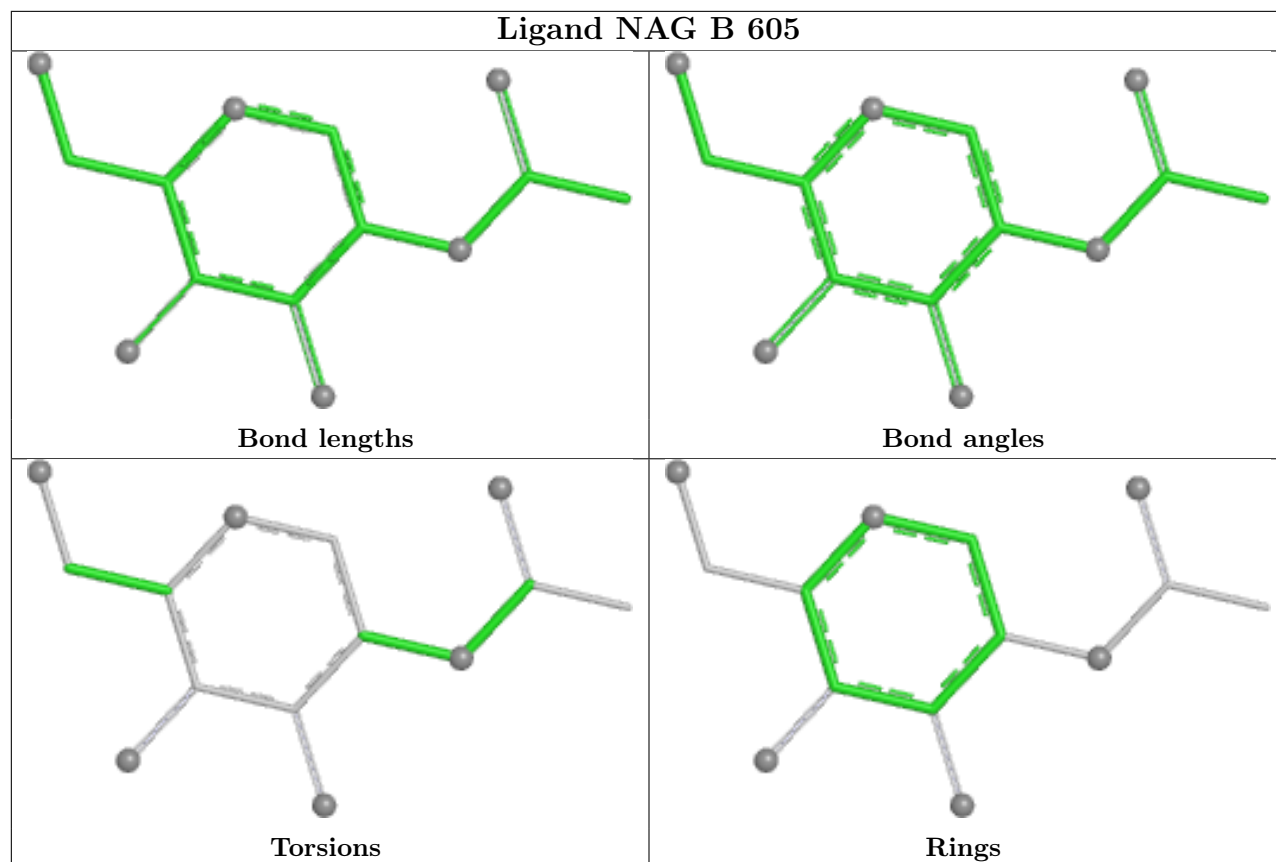
Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	D	601	GLC	2	0
5	C	608	FRU	2	0
9	B	607	GOL	1	0
4	C	607	GLC	2	0
4	B	601	GLC	4	0
5	A	602	FRU	2	0
4	A	601	GLC	2	0
5	B	602	FRU	3	0
5	D	602	FRU	3	0

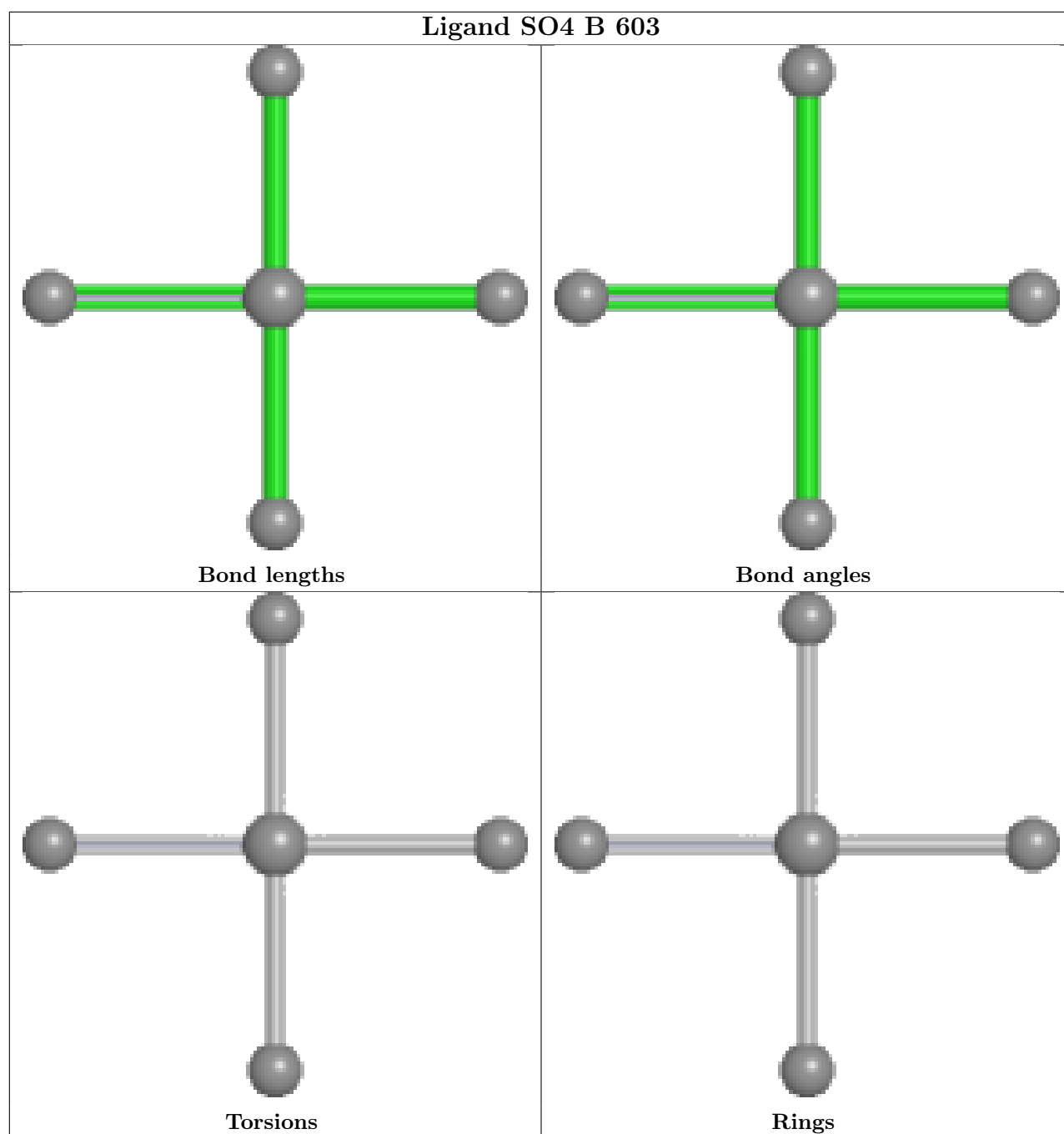
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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



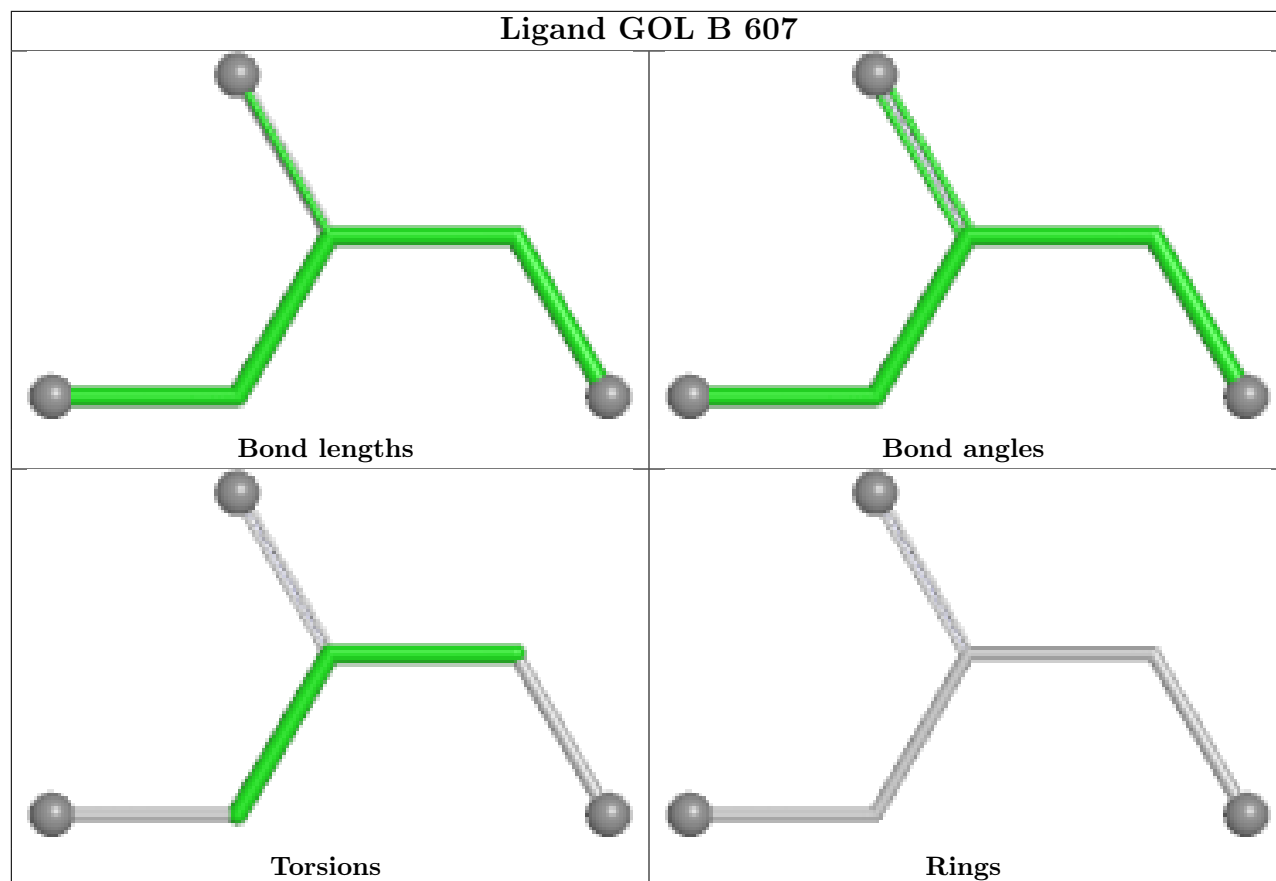




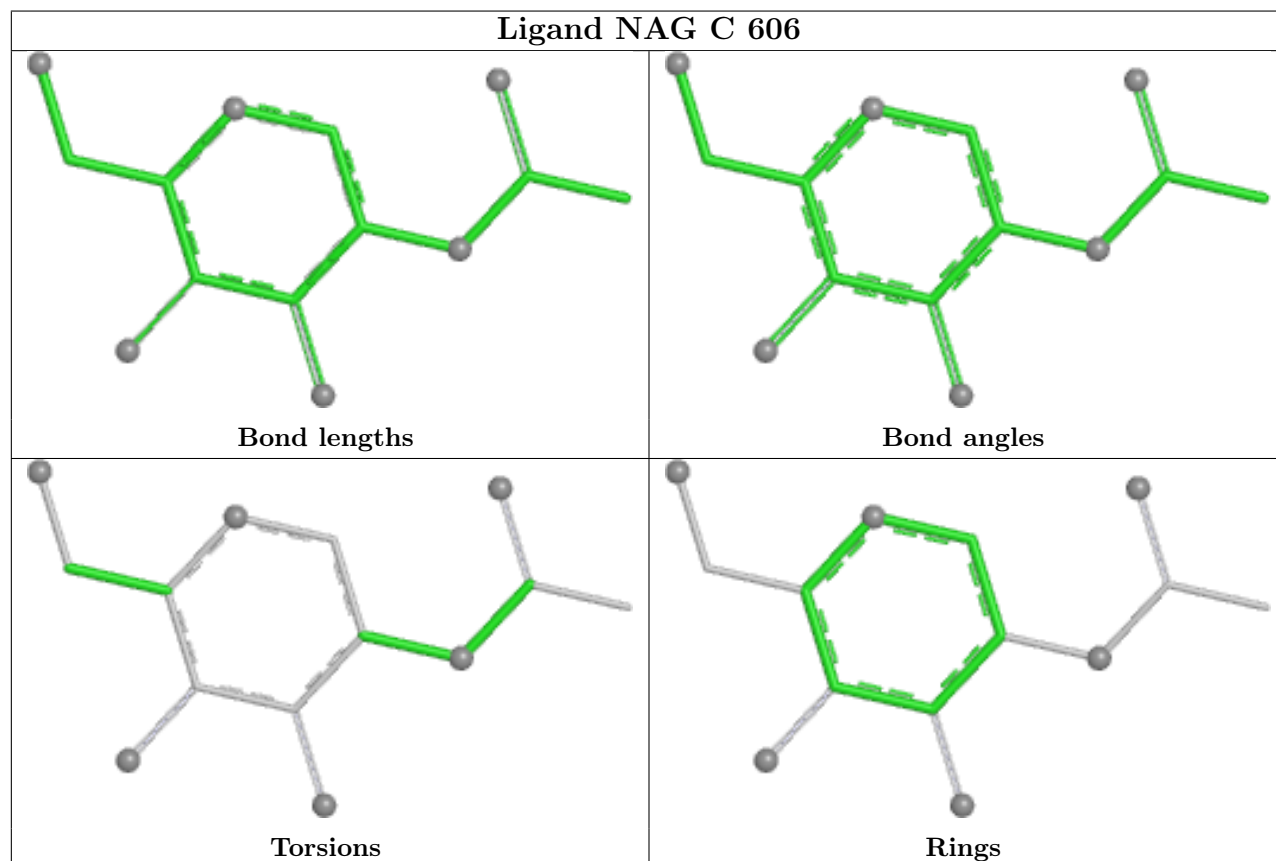


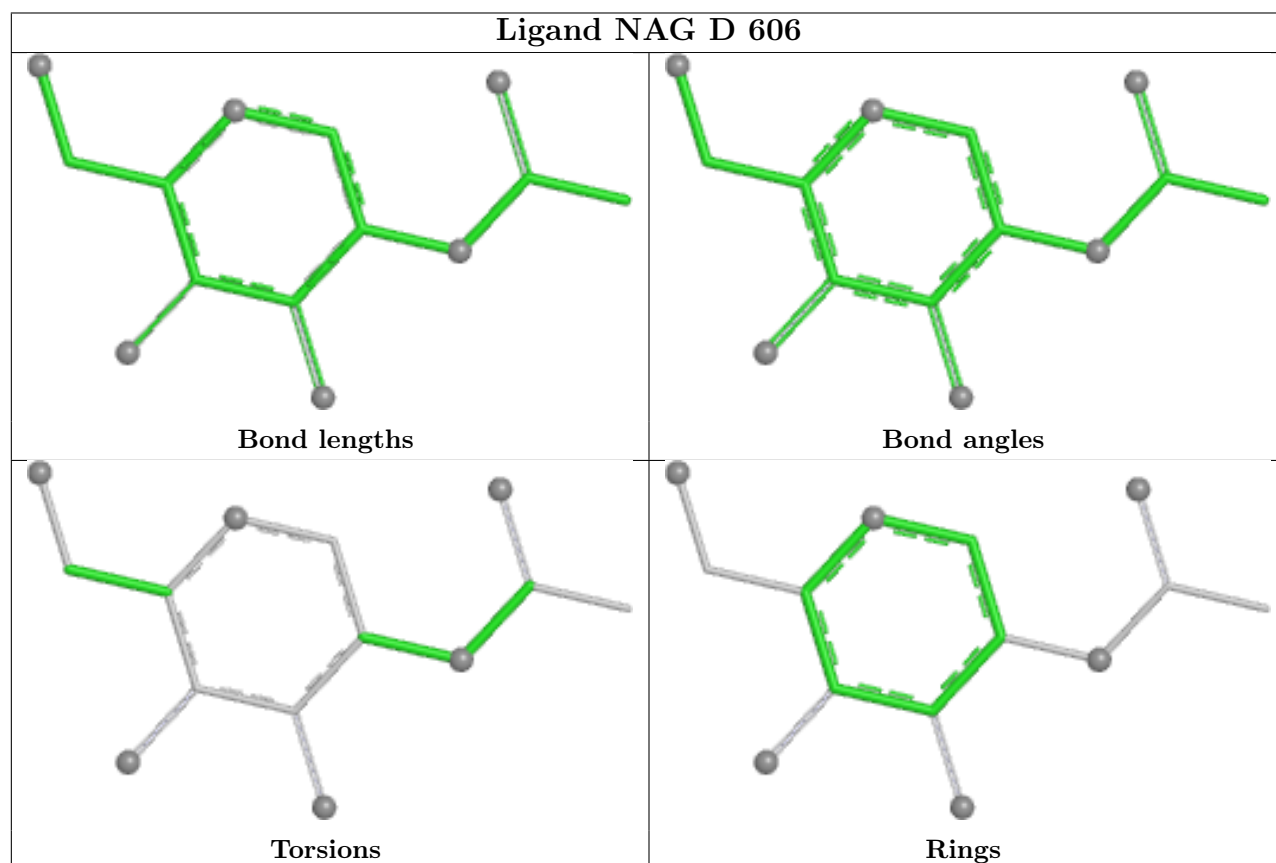
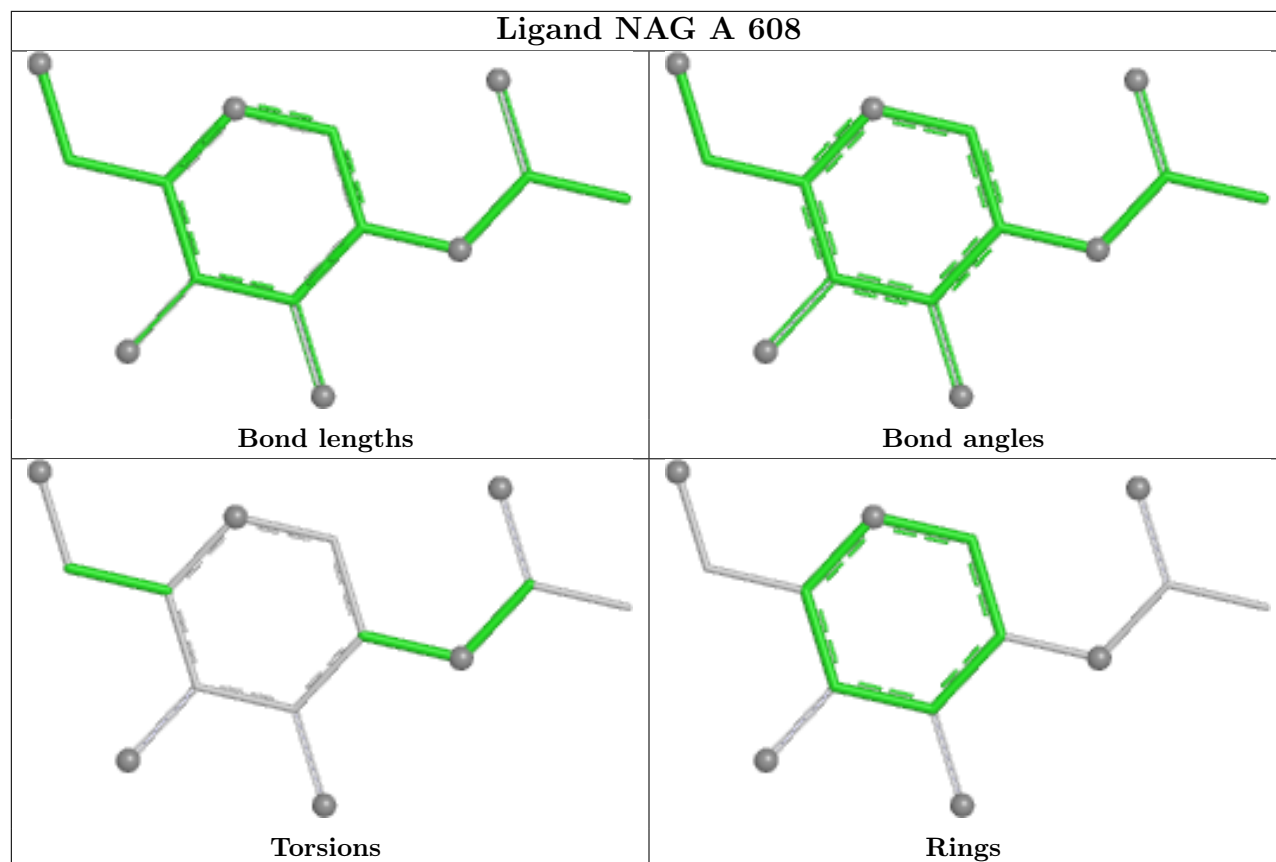


Ligand GOL B 607

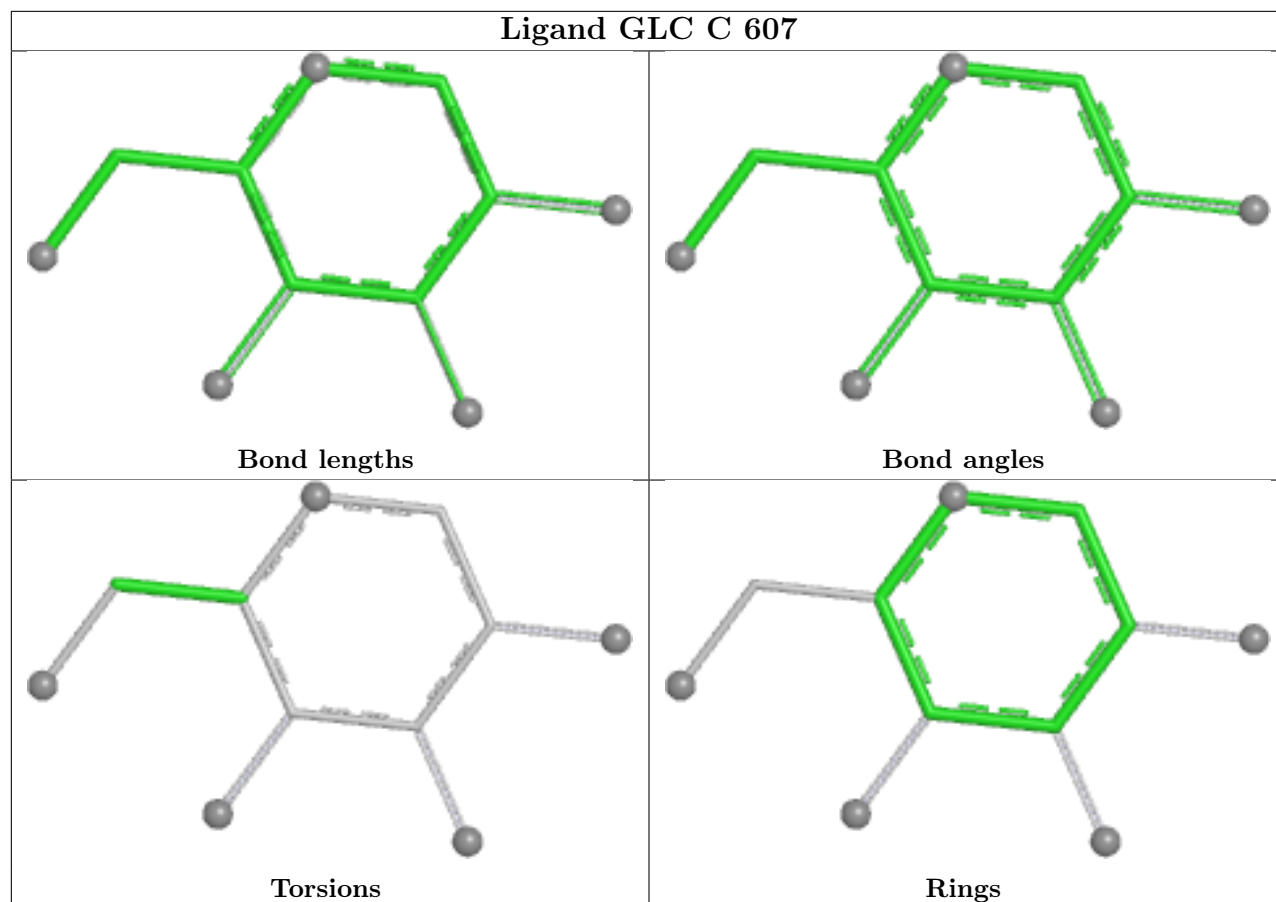


Ligand NAG C 606

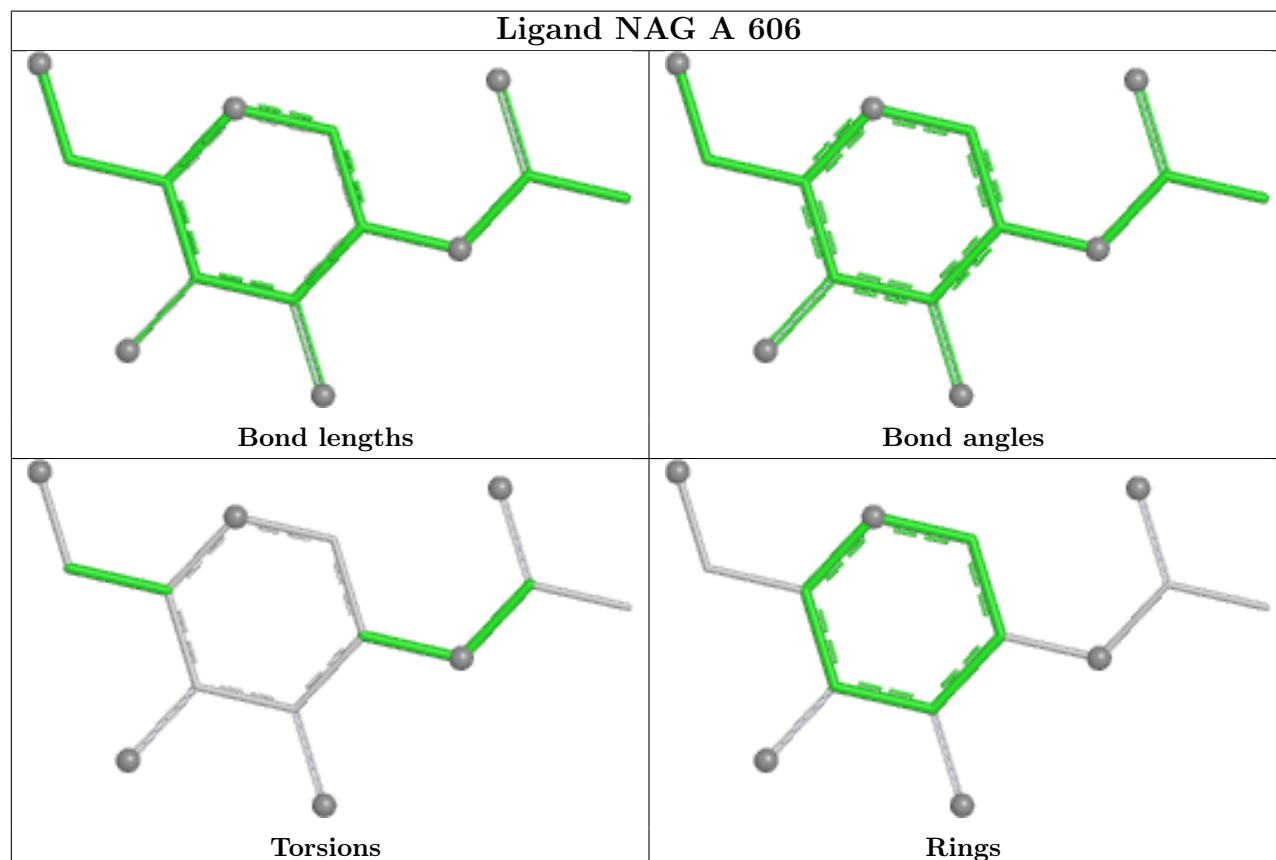


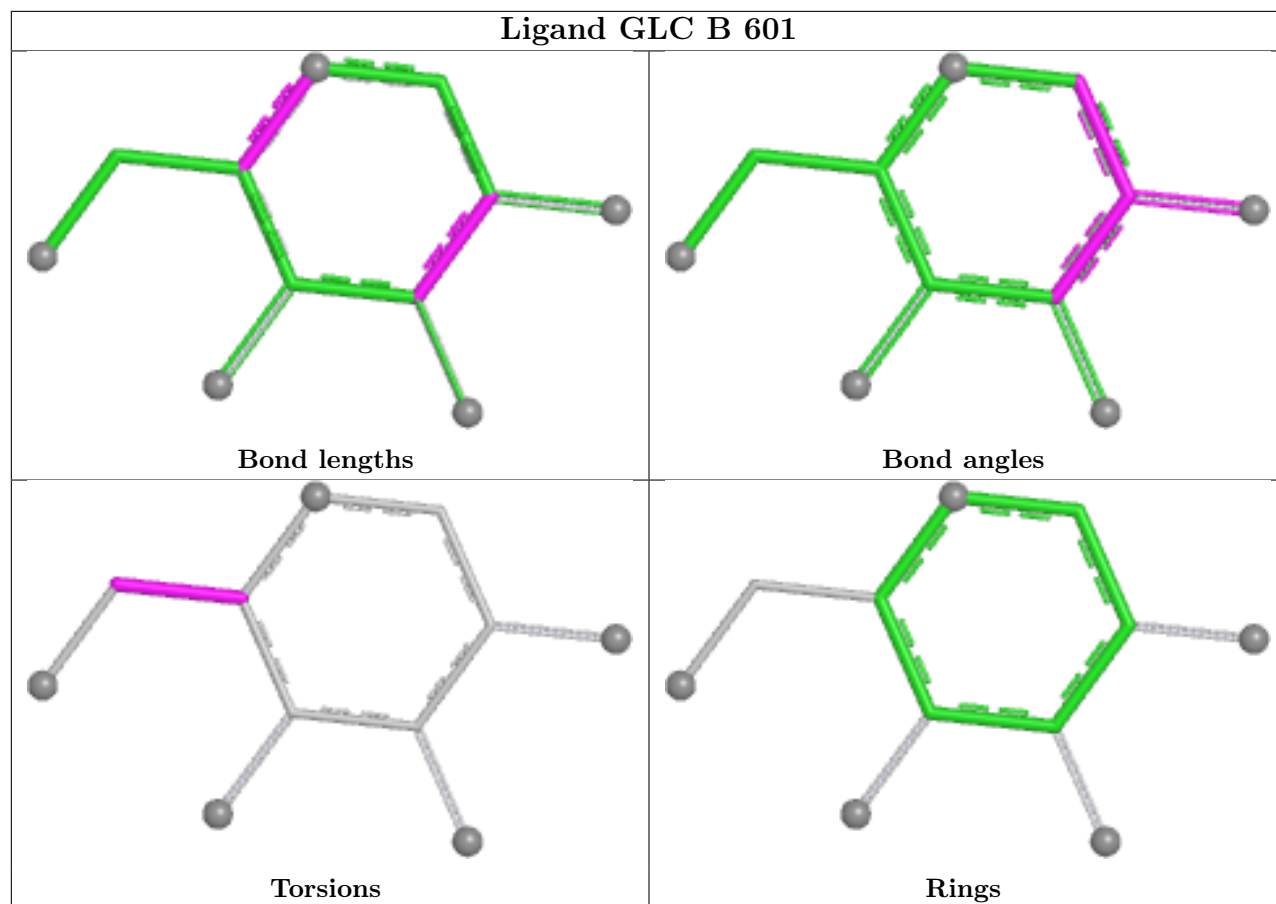


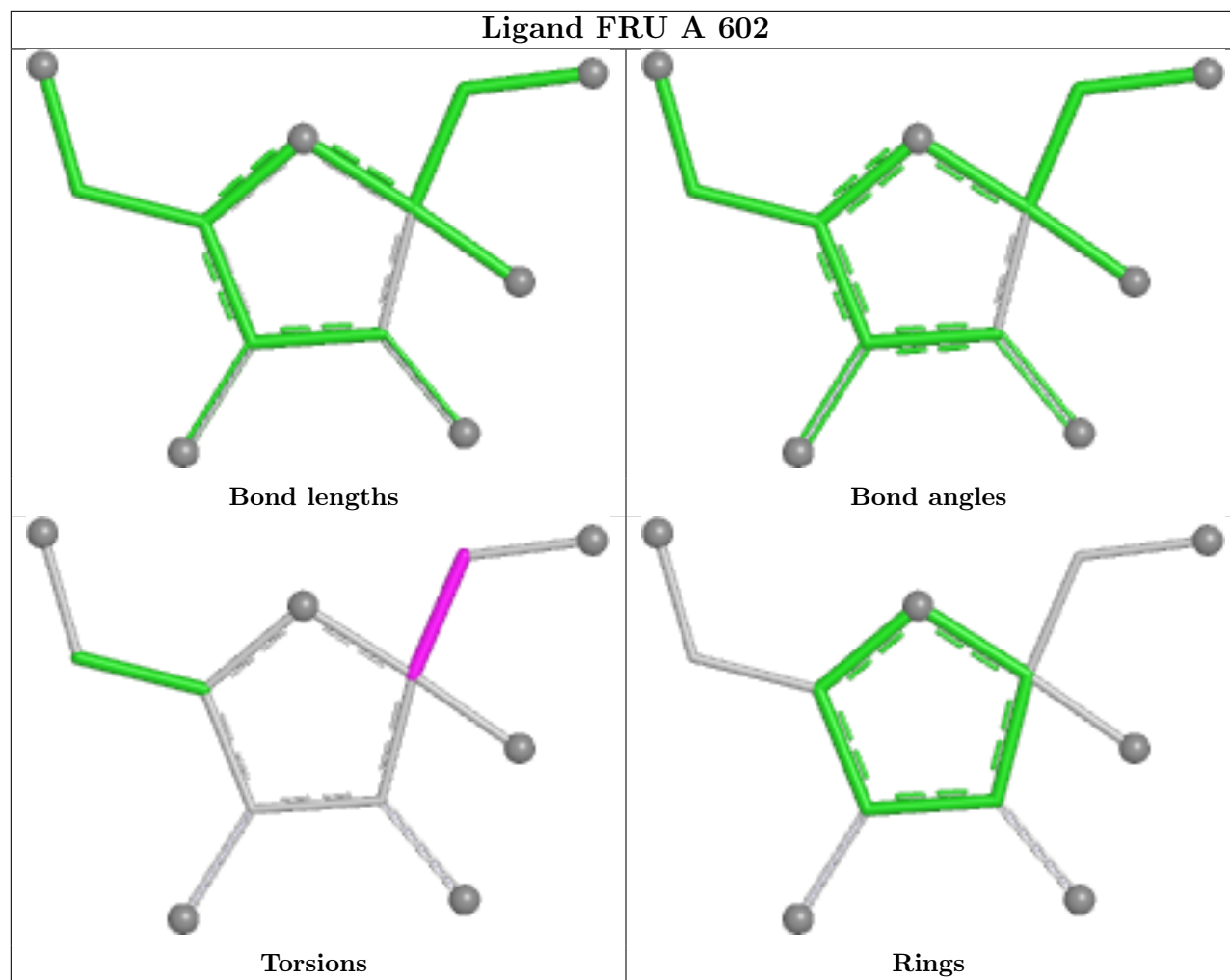
Ligand GLC C 607

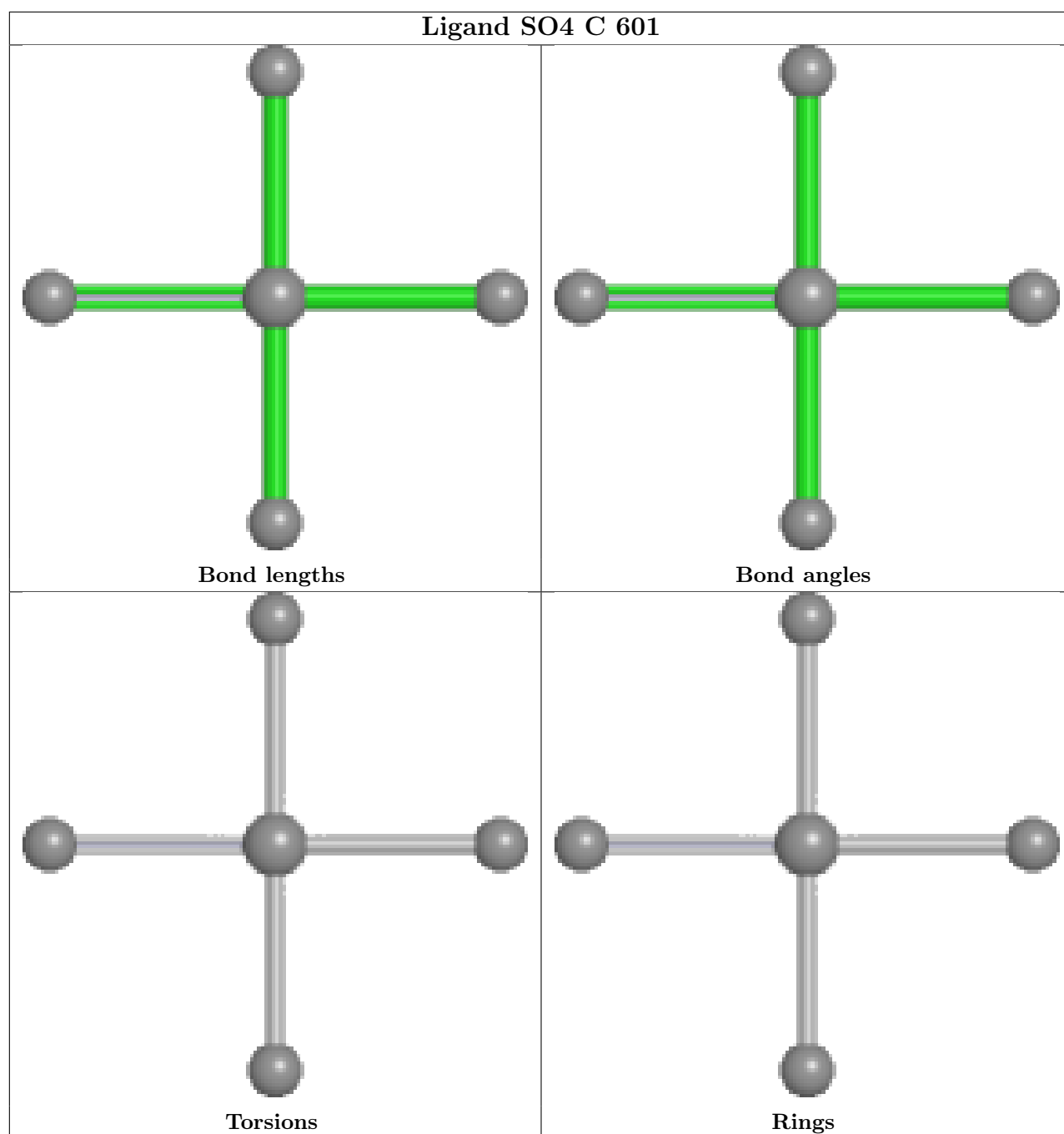


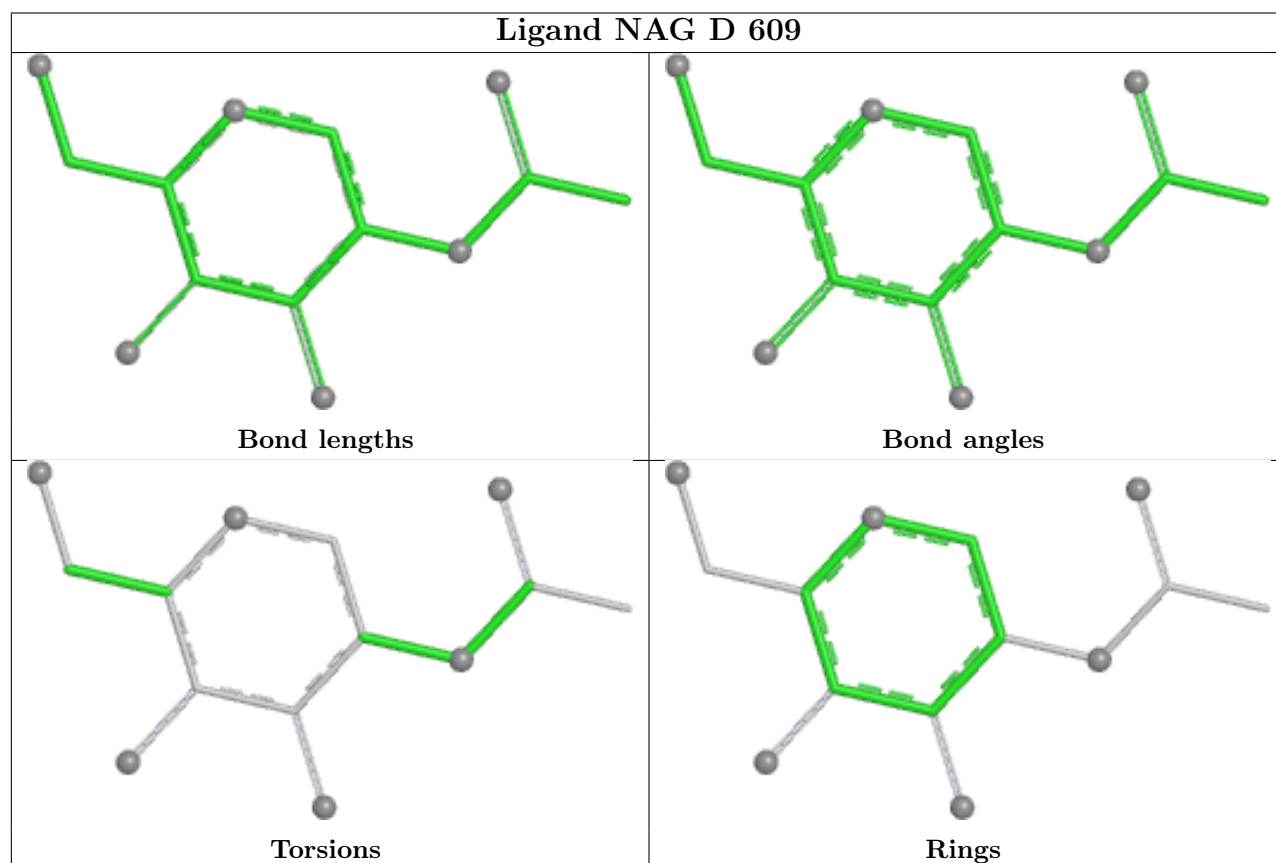
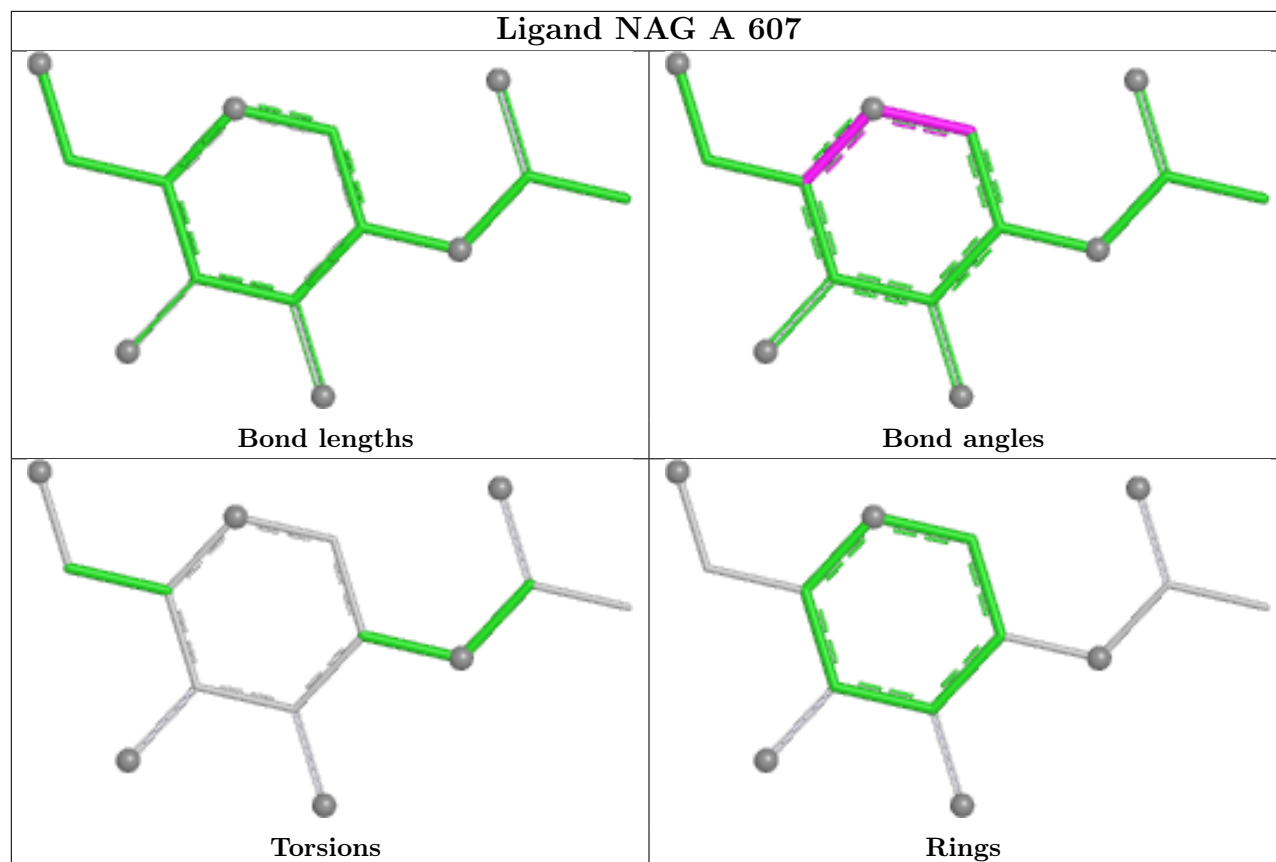
Ligand NAG A 606

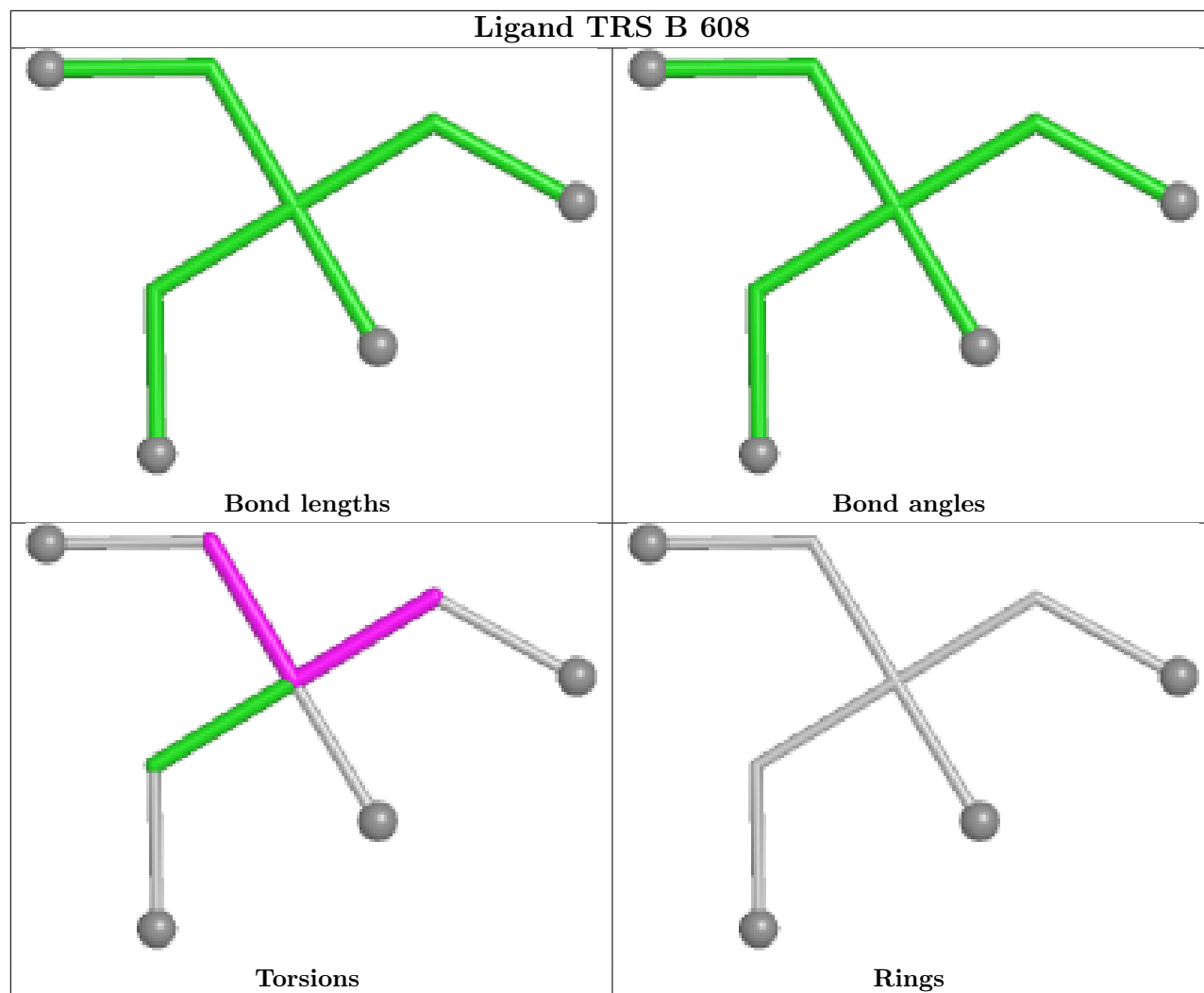


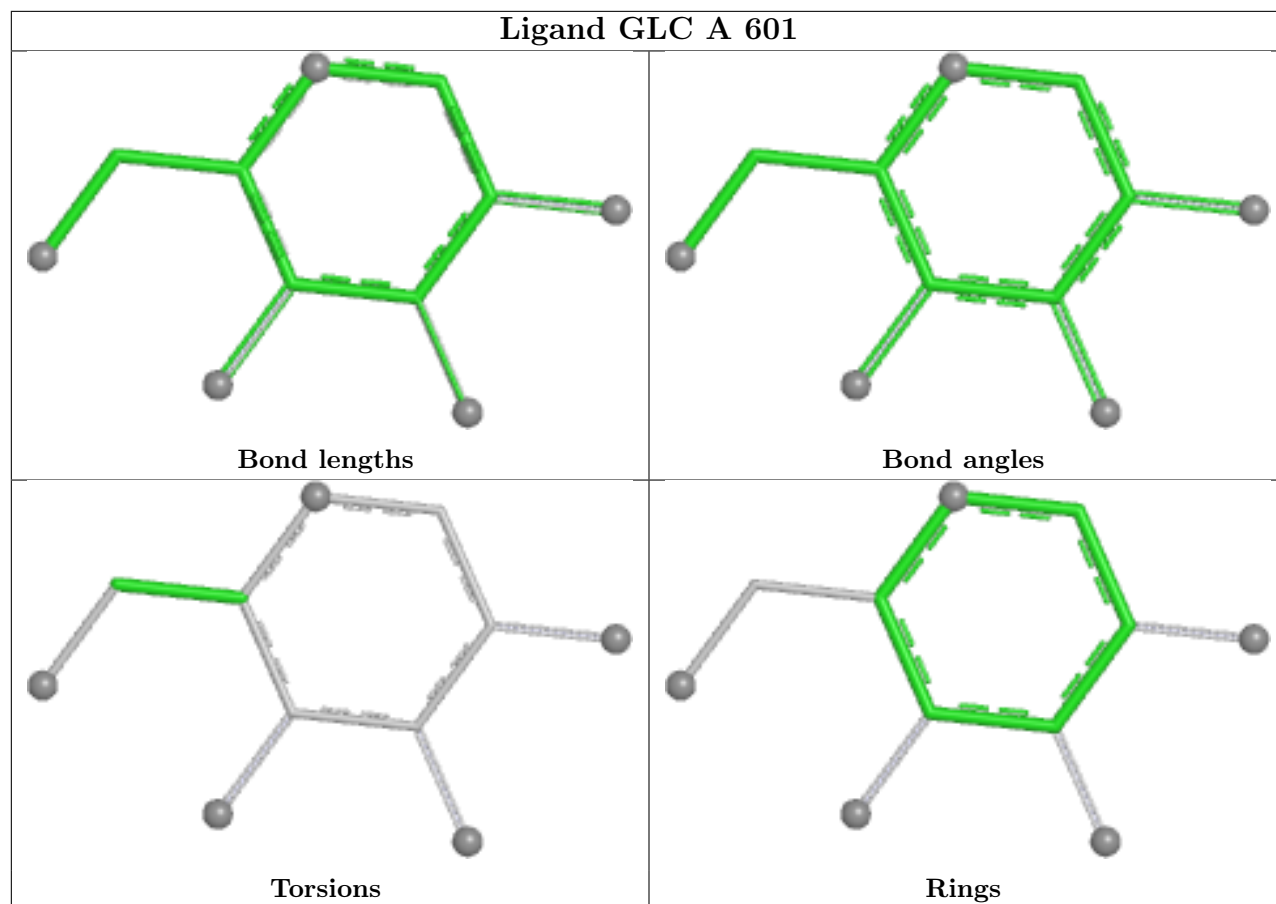


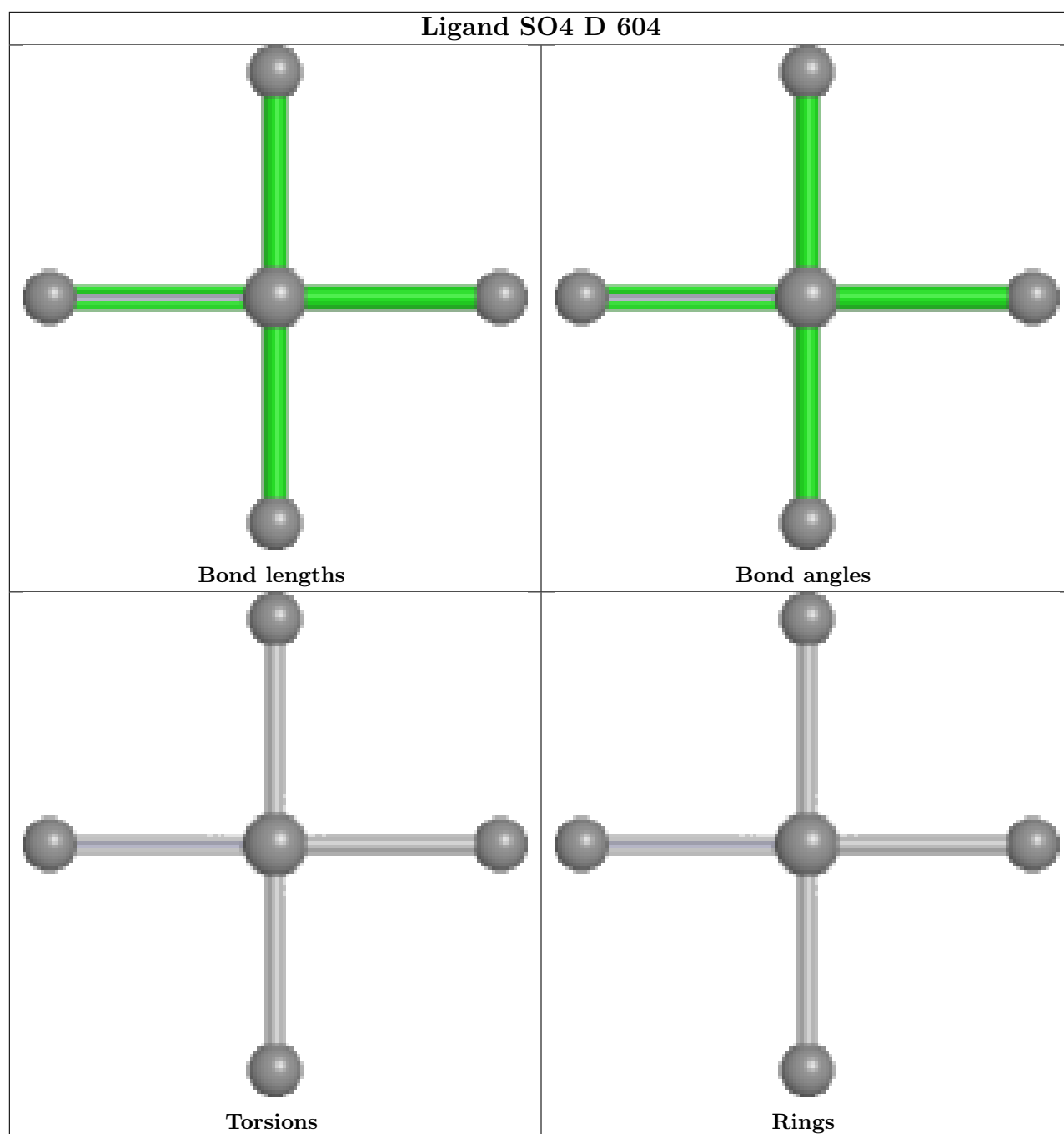


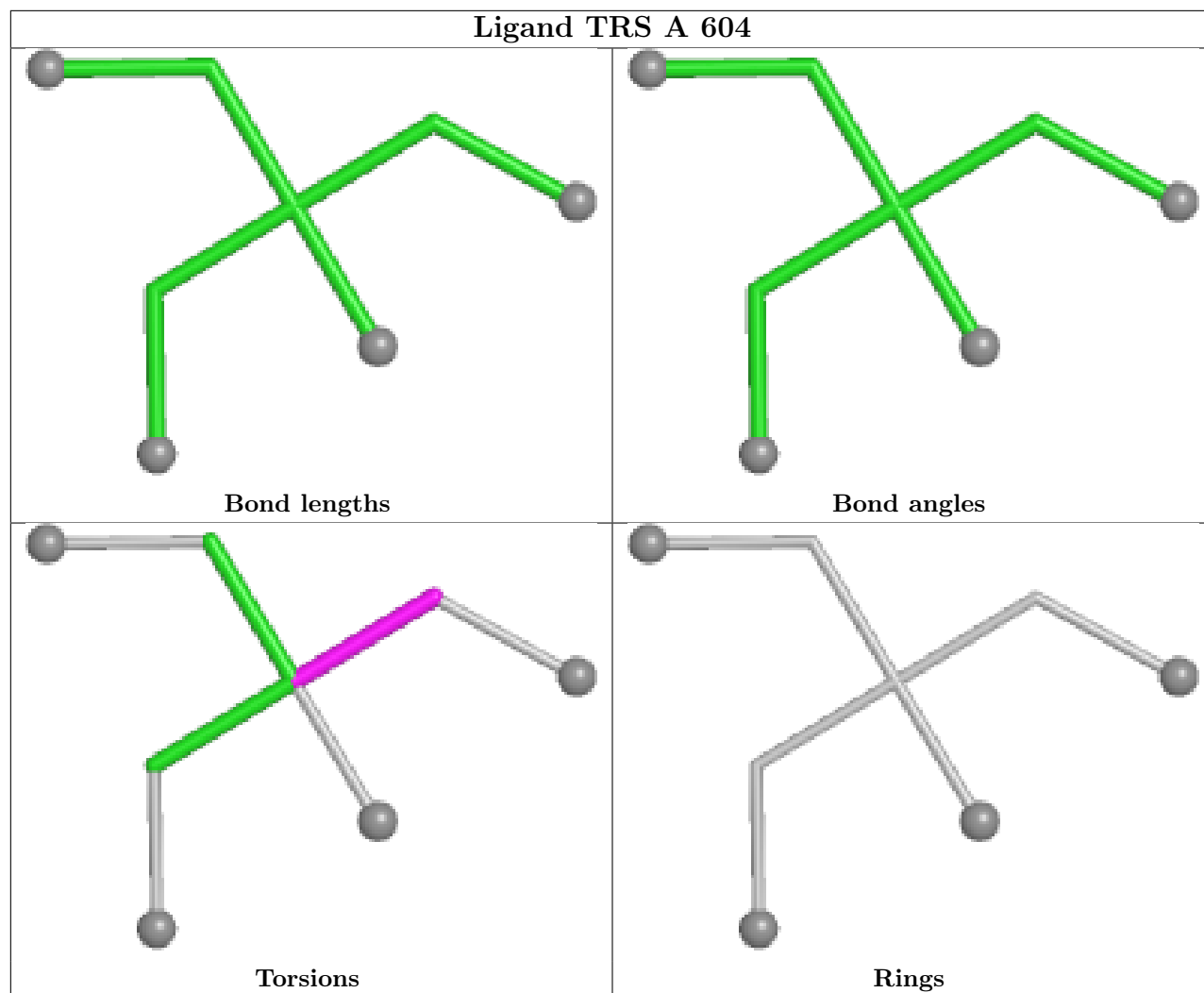


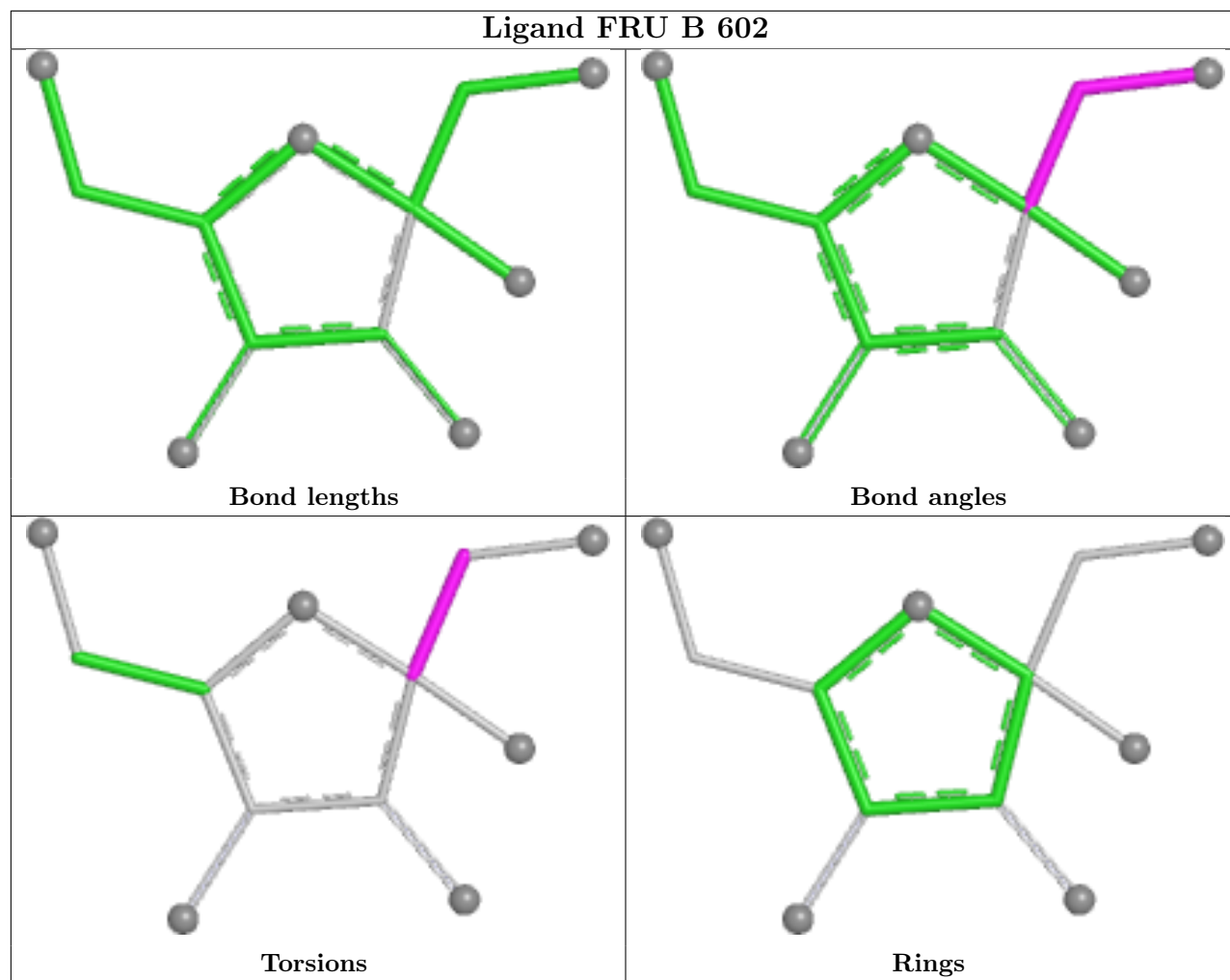


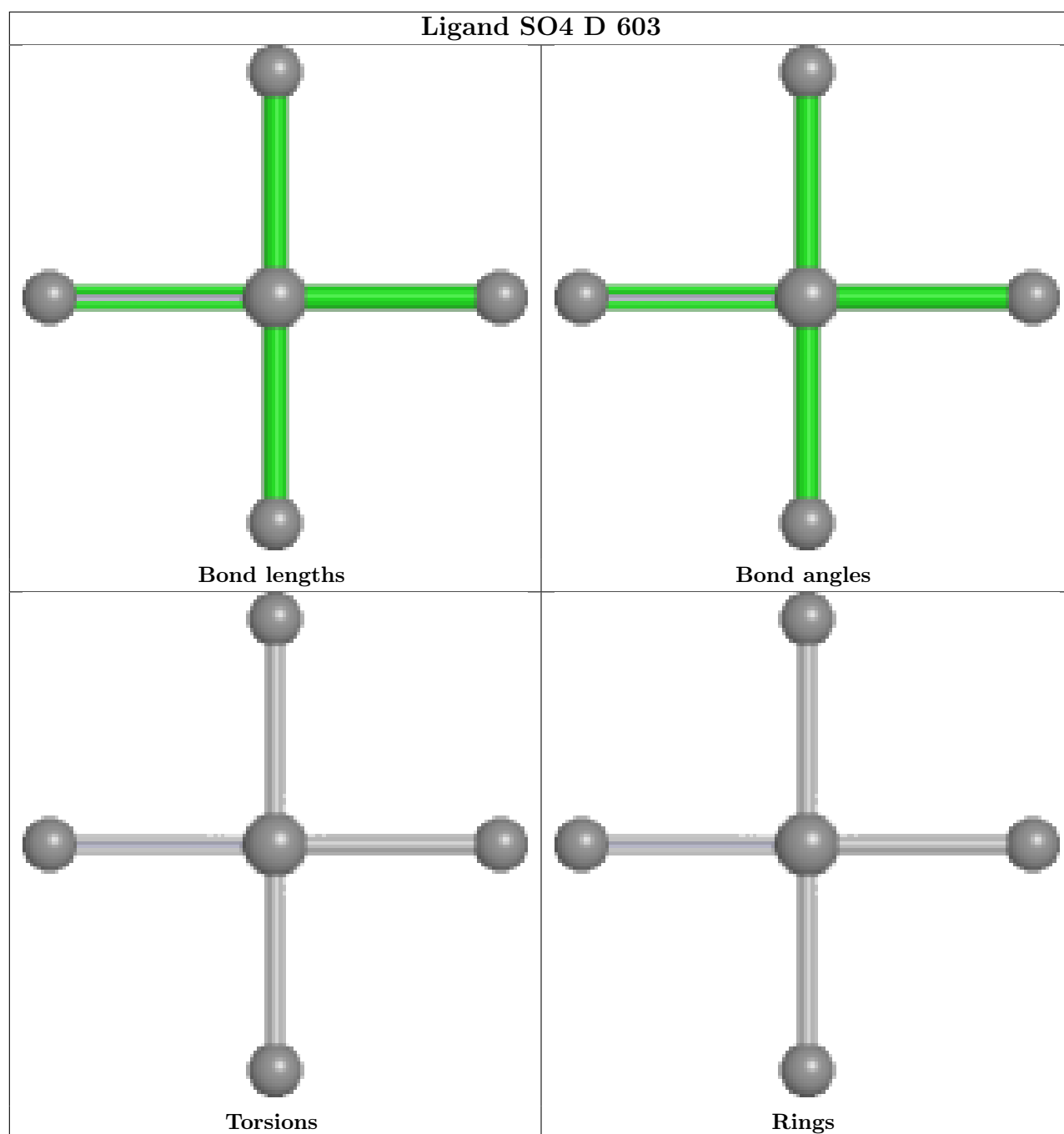


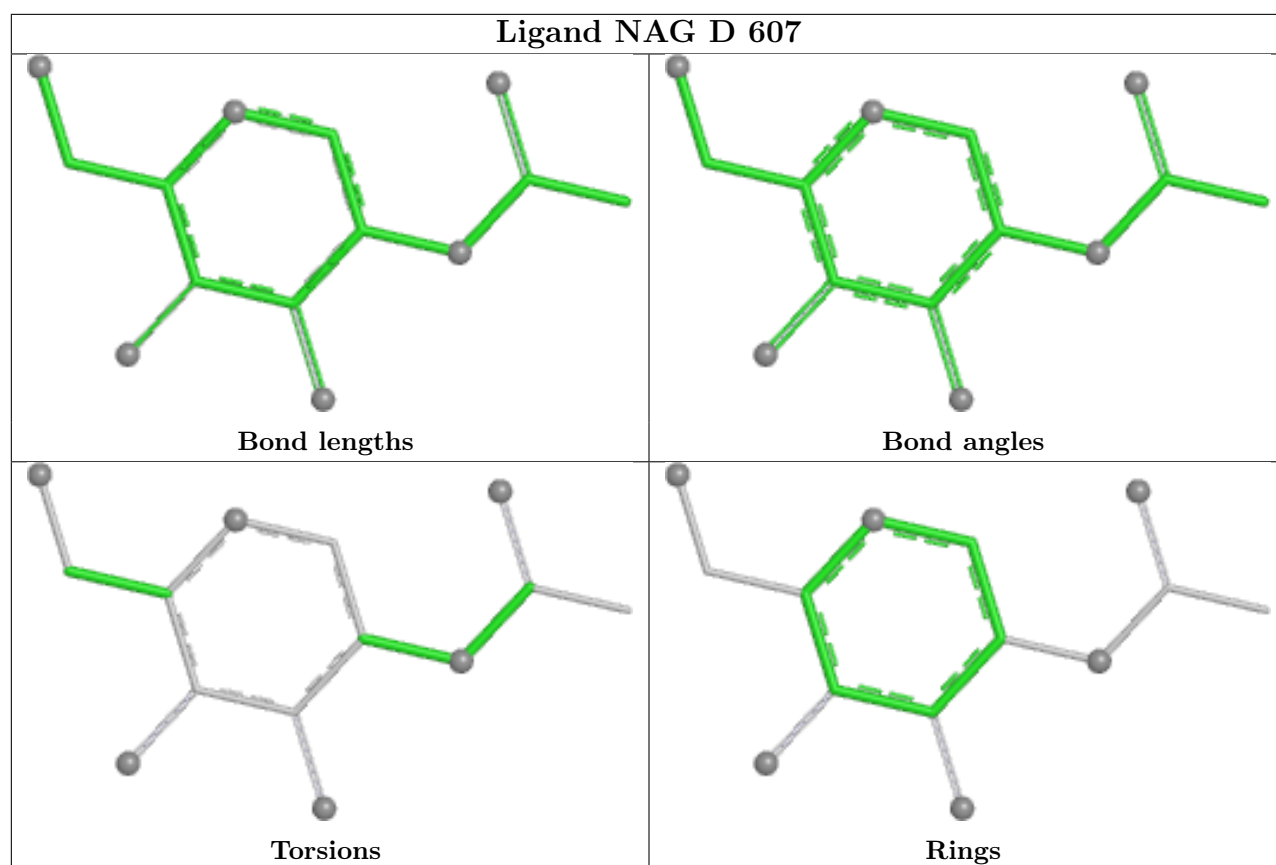


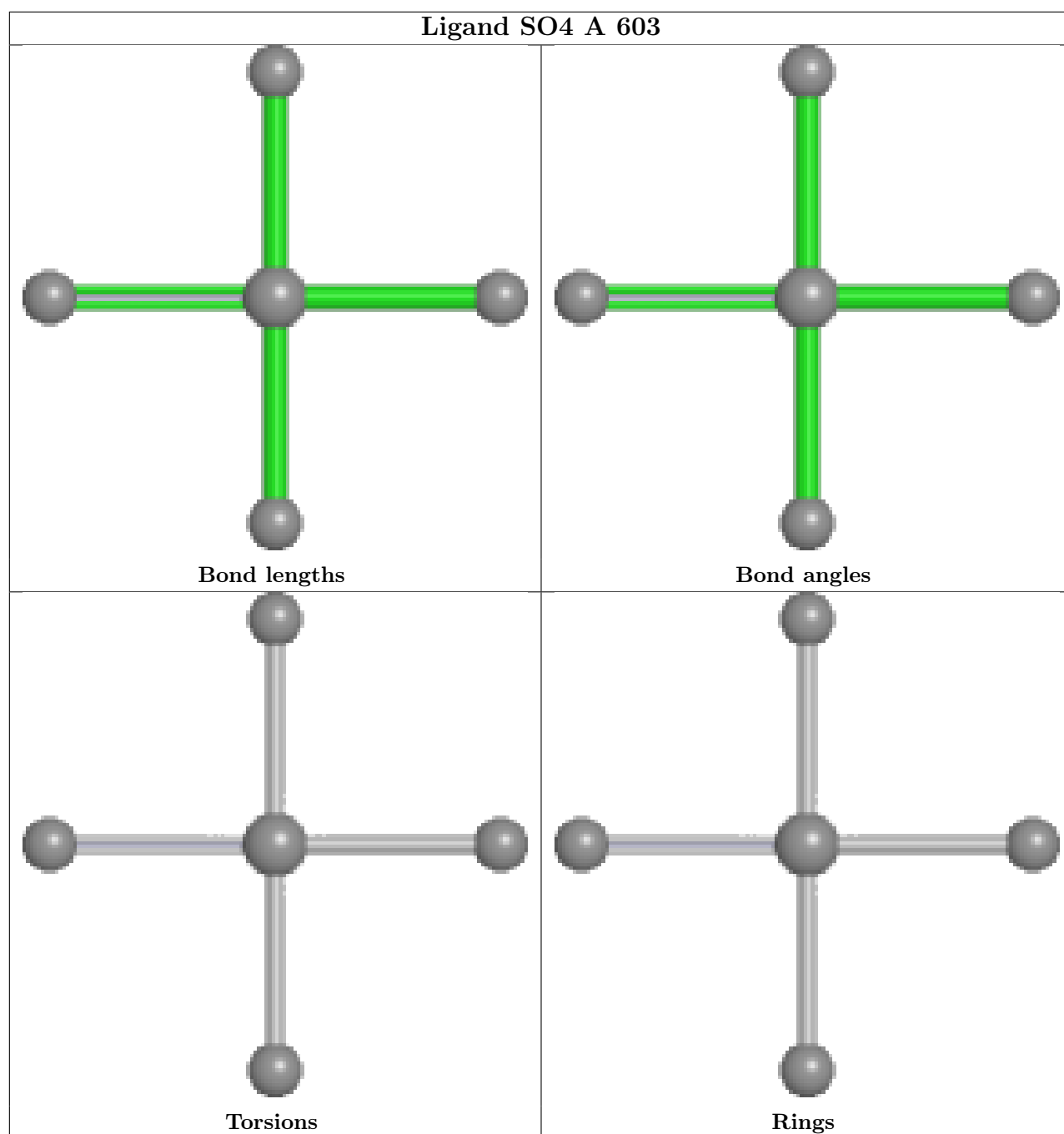


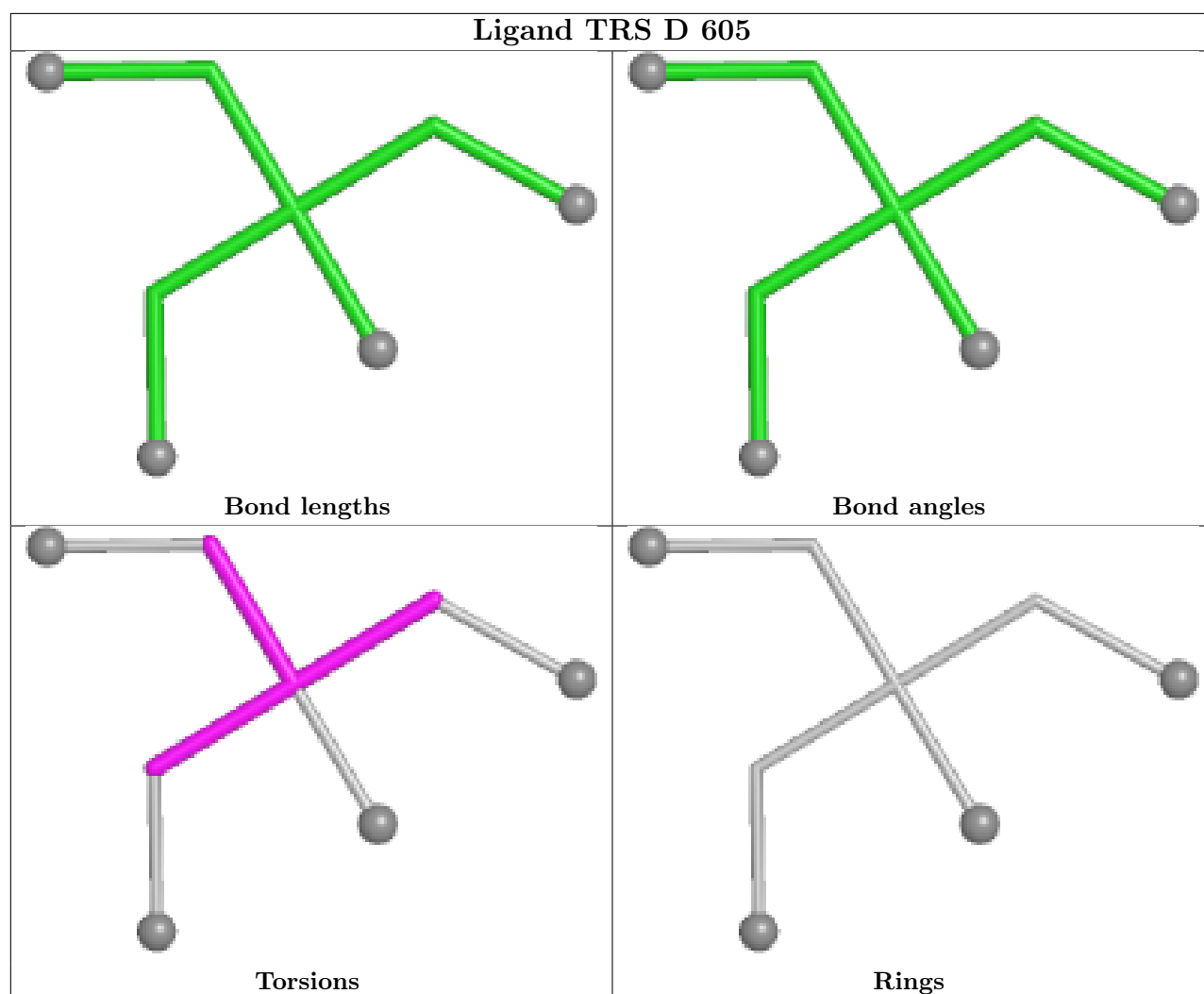


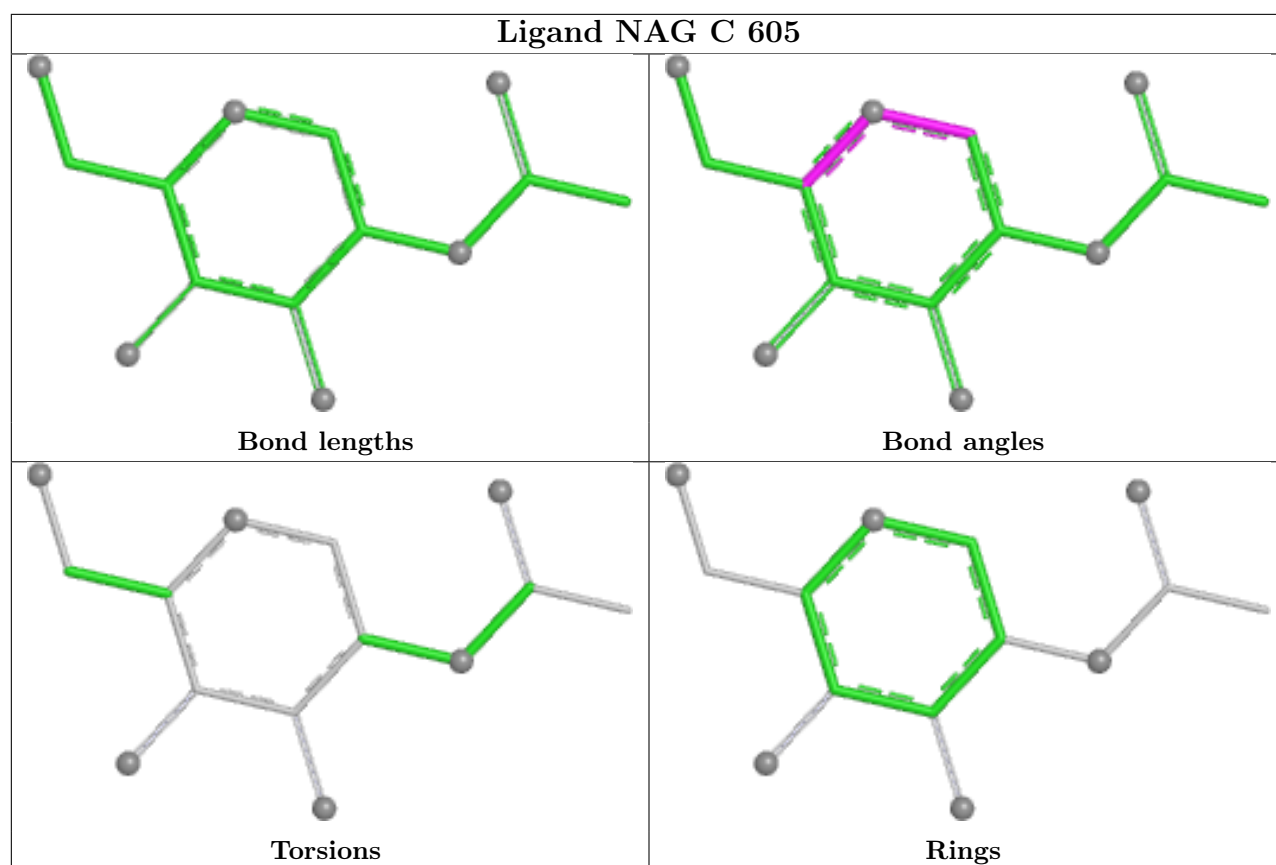


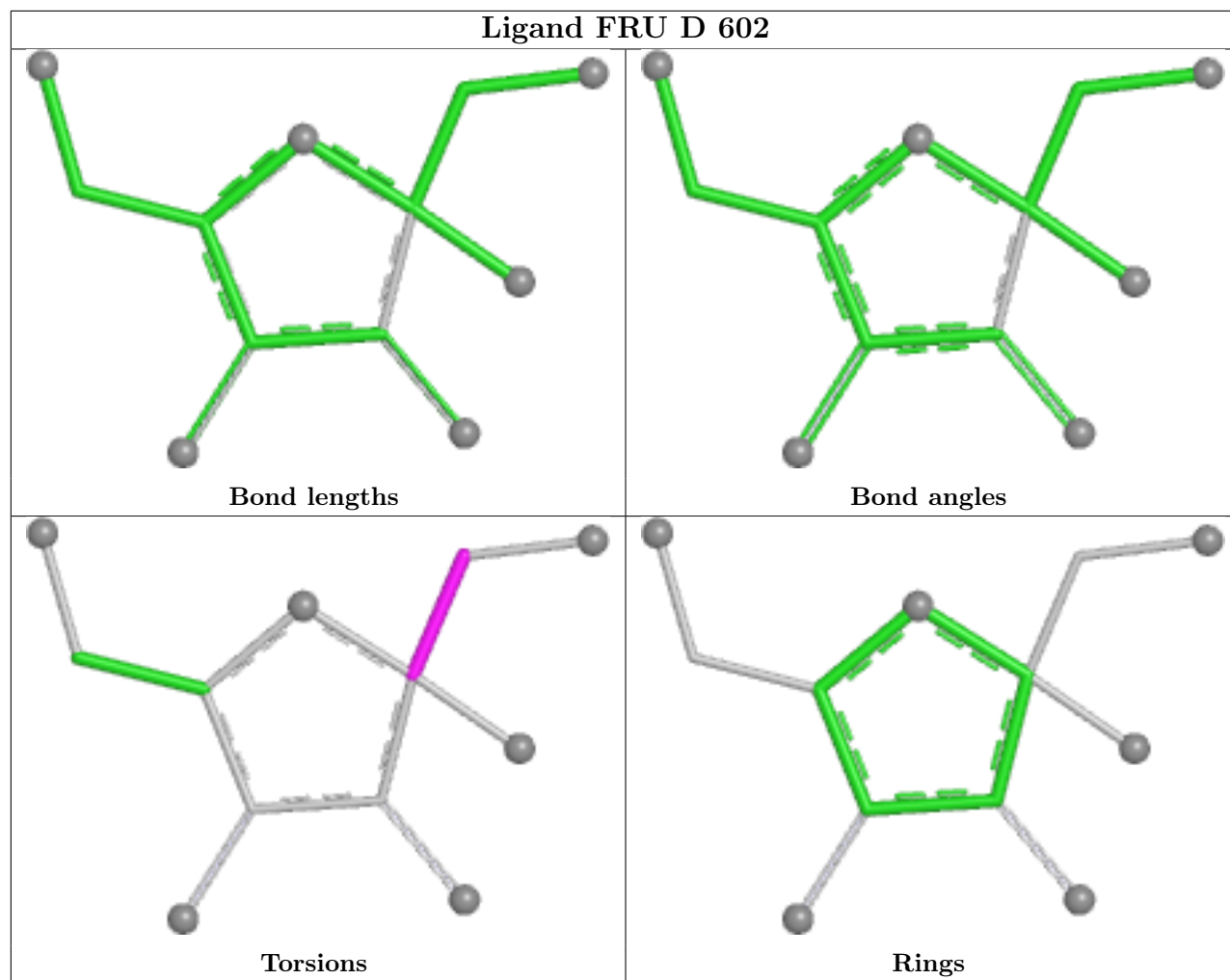


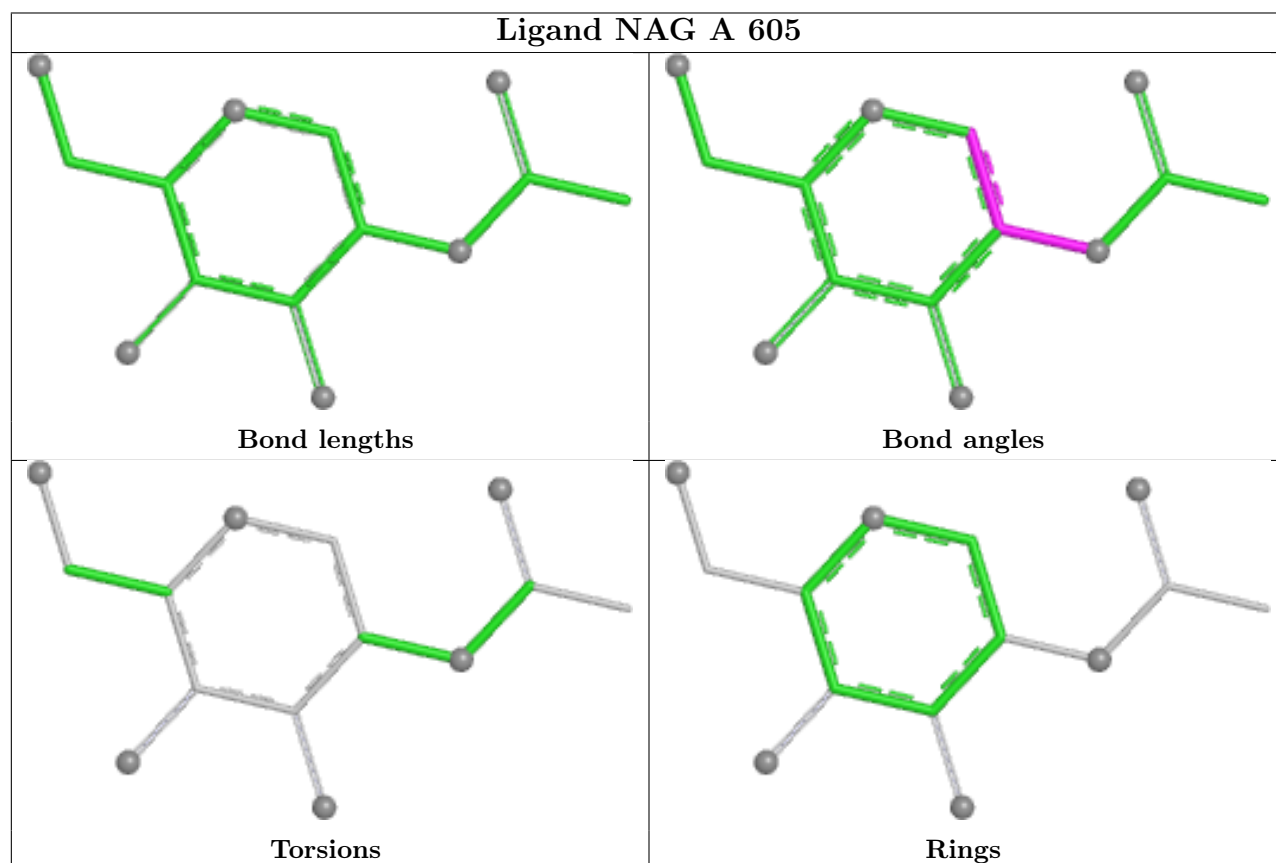
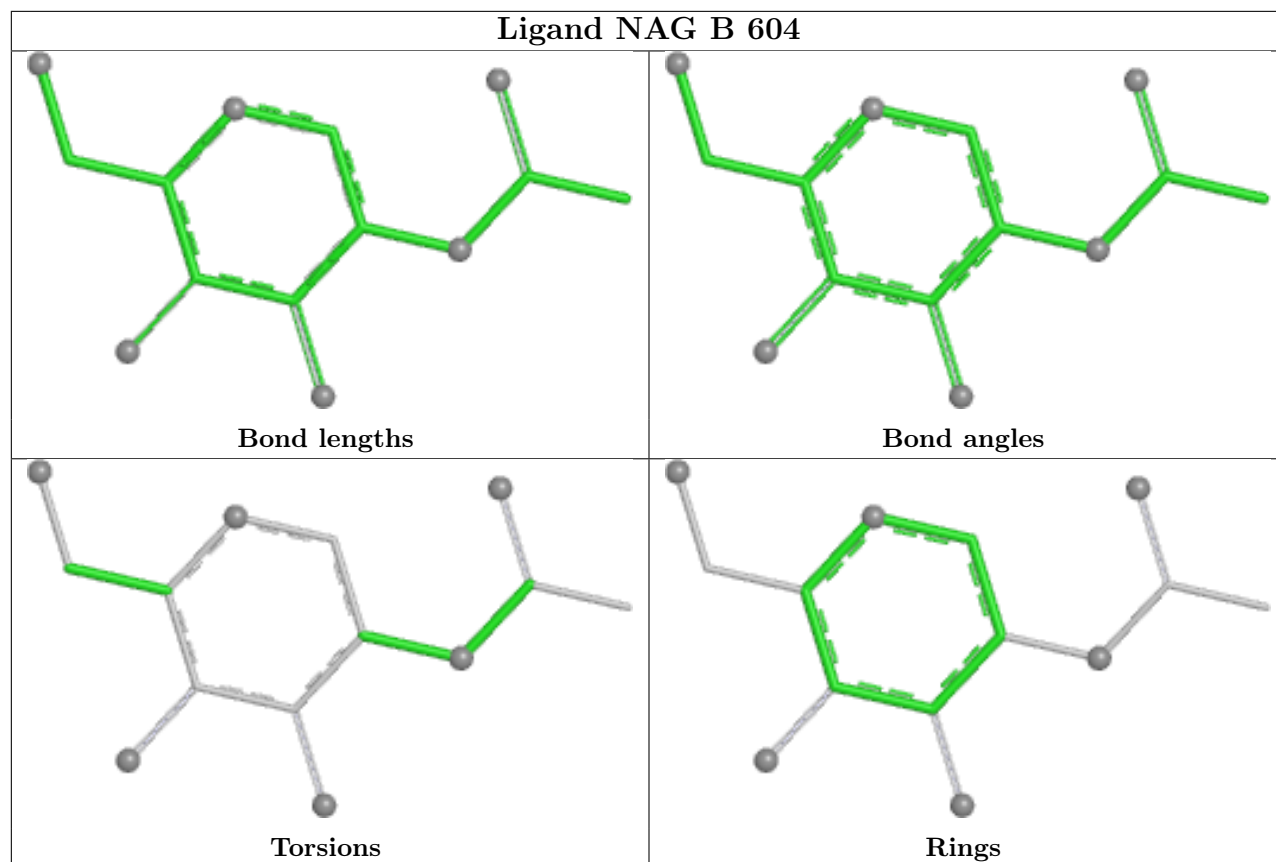


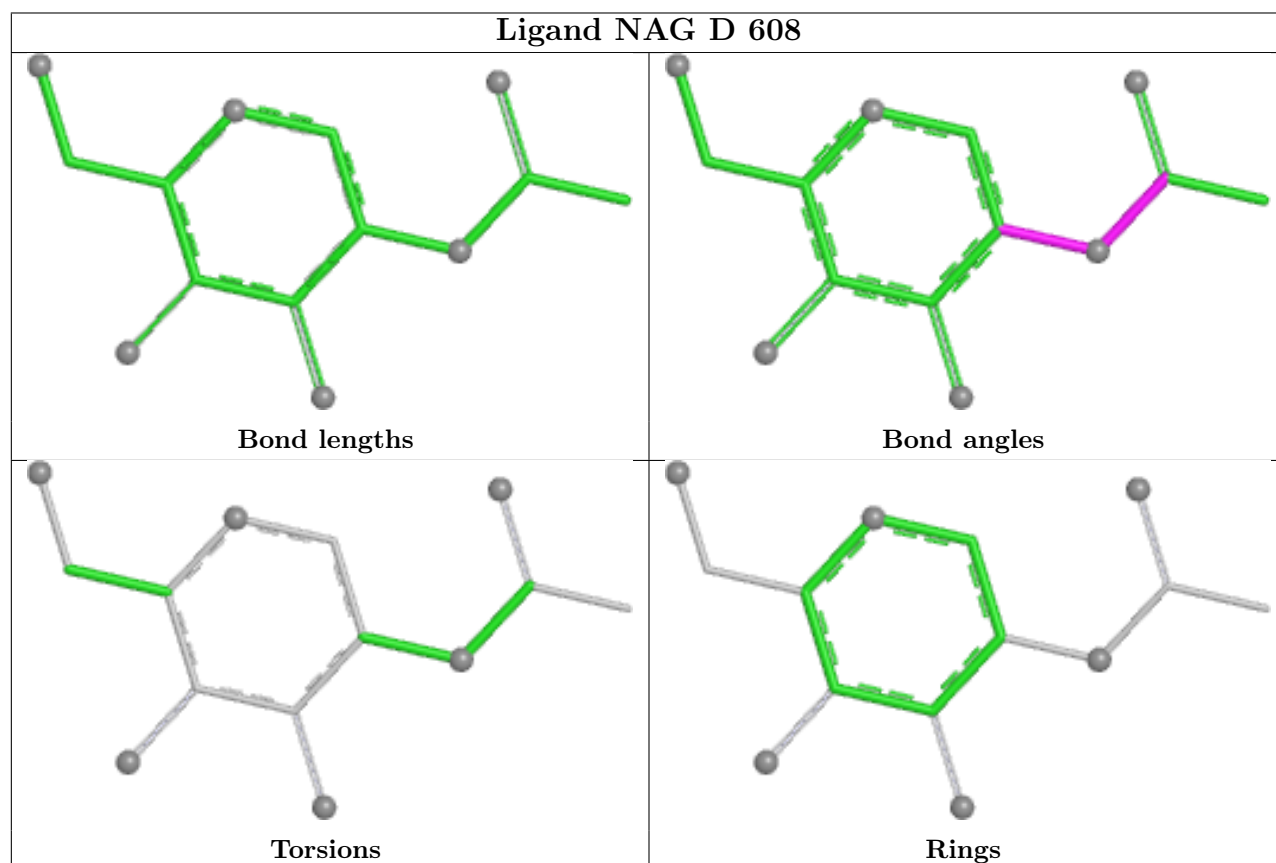
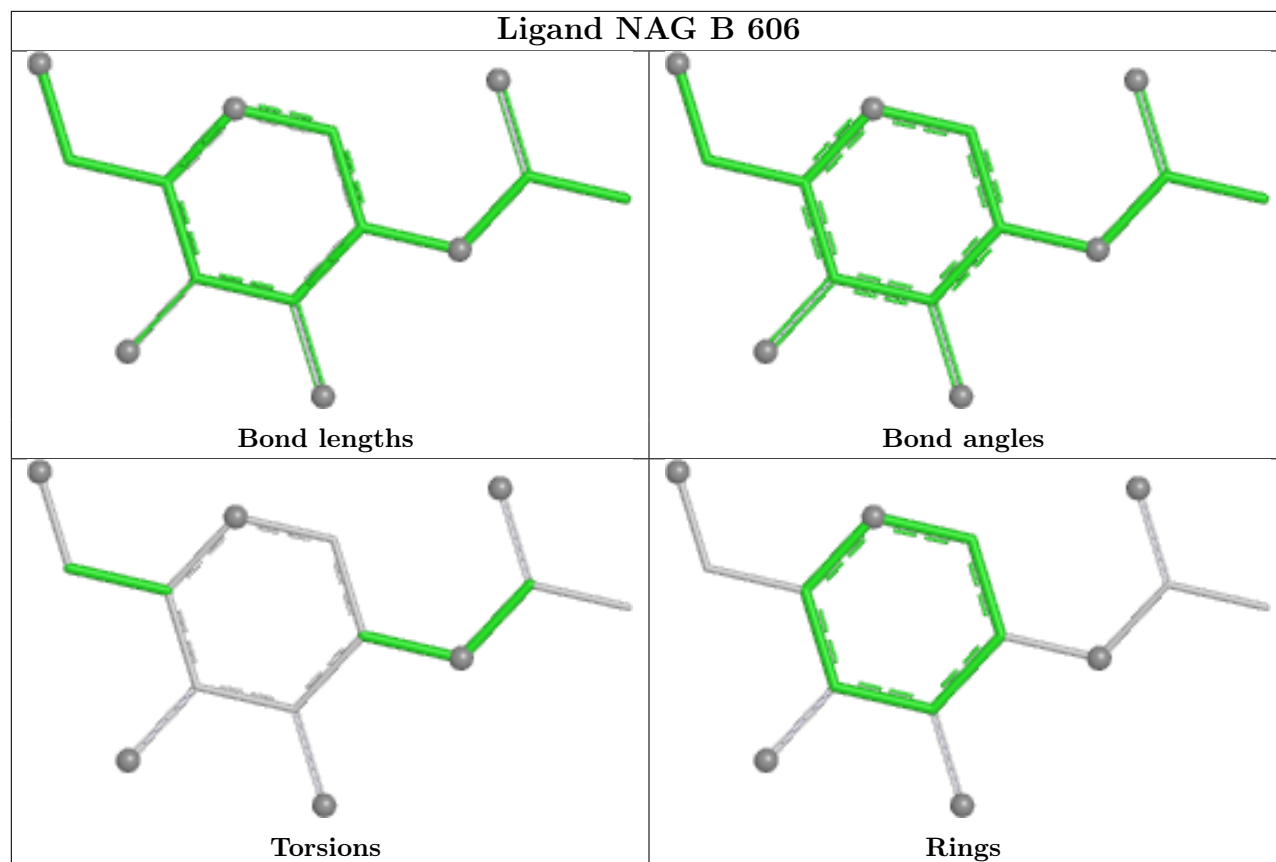


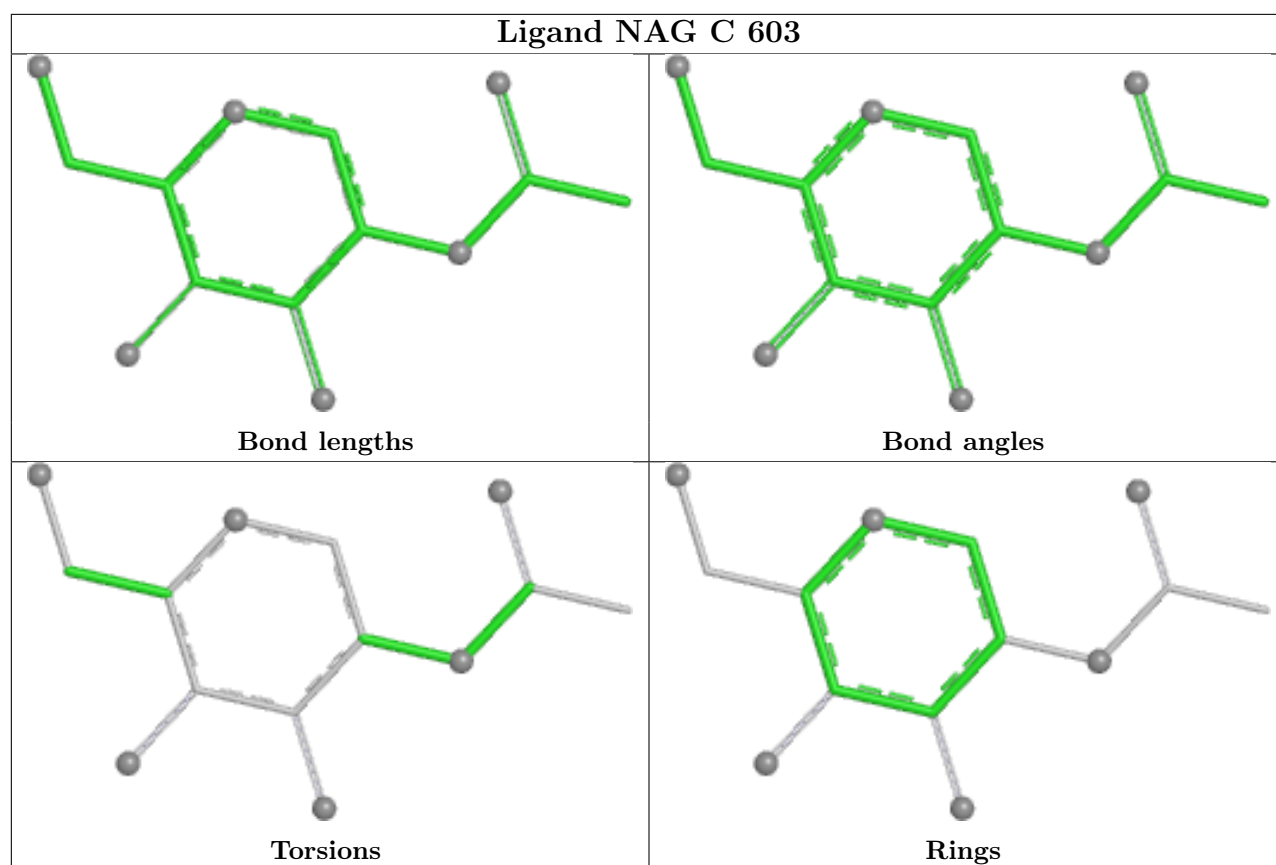


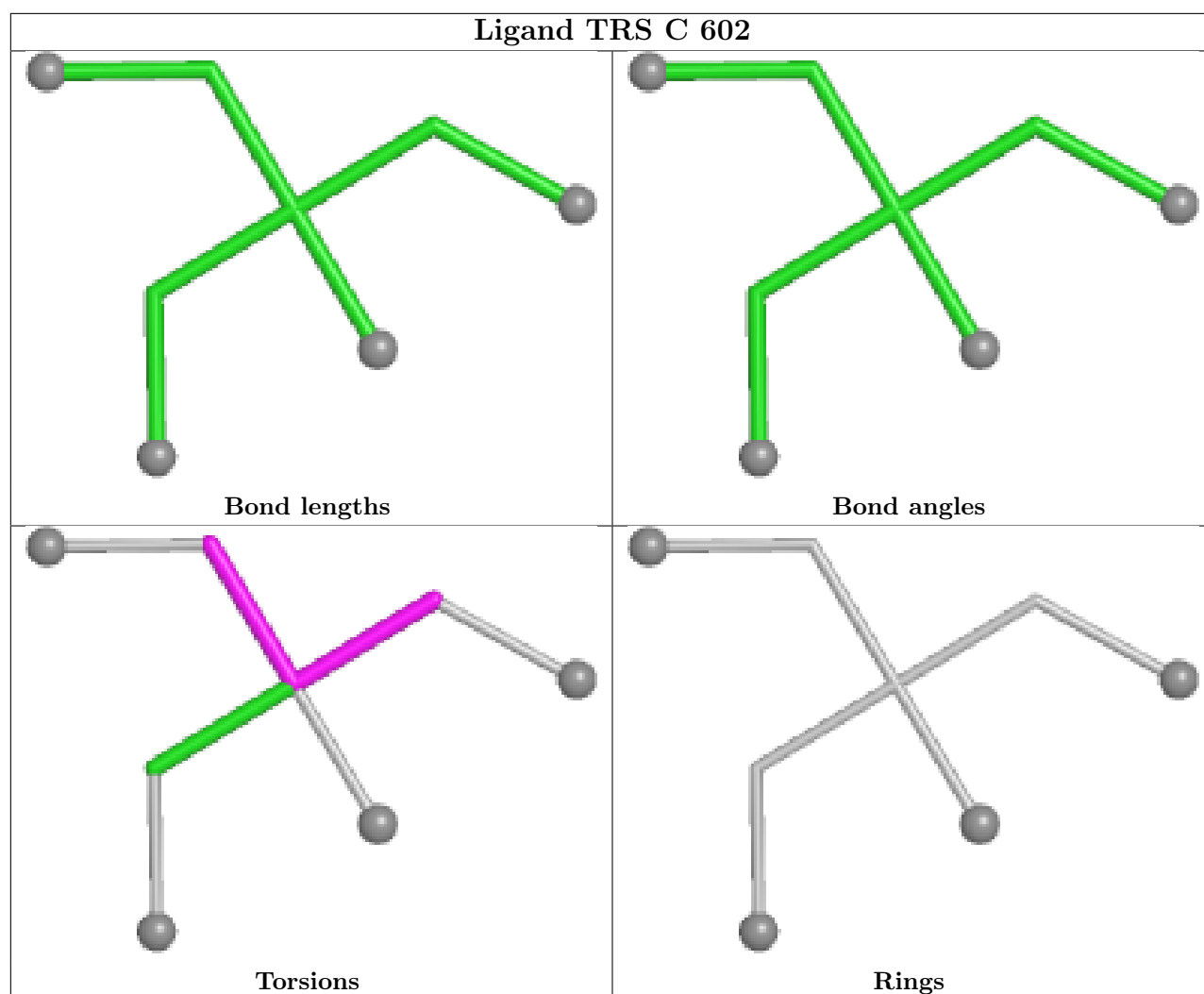












5.7 Other polymers [i](#)

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, 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	535/576 (92%)	-0.08	20 (3%)	45 44	20, 31, 62, 98	1 (0%)
1	B	536/576 (93%)	0.14	21 (3%)	43 42	24, 35, 62, 96	0
1	C	536/576 (93%)	0.01	24 (4%)	38 37	24, 32, 62, 94	0
1	D	535/576 (92%)	0.34	25 (4%)	36 35	24, 37, 68, 98	0
All	All	2142/2304 (92%)	0.10	90 (4%)	40 39	20, 34, 64, 98	1 (0%)

All (90) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	408	ILE	6.9
1	C	336	SER	6.5
1	A	408	ILE	6.4
1	D	336	SER	6.3
1	B	336	SER	6.3
1	C	340	ALA	5.6
1	D	340	ALA	5.5
1	B	408	ILE	5.4
1	A	336	SER	4.8
1	A	39	THR	4.8
1	D	39	THR	4.7
1	B	409	ALA	4.6
1	B	38	THR	4.5
1	D	409	ALA	4.5
1	A	340	ALA	4.4
1	D	41	THR	4.3
1	B	39	THR	4.3
1	A	40	THR	4.1
1	A	41	THR	3.9
1	C	38	THR	3.9
1	B	340	ALA	3.7

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Mol	Chain	Res	Type	RSRZ
1	C	40	THR	3.7
1	D	40	THR	3.7
1	A	335	GLY	3.6
1	D	335	GLY	3.6
1	B	40	THR	3.5
1	B	41	THR	3.5
1	A	409	ALA	3.5
1	C	39	THR	3.5
1	C	156	ASN	3.4
1	C	576	SER	3.4
1	B	576	SER	3.3
1	C	409	ALA	3.3
1	B	335	GLY	3.2
1	D	155	PRO	3.2
1	C	575	GLN	3.2
1	D	407	SER	3.2
1	C	41	THR	3.2
1	D	172	SER	3.1
1	C	407	SER	2.9
1	A	273	GLU	2.9
1	C	197	HIS	2.9
1	D	197	HIS	2.9
1	D	42	THR	2.8
1	D	242	ILE	2.8
1	A	548	ARG	2.8
1	C	50	PRO	2.8
1	D	205	ASP	2.7
1	C	408	ILE	2.7
1	C	173	ASP	2.7
1	B	147	ARG	2.7
1	B	277	ARG	2.7
1	D	219	GLY	2.6
1	A	42	THR	2.6
1	A	89	ASP	2.6
1	C	406	ASP	2.6
1	C	275	GLY	2.6
1	A	437	SER	2.6
1	B	42	THR	2.6
1	C	546	GLY	2.5
1	D	277	ARG	2.5
1	D	437	SER	2.5
1	D	576	SER	2.5

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Mol	Chain	Res	Type	RSRZ
1	C	335	GLY	2.5
1	A	173	ASP	2.4
1	D	274	LYS	2.4
1	B	406	ASP	2.4
1	A	274	LYS	2.4
1	A	565	SER	2.4
1	C	274	LYS	2.3
1	B	197	HIS	2.3
1	D	548	ARG	2.3
1	B	89	ASP	2.3
1	A	497	SER	2.2
1	A	133[A]	ARG	2.2
1	D	156	ASN	2.2
1	B	407	SER	2.2
1	A	172	SER	2.2
1	B	575	GLN	2.1
1	C	438	GLY	2.1
1	D	220	HIS	2.1
1	C	277	ARG	2.1
1	B	273	GLU	2.1
1	B	321	ALA	2.1
1	D	273	GLU	2.0
1	C	147	ARG	2.0
1	D	147	ARG	2.0
1	B	50	PRO	2.0
1	A	439	ASP	2.0
1	C	42	THR	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](#)

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
2	MAN	I	8	11/12	0.61	0.18	40,41,43,43	11

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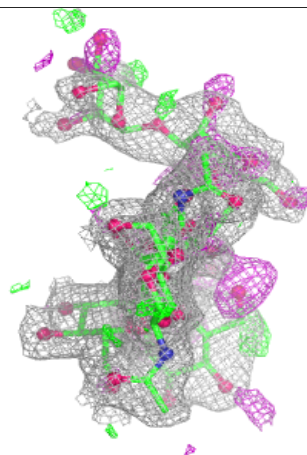
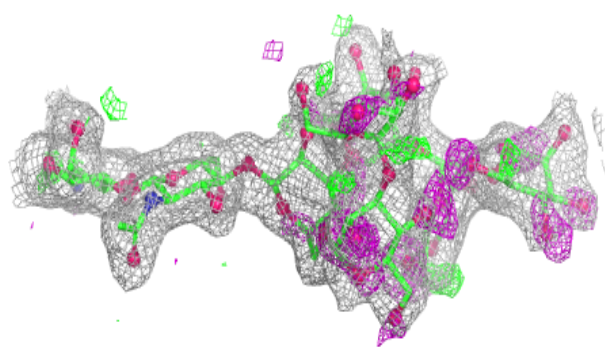
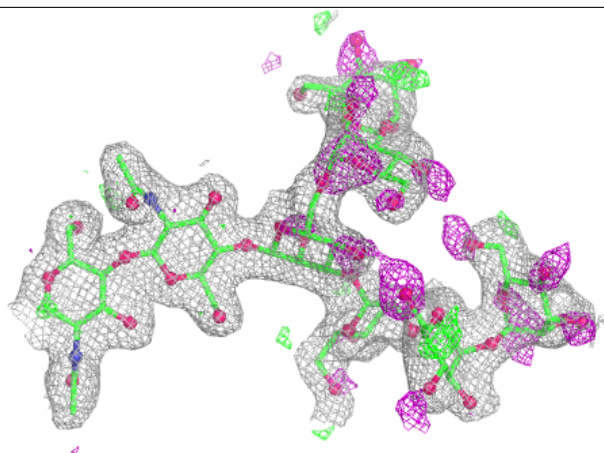
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	MAN	G	8	11/12	0.68	0.17	38,39,41,42	11
3	NAG	F	2	14/15	0.71	0.16	37,39,42,43	14
2	MAN	G	7	11/12	0.72	0.14	39,40,41,42	0
2	MAN	H	8	11/12	0.72	0.15	38,39,40,40	11
2	MAN	I	7	11/12	0.73	0.15	38,40,42,42	0
2	MAN	E	8	11/12	0.75	0.14	38,39,40,42	11
2	MAN	E	6	11/12	0.77	0.14	35,37,40,41	0
2	MAN	E	7	11/12	0.78	0.14	37,39,40,41	0
2	MAN	H	7	11/12	0.78	0.14	38,39,39,40	0
2	MAN	E	5	11/12	0.82	0.11	34,36,39,40	0
2	BMA	I	3	11/12	0.82	0.13	33,35,40,40	0
2	MAN	I	6	11/12	0.82	0.13	33,35,38,40	0
2	MAN	H	5	11/12	0.83	0.11	35,36,39,40	0
2	MAN	H	6	11/12	0.83	0.12	34,36,39,39	0
2	MAN	H	4	11/12	0.83	0.11	34,36,37,40	0
2	BMA	E	3	11/12	0.85	0.11	33,36,38,38	0
2	BMA	H	3	11/12	0.86	0.10	31,35,38,38	0
2	BMA	G	3	11/12	0.88	0.10	31,35,38,38	0
2	MAN	G	4	11/12	0.88	0.10	32,34,38,40	0
2	MAN	G	6	11/12	0.88	0.10	33,35,39,40	0
2	MAN	I	4	11/12	0.88	0.10	33,34,37,39	0
2	MAN	I	5	11/12	0.90	0.10	34,36,39,40	0
2	MAN	G	5	11/12	0.91	0.08	34,35,39,40	0
3	NAG	F	1	14/15	0.92	0.08	33,34,36,38	0
2	MAN	E	4	11/12	0.92	0.09	34,35,37,40	0
2	NAG	I	2	14/15	0.93	0.09	29,31,34,34	0
2	NAG	E	2	14/15	0.94	0.08	30,32,34,34	0
2	NAG	G	2	14/15	0.94	0.08	28,31,33,33	0
2	NAG	H	2	14/15	0.94	0.07	30,32,35,36	0
2	NAG	E	1	14/15	0.96	0.10	27,29,30,32	0
2	NAG	H	1	14/15	0.97	0.08	28,30,32,32	0
2	NAG	G	1	14/15	0.97	0.08	27,30,32,35	0
2	NAG	I	1	14/15	0.97	0.07	27,29,31,31	0

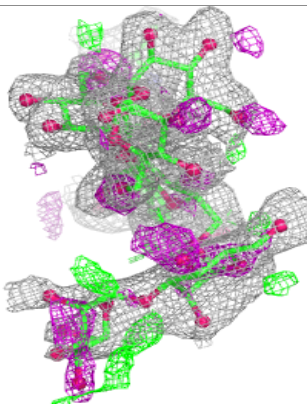
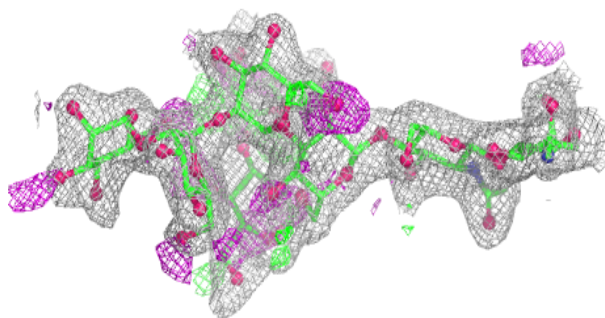
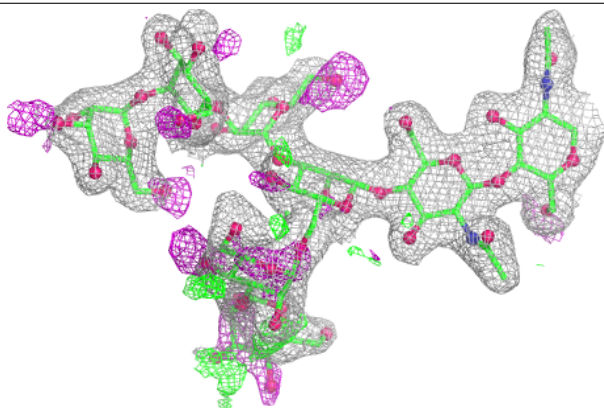
The following is a graphical depiction of the model fit to experimental electron density for oligosaccharide. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around Chain E:

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

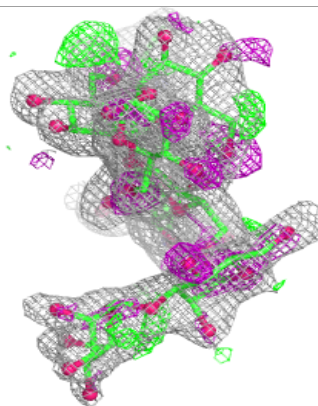
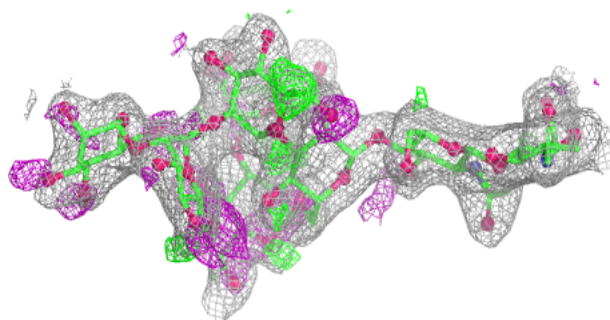
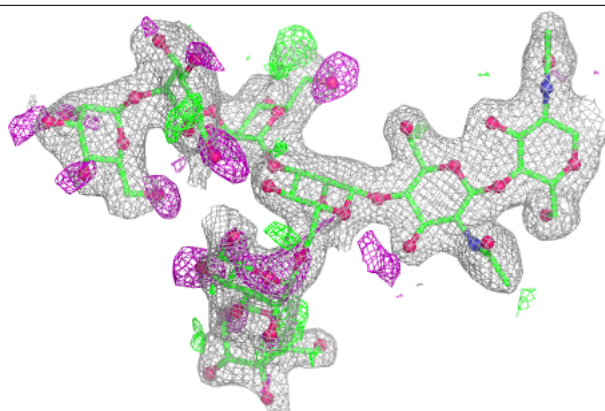
**Electron density around Chain G:**

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

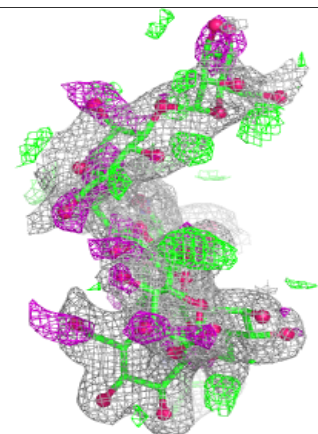
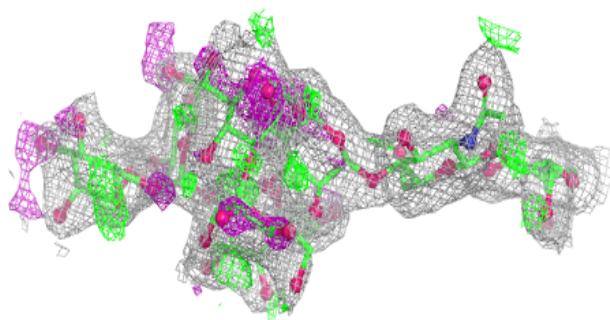
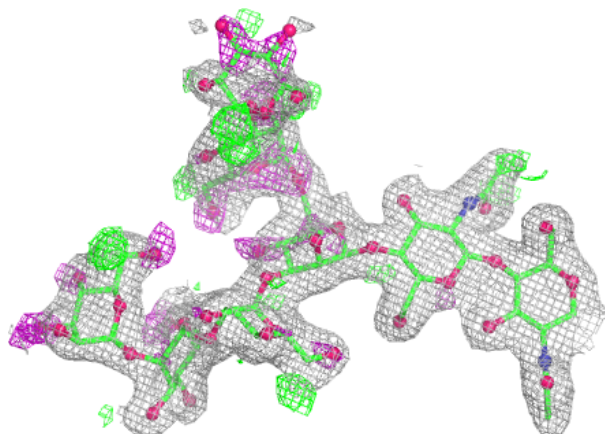


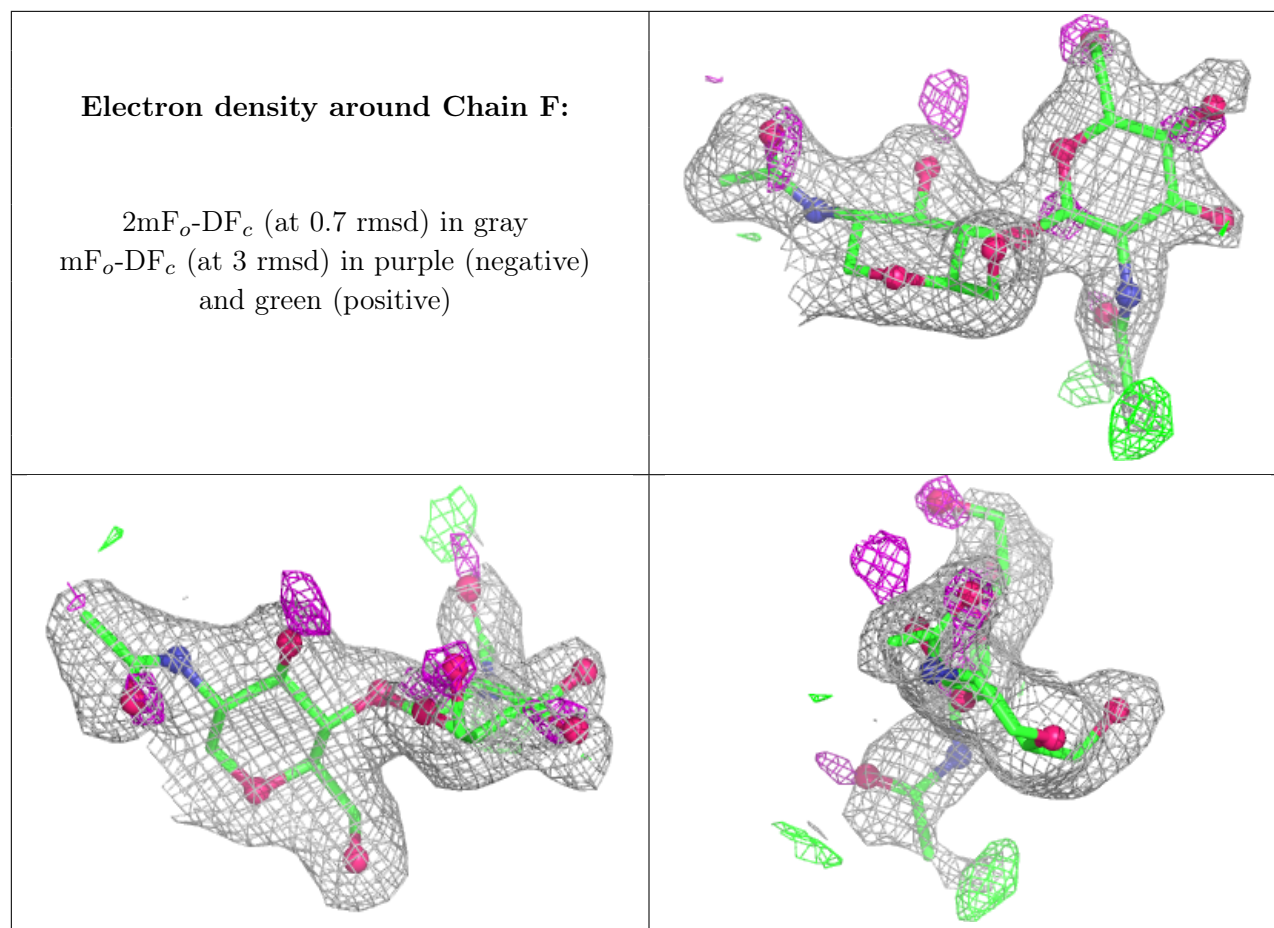
Electron density around Chain H:

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

**Electron density around Chain I:**

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





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($\text{\AA}^2$)	Q<0.9
7	TRS	D	605	8/8	0.59	0.32	24,27,28,28	8
7	TRS	C	602	8/8	0.62	0.35	24,27,28,29	8
7	TRS	A	604	8/8	0.66	0.34	25,27,28,28	8
8	NAG	B	604	14/15	0.70	0.16	35,39,41,42	14
7	TRS	B	608	8/8	0.71	0.30	25,28,29,29	8
8	NAG	A	606	14/15	0.75	0.13	37,39,40,41	0
8	NAG	D	607	14/15	0.75	0.13	38,39,41,42	0
8	NAG	D	609	14/15	0.75	0.13	37,39,41,41	0
8	NAG	C	604	14/15	0.76	0.13	36,39,40,42	0
8	NAG	C	605	14/15	0.79	0.12	34,36,39,40	0
8	NAG	B	605	14/15	0.80	0.11	35,37,40,41	0
8	NAG	B	606	14/15	0.80	0.13	36,39,41,42	0

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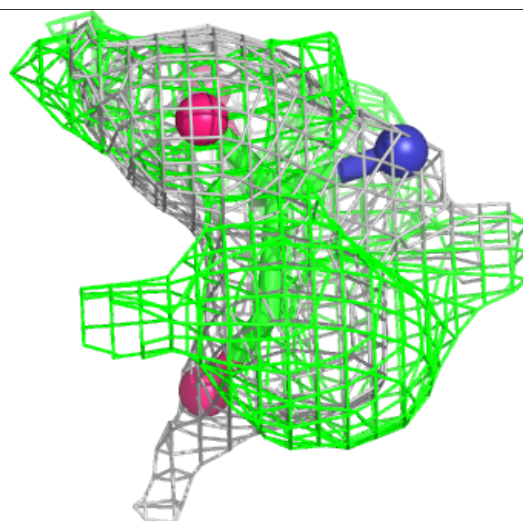
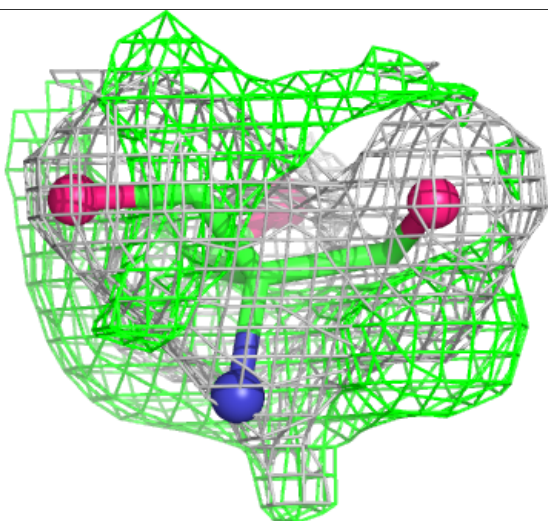
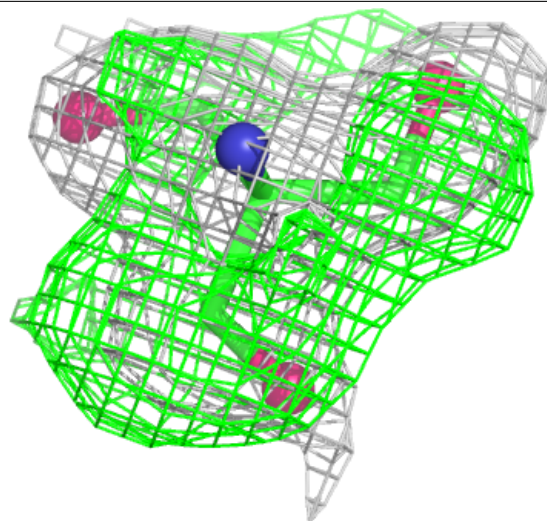
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
9	GOL	B	607	6/6	0.80	0.34	24,26,28,28	6
4	GLC	B	601	11/12	0.81	0.17	47,54,60,63	0
8	NAG	C	606	14/15	0.84	0.13	36,39,40,41	0
8	NAG	D	608	14/15	0.84	0.11	36,37,39,41	0
8	NAG	A	607	14/15	0.85	0.10	34,36,39,40	0
5	FRU	B	602	12/12	0.87	0.11	33,39,46,48	0
8	NAG	A	608	14/15	0.87	0.11	34,37,39,41	0
6	SO4	D	603	5/5	0.88	0.13	25,25,28,29	5
5	FRU	D	602	12/12	0.89	0.12	32,41,52,57	0
6	SO4	A	603	5/5	0.89	0.16	20,20,22,23	5
8	NAG	C	603	14/15	0.90	0.09	32,35,36,36	0
8	NAG	D	606	14/15	0.90	0.09	35,36,39,39	0
4	GLC	C	607	11/12	0.92	0.09	36,47,50,51	0
4	GLC	D	601	11/12	0.92	0.09	39,49,51,53	0
6	SO4	A	609	5/5	0.92	0.13	22,27,27,27	5
4	GLC	A	601	11/12	0.92	0.09	36,47,53,53	0
8	NAG	A	605	14/15	0.92	0.08	32,33,35,37	0
6	SO4	C	601	5/5	0.95	0.12	23,23,24,26	5
5	FRU	A	602	12/12	0.95	0.07	29,36,41,43	0
6	SO4	D	604	5/5	0.96	0.14	20,22,24,24	5
5	FRU	C	608	12/12	0.96	0.07	26,35,46,50	0
6	SO4	B	603	5/5	0.96	0.18	19,22,23,24	5

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.

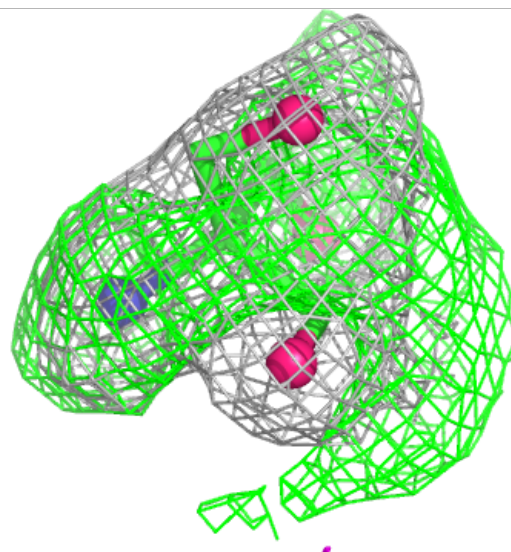
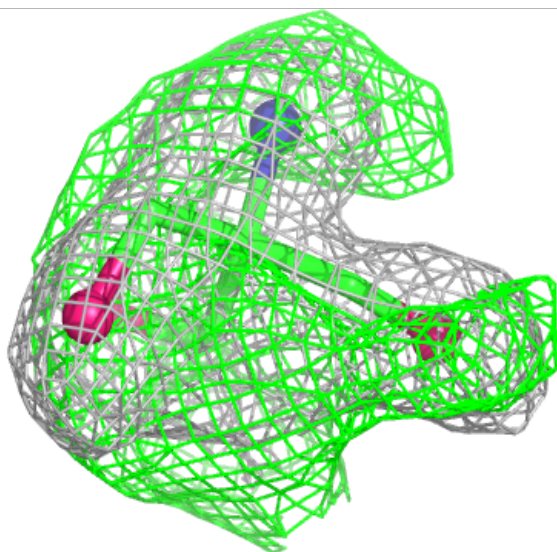
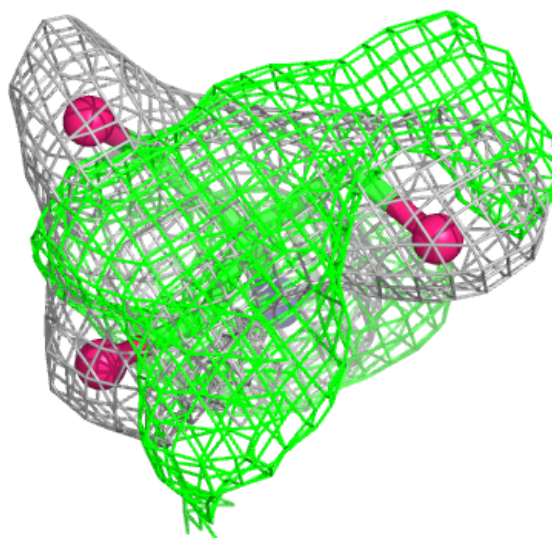
Electron density around TRS D 605:

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



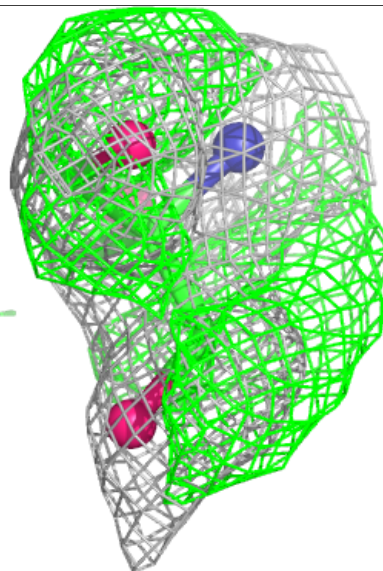
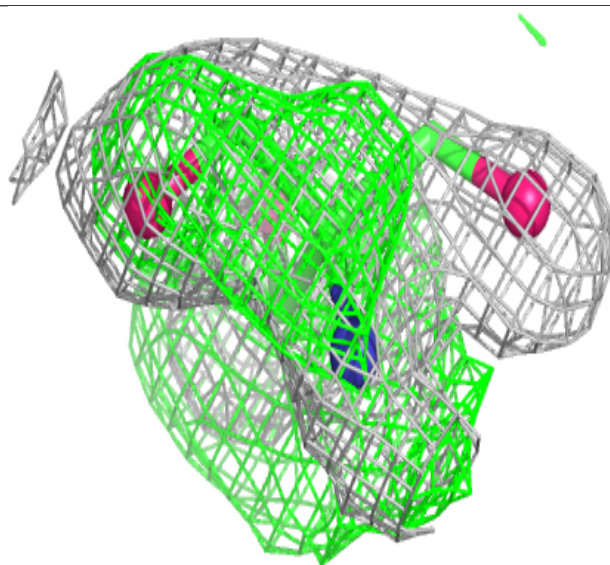
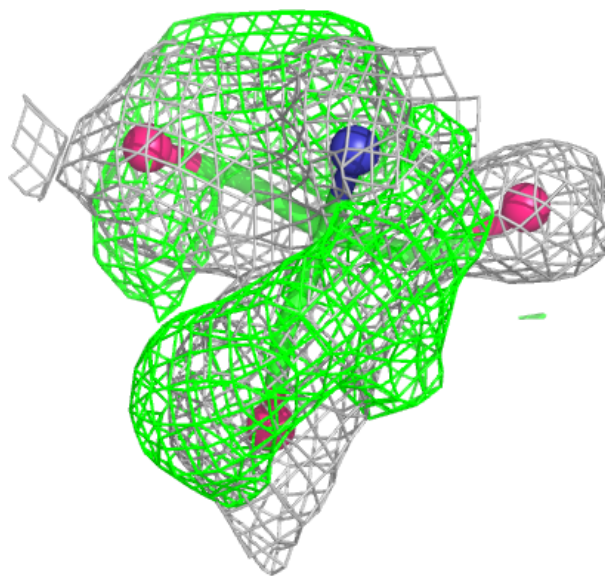
Electron density around TRS C 602:

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



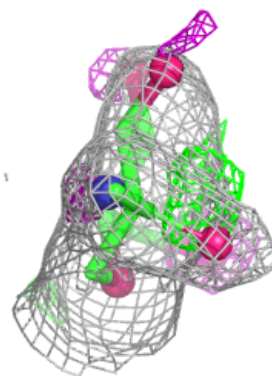
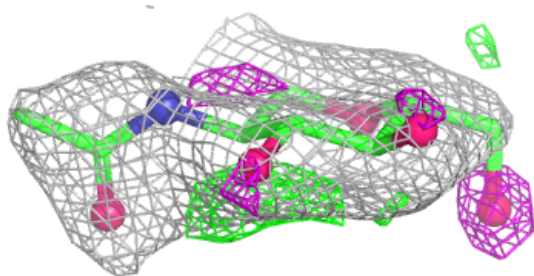
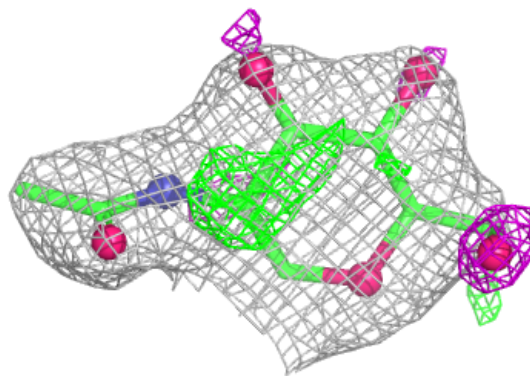
Electron density around TRS A 604:

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



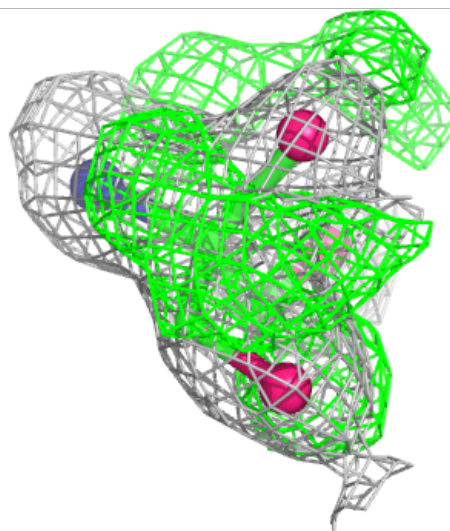
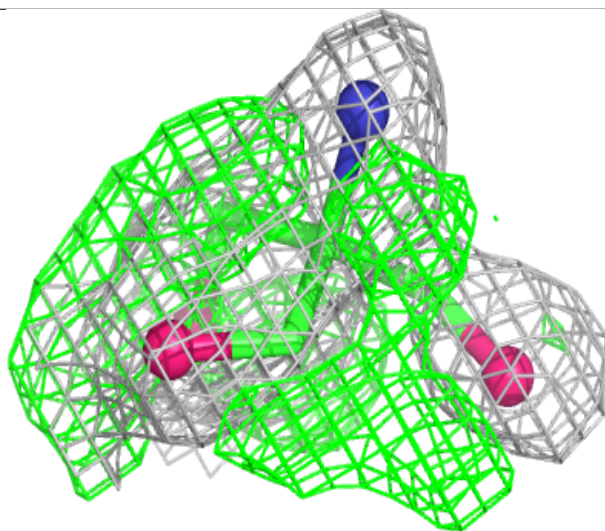
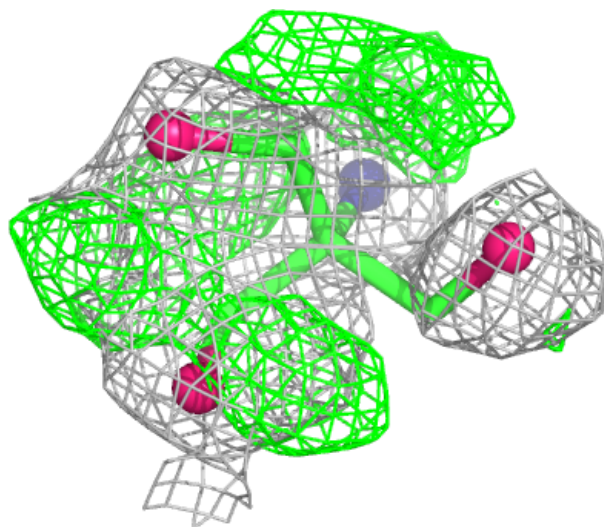
Electron density around NAG B 604:

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



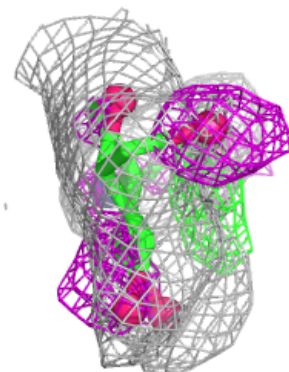
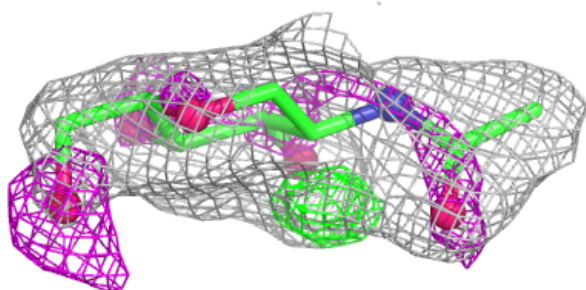
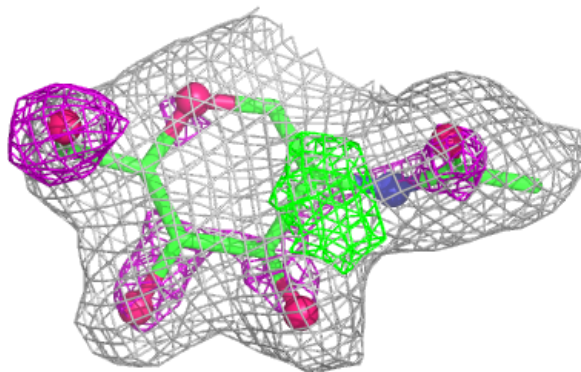
Electron density around TRS B 608:

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

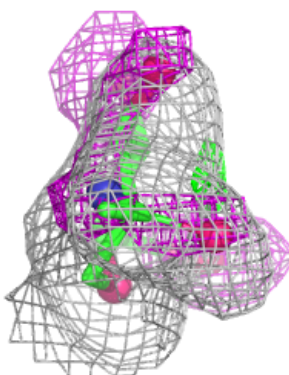
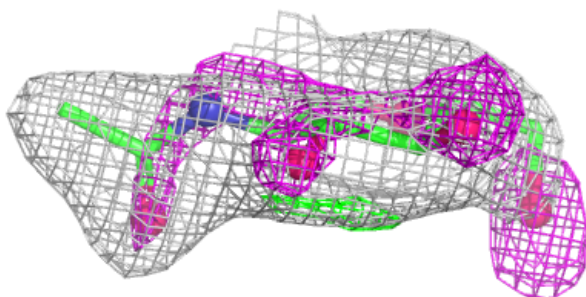
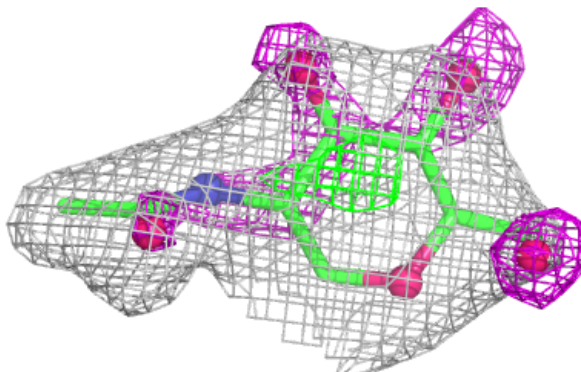


Electron density around NAG A 606:

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

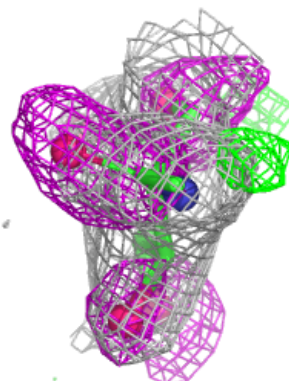
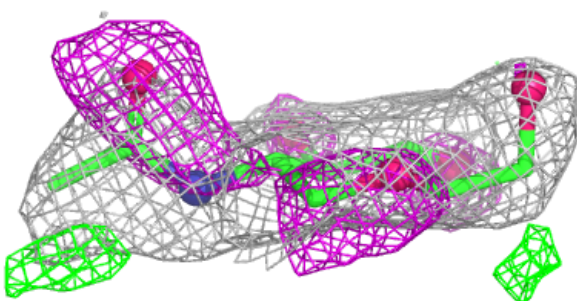
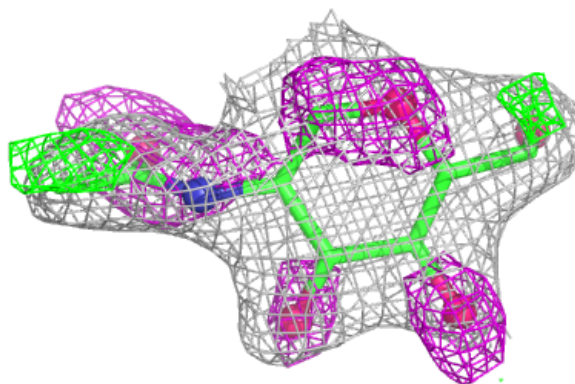
**Electron density around NAG D 607:**

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

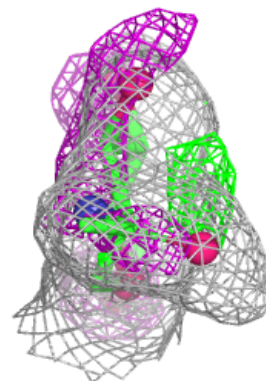
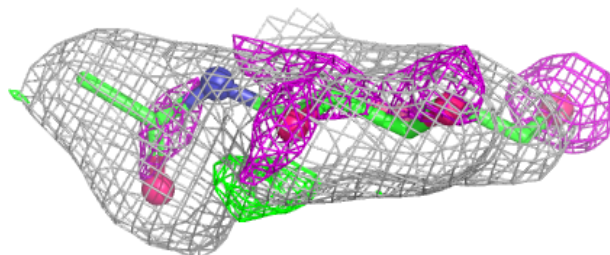
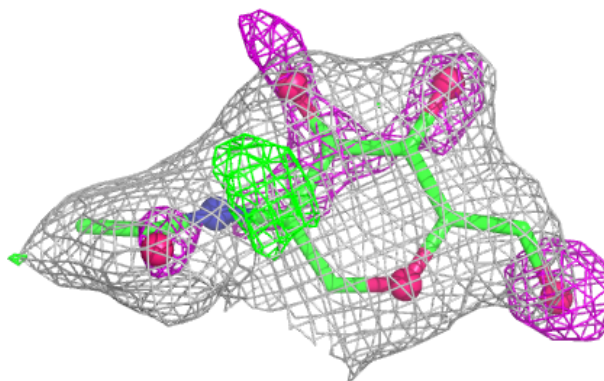


Electron density around NAG D 609:

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

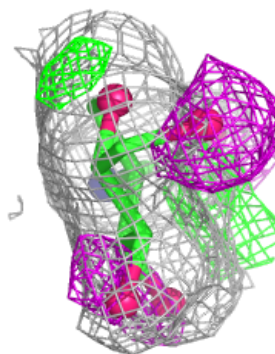
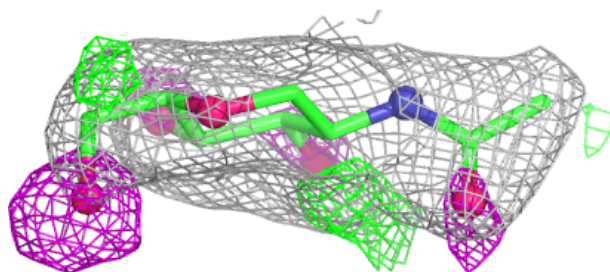
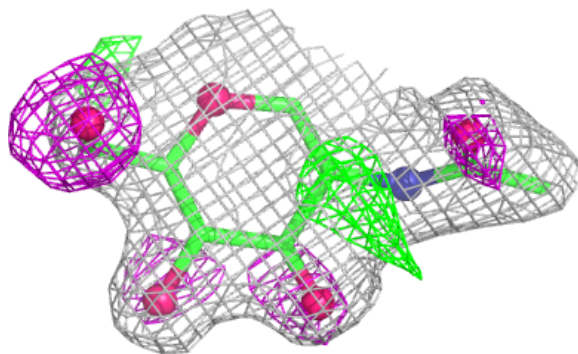
**Electron density around NAG C 604:**

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and green (positive)

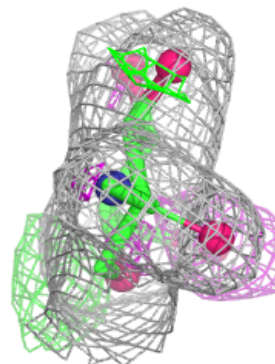
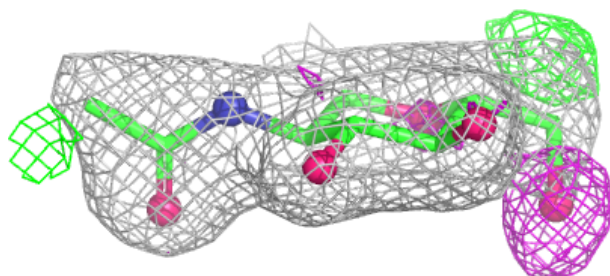
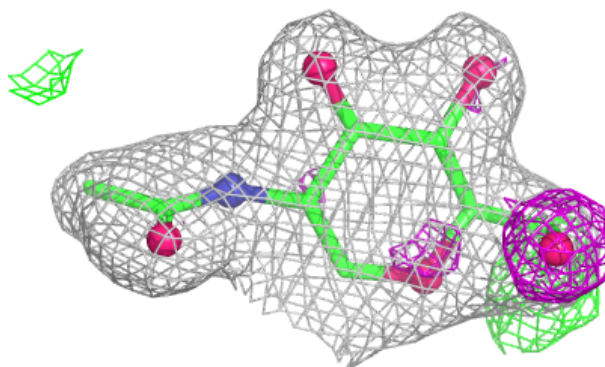


Electron density around NAG C 605:

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

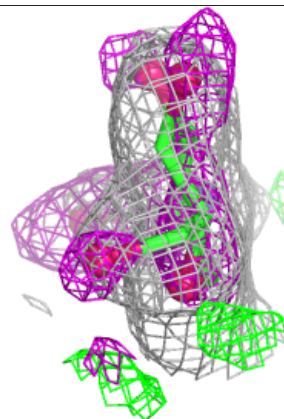
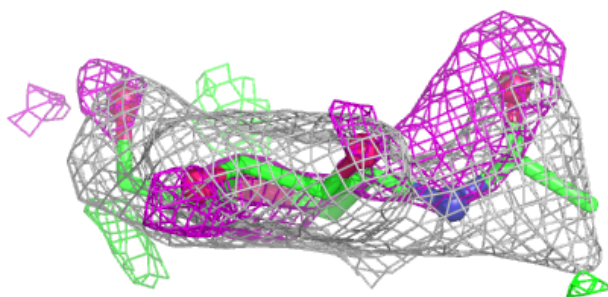
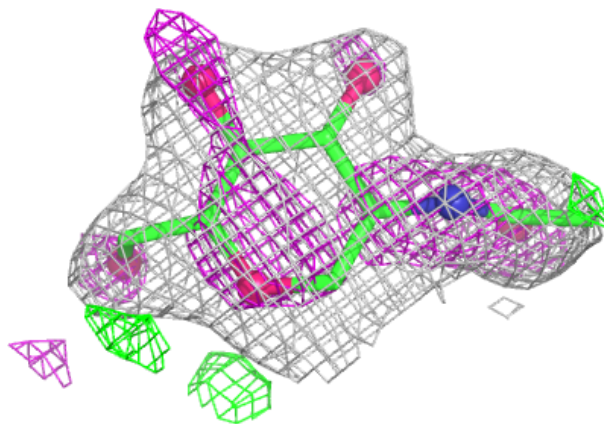
**Electron density around NAG B 605:**

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



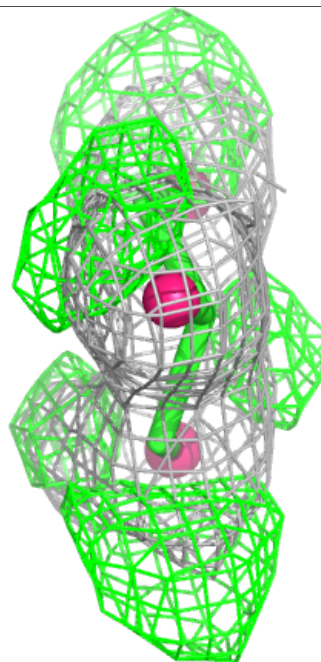
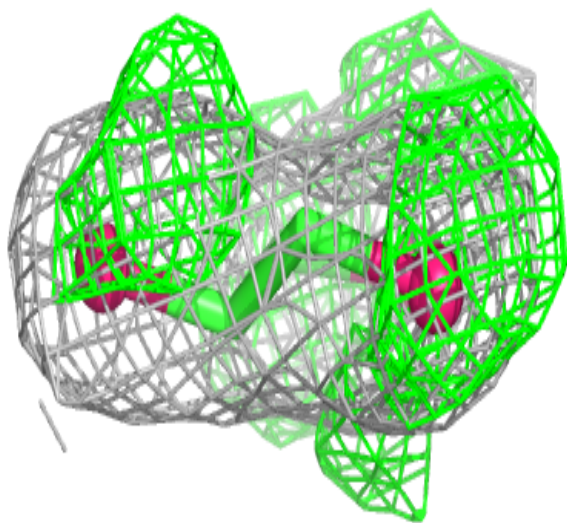
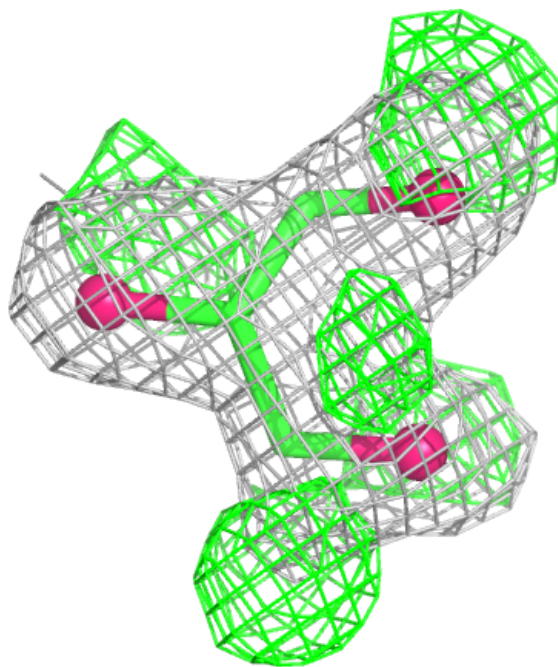
Electron density around NAG B 606:

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



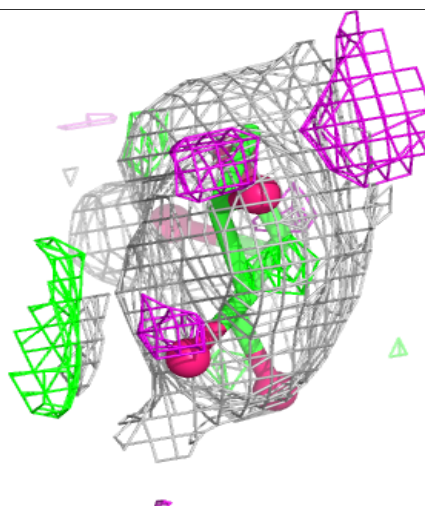
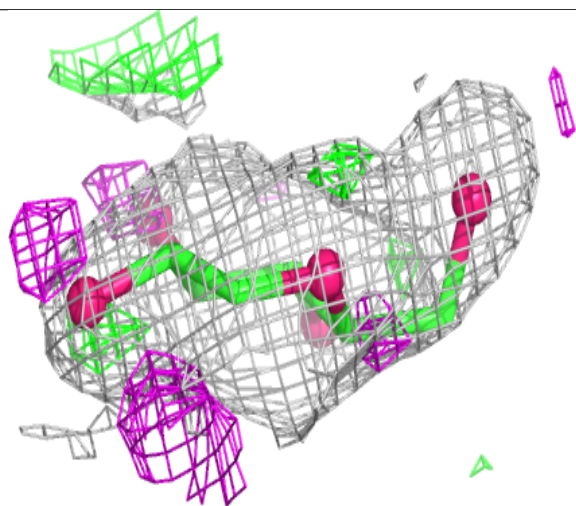
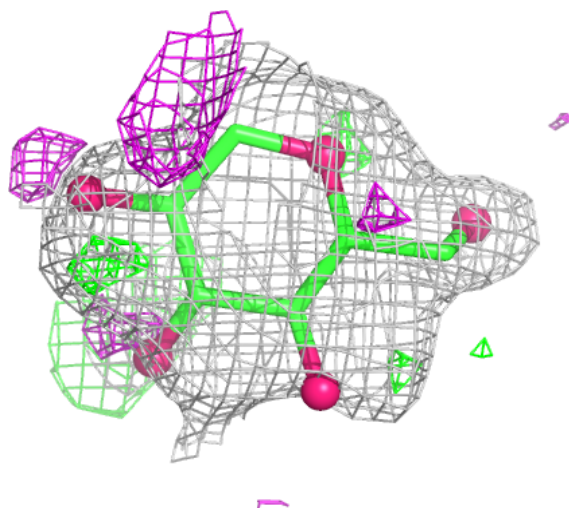
Electron density around GOL B 607:

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



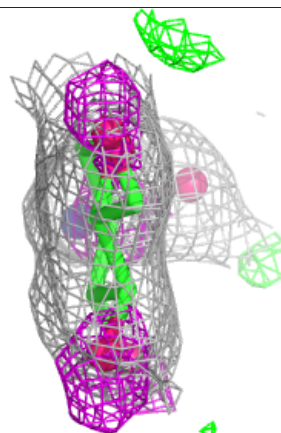
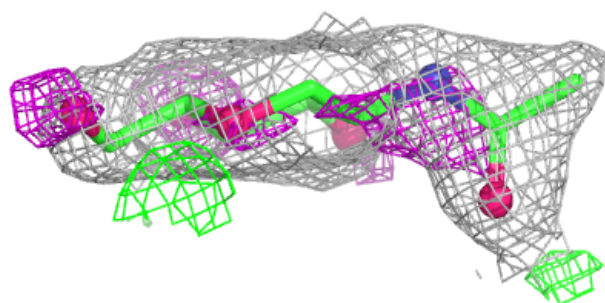
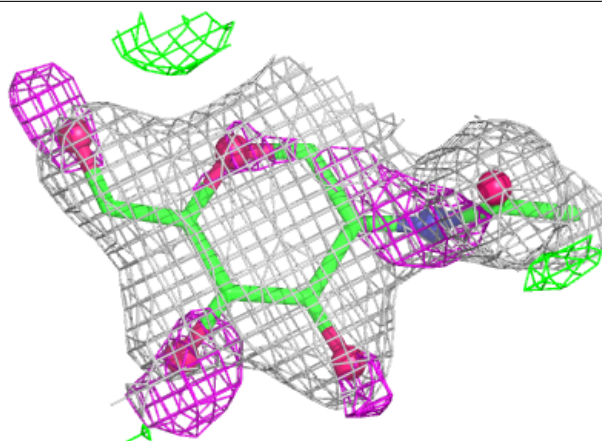
Electron density around GLC B 601:

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

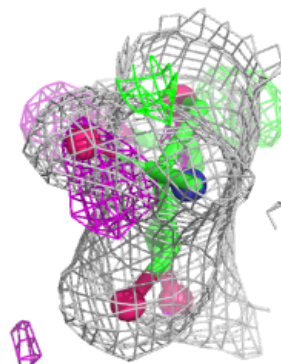
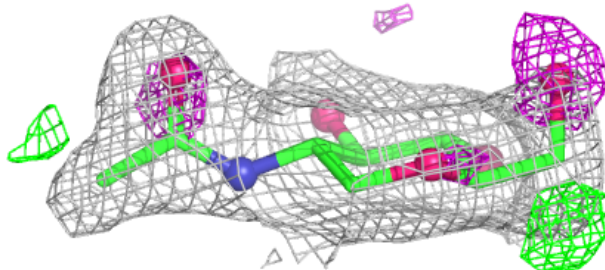
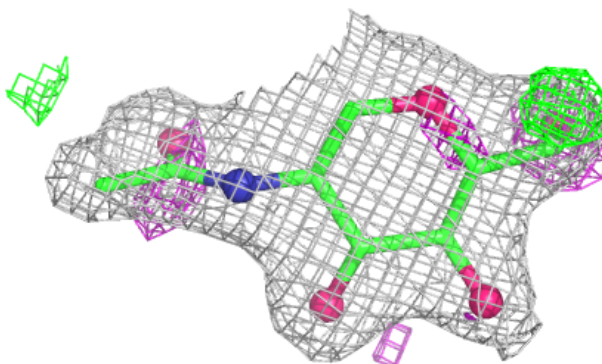


Electron density around NAG C 606:

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

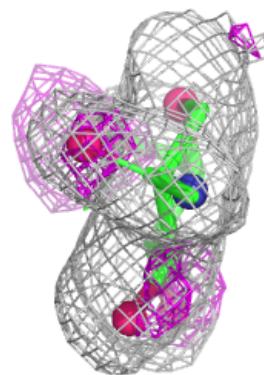
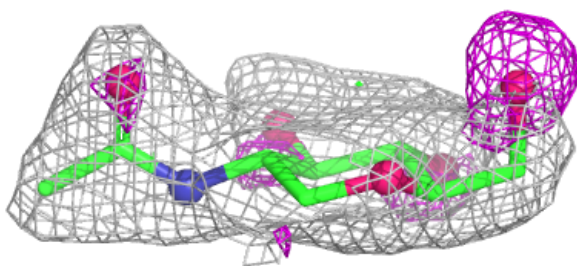
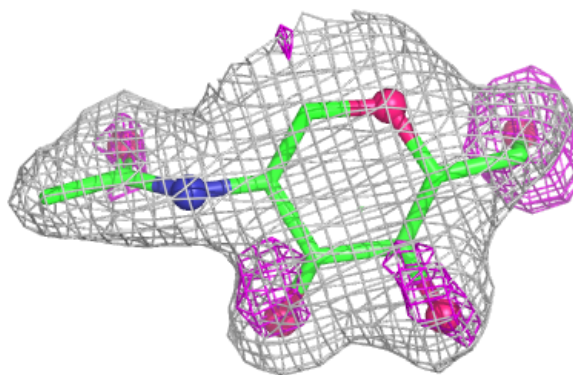
**Electron density around NAG D 608:**

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



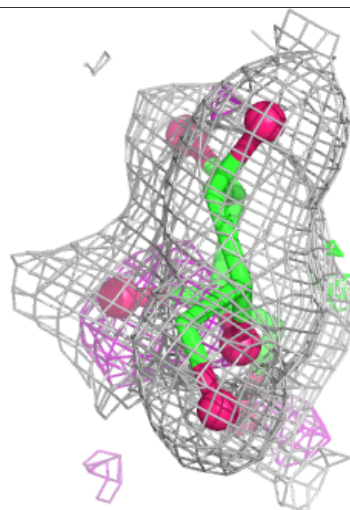
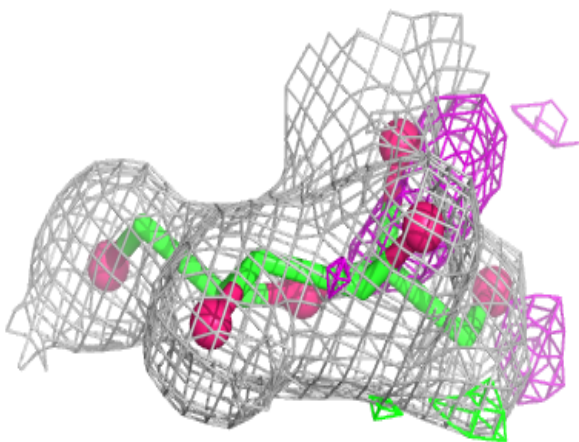
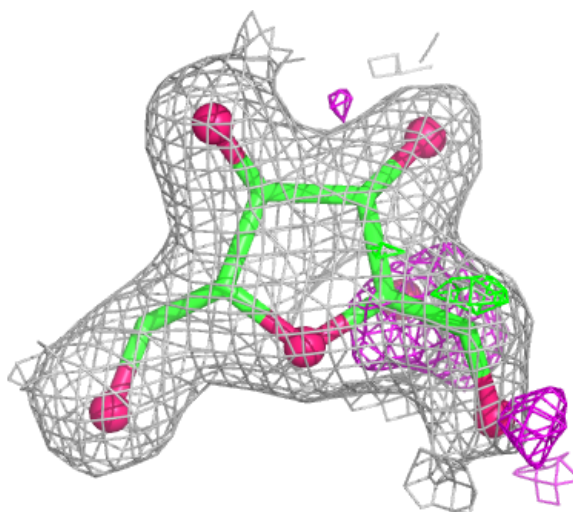
Electron density around NAG A 607:

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



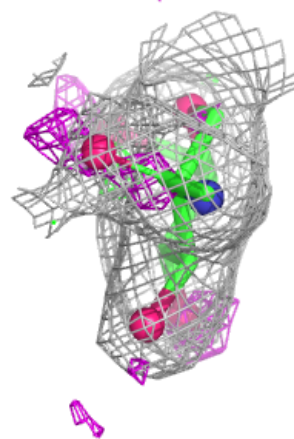
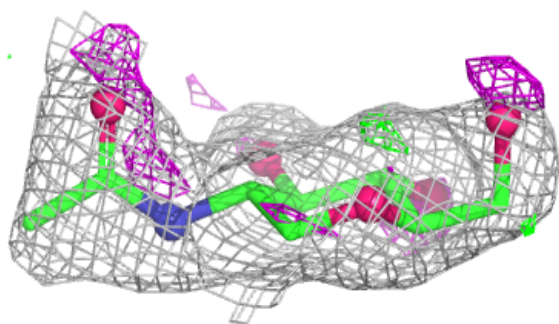
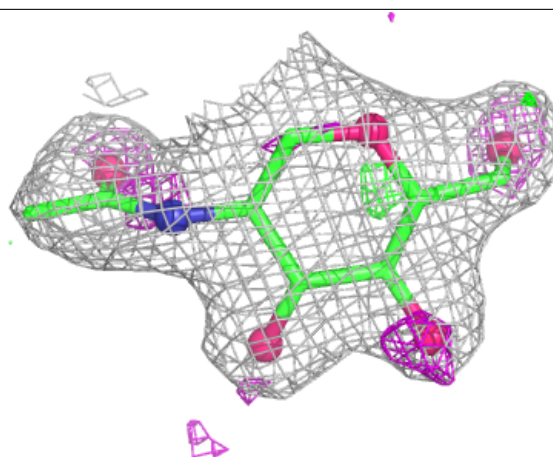
Electron density around FRU B 602:

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



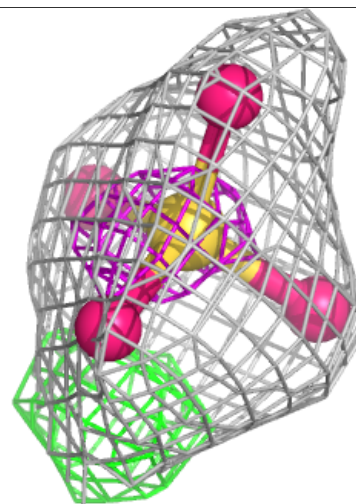
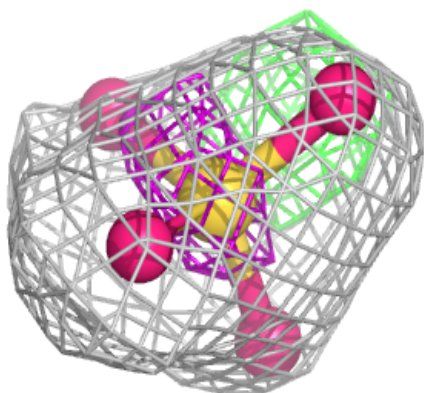
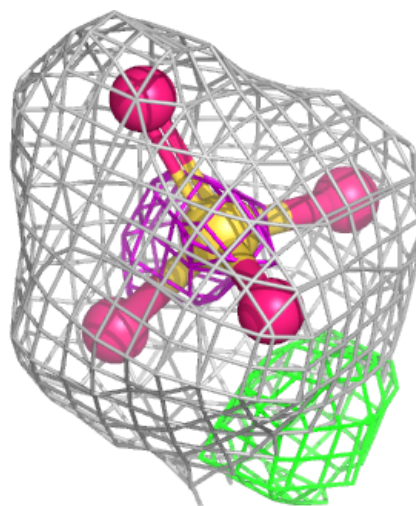
Electron density around NAG A 608:

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



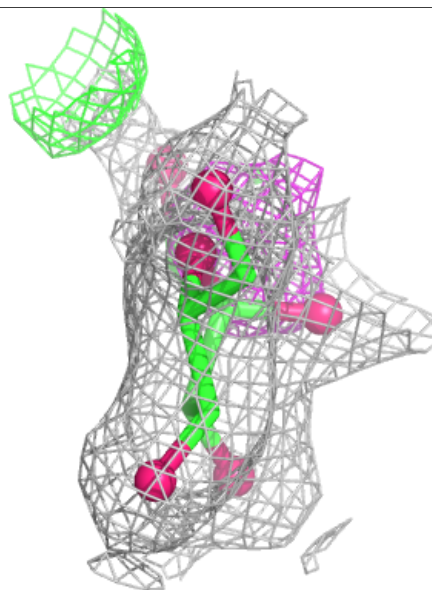
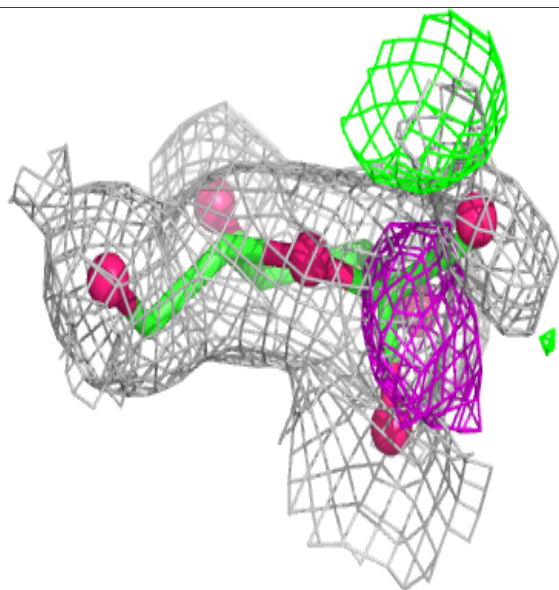
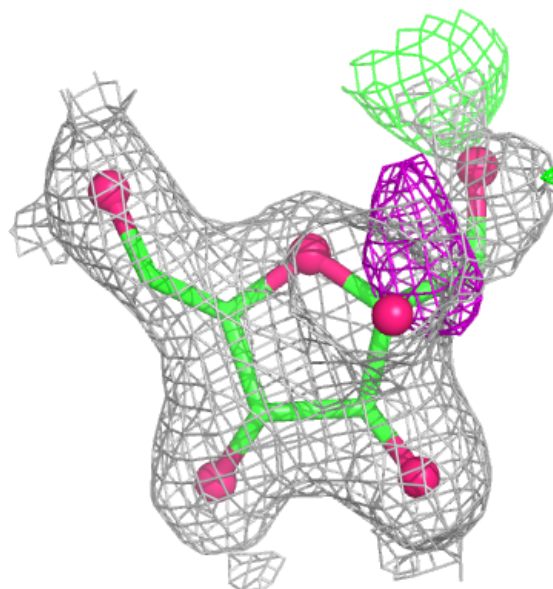
Electron density around SO4 D 603:

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



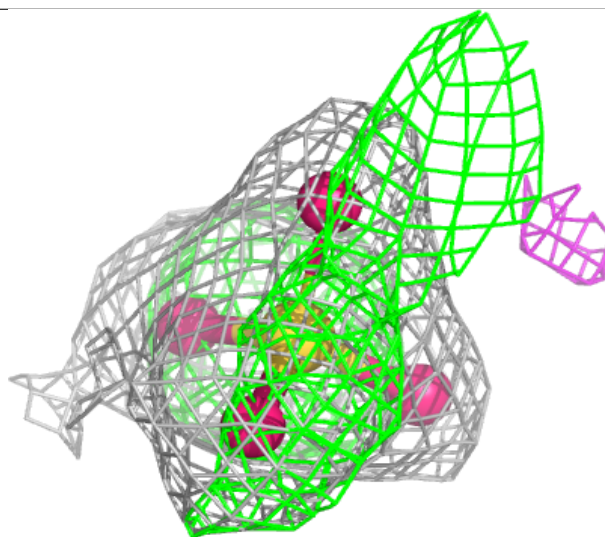
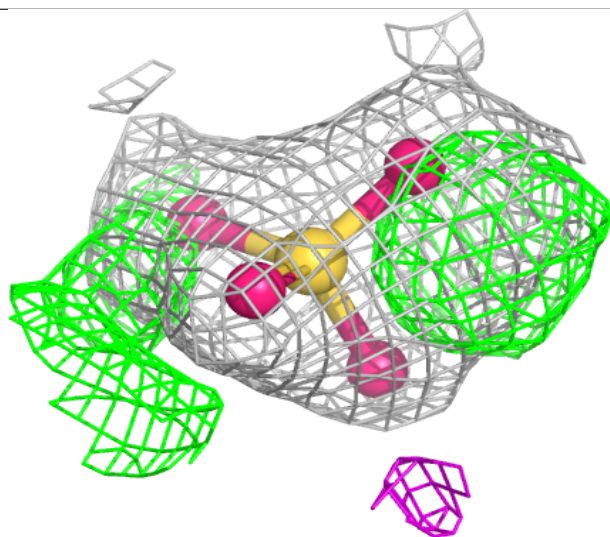
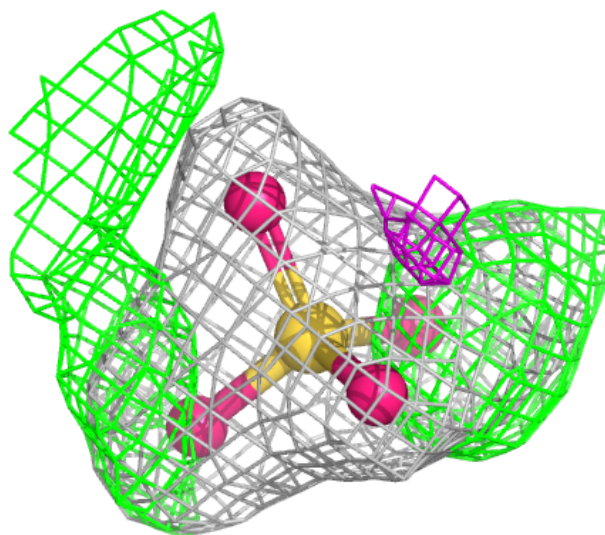
Electron density around FRU D 602:

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



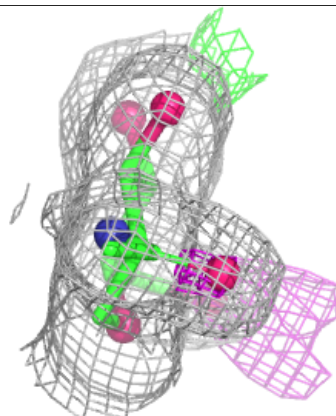
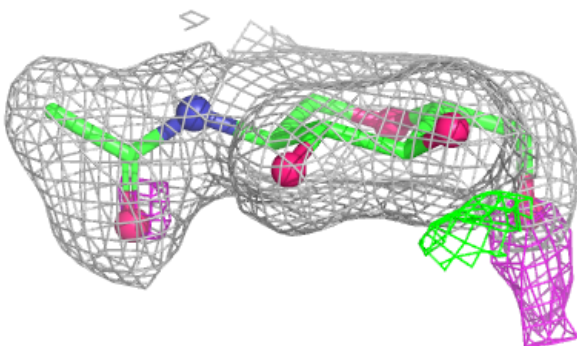
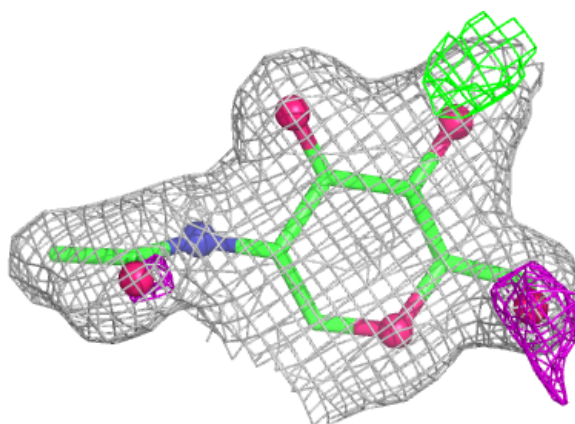
Electron density around SO4 A 603:

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

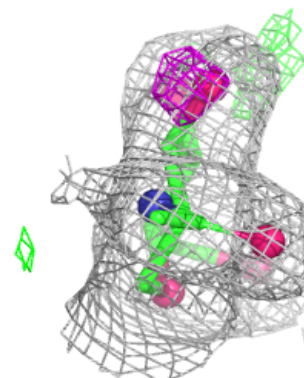
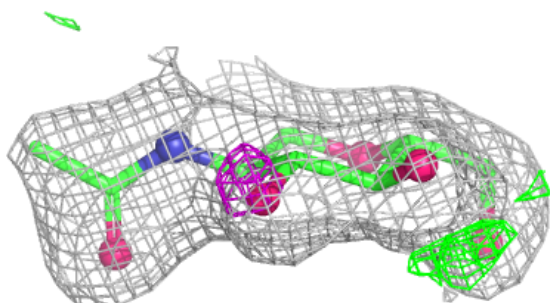
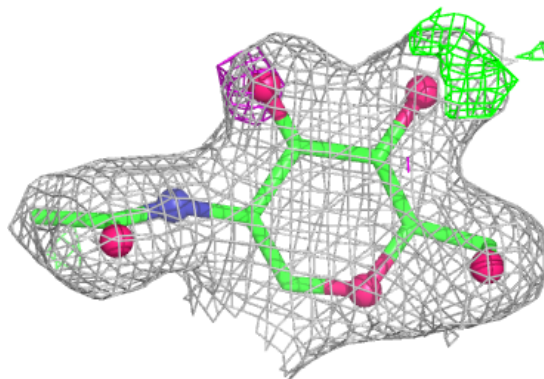


Electron density around NAG C 603:

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

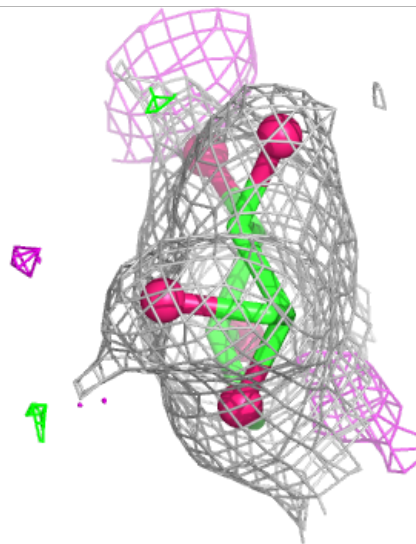
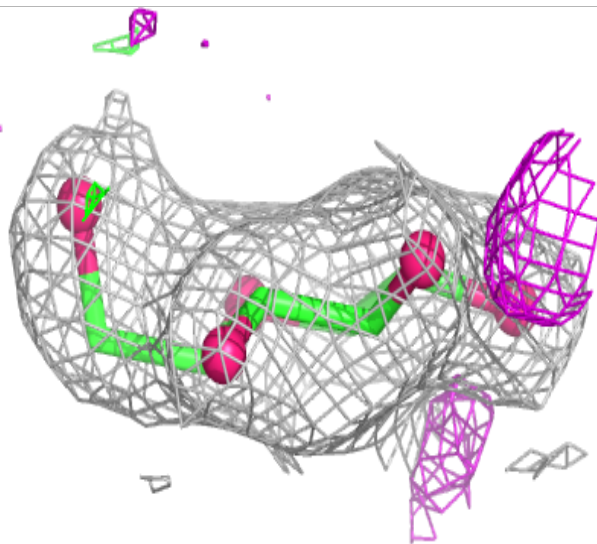
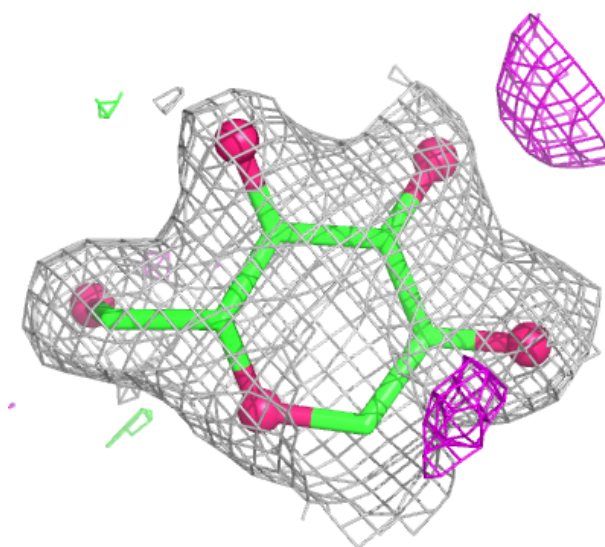
**Electron density around NAG D 606:**

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



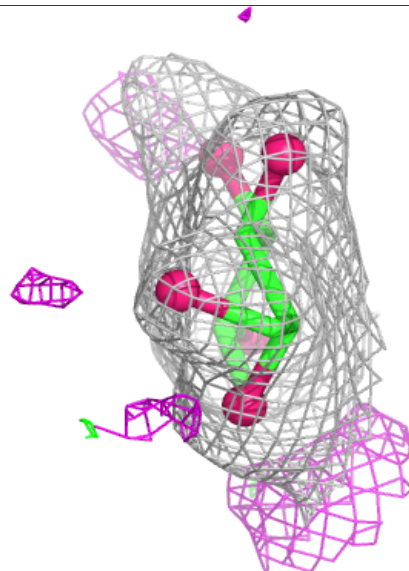
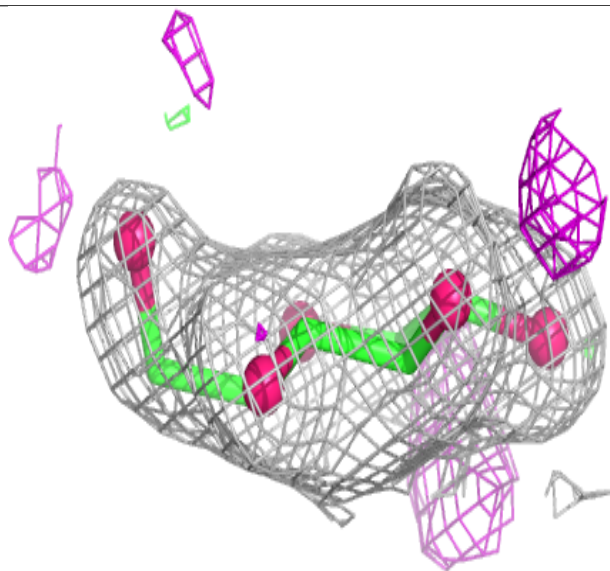
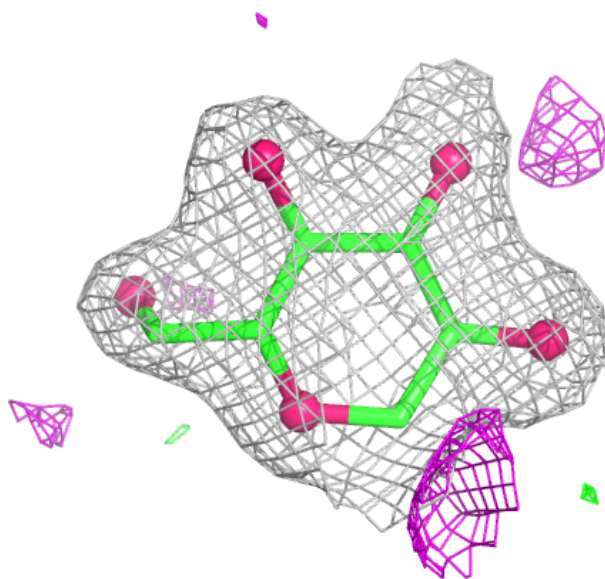
Electron density around GLC C 607:

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



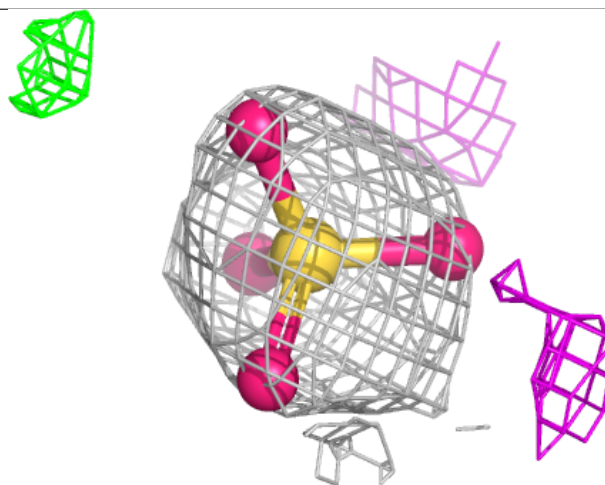
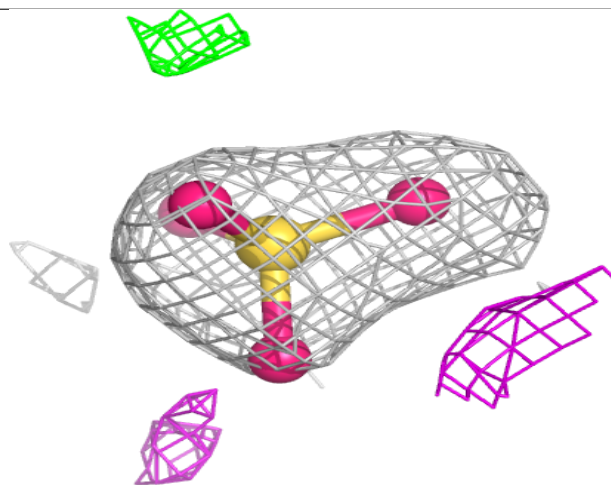
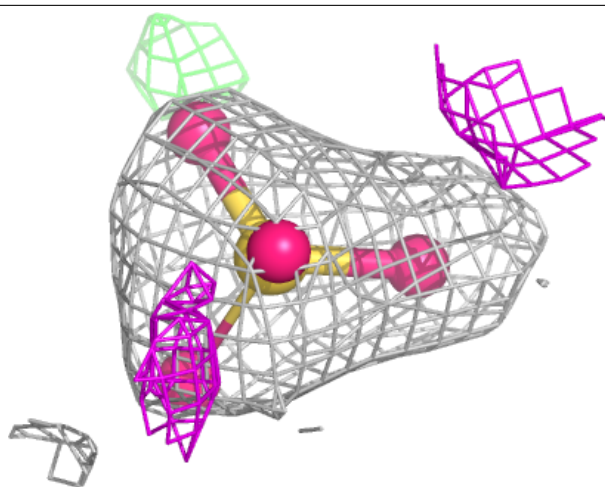
Electron density around GLC D 601:

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



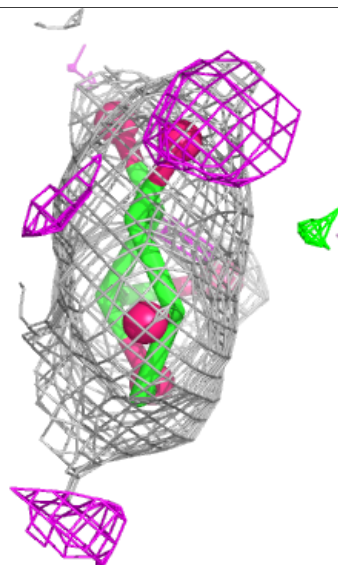
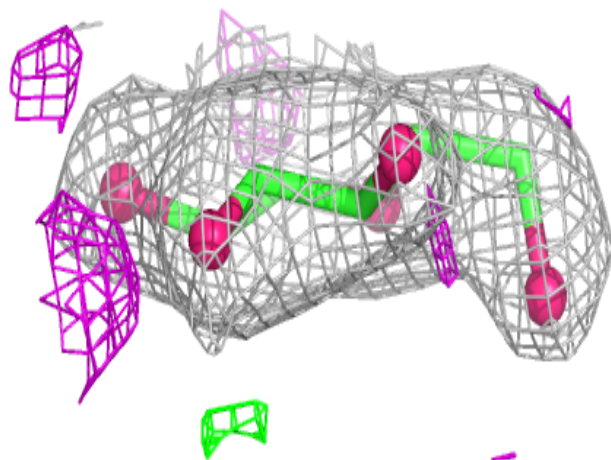
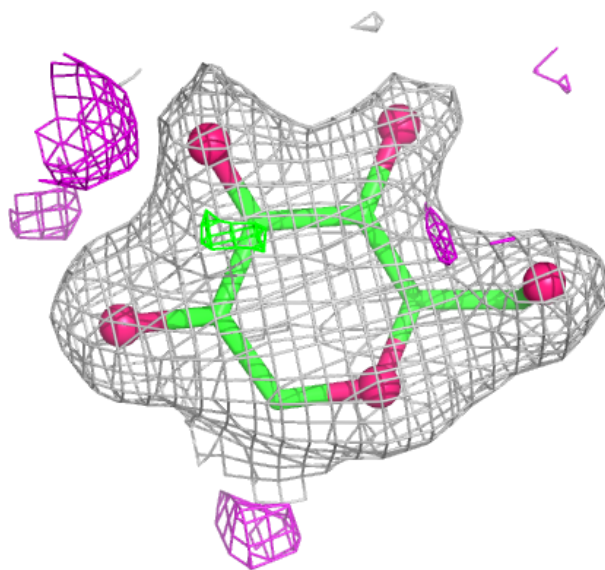
Electron density around SO4 A 609:

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



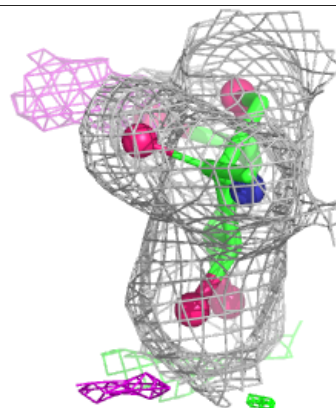
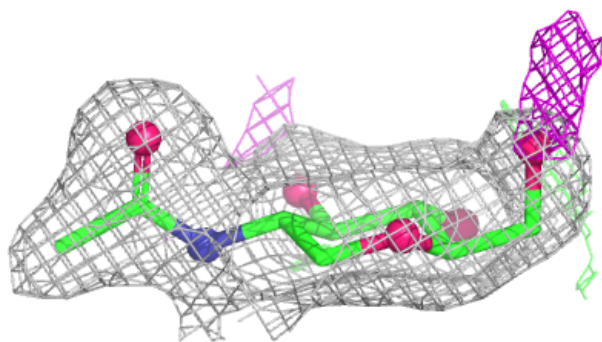
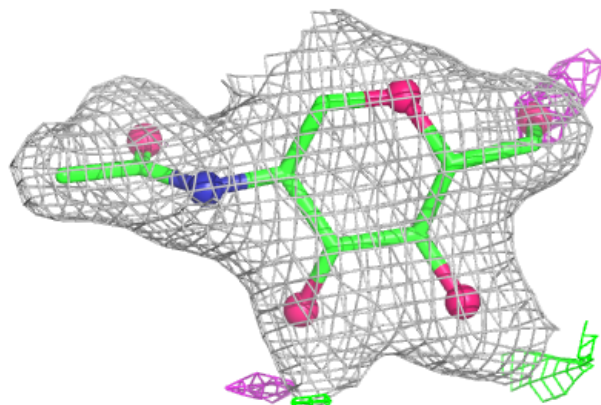
Electron density around GLC A 601:

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



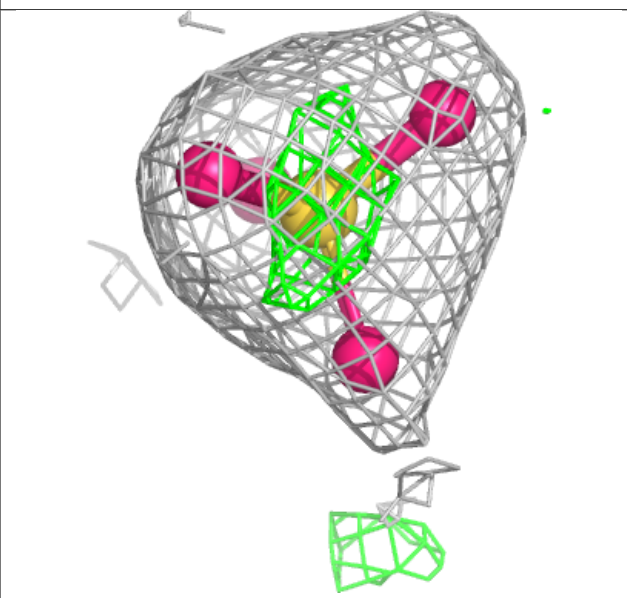
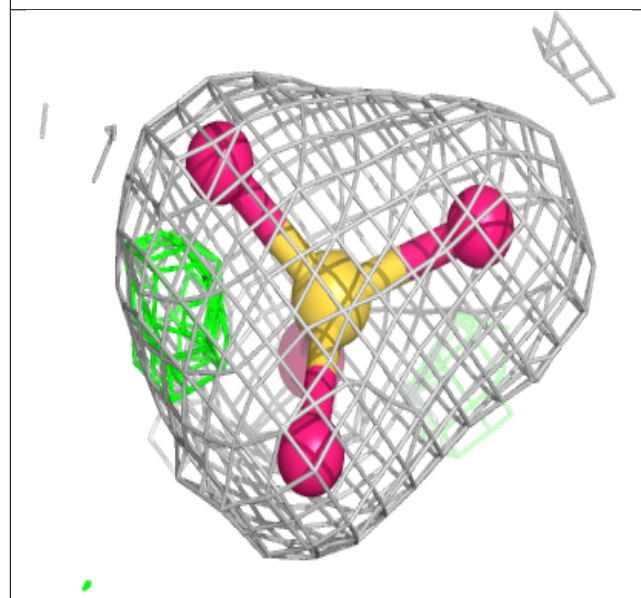
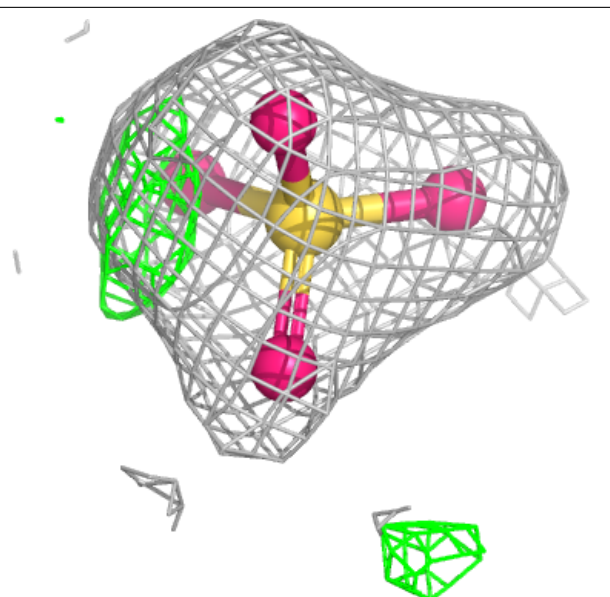
Electron density around NAG A 605:

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



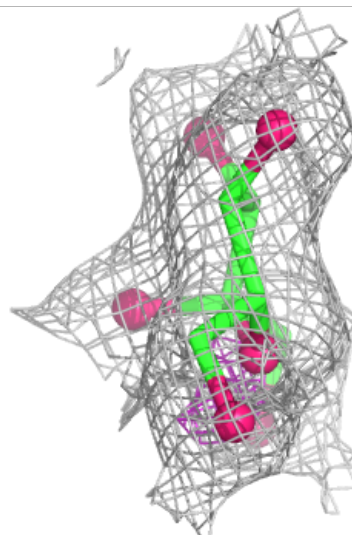
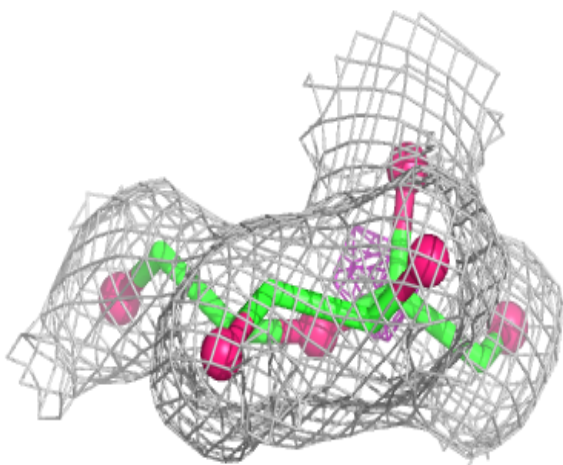
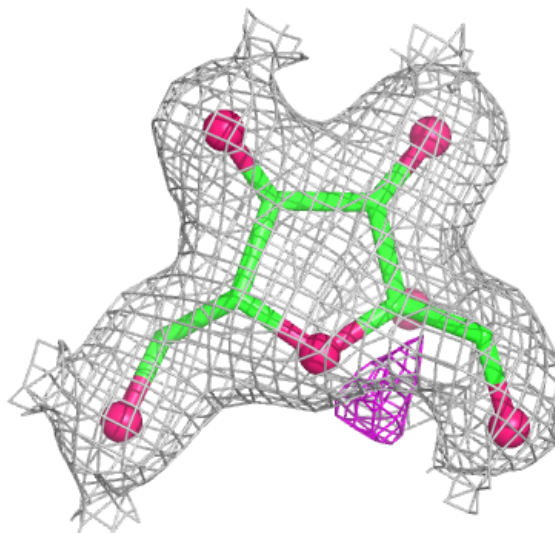
Electron density around SO4 C 601:

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



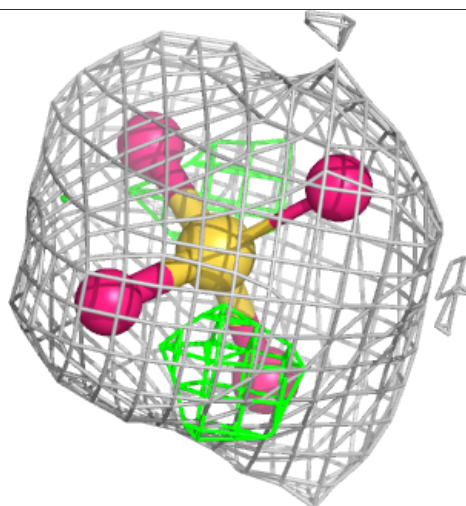
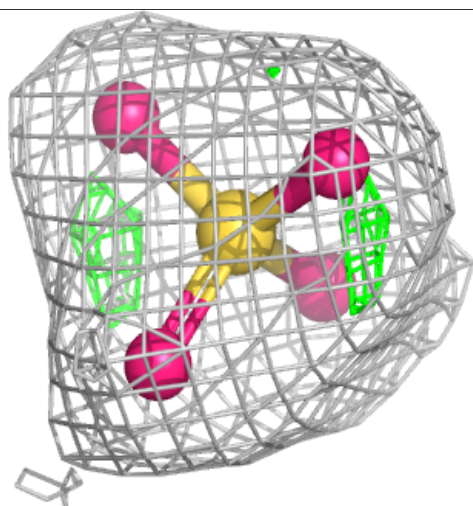
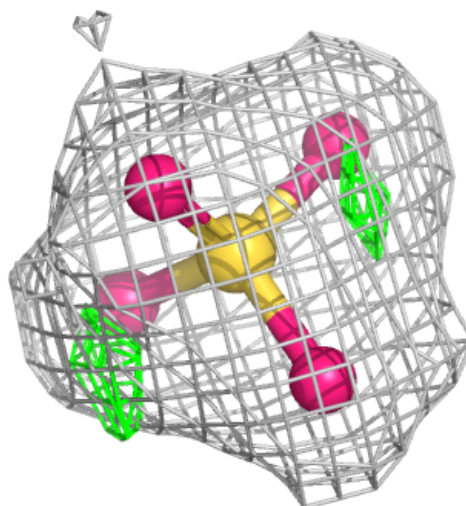
Electron density around FRU A 602:

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



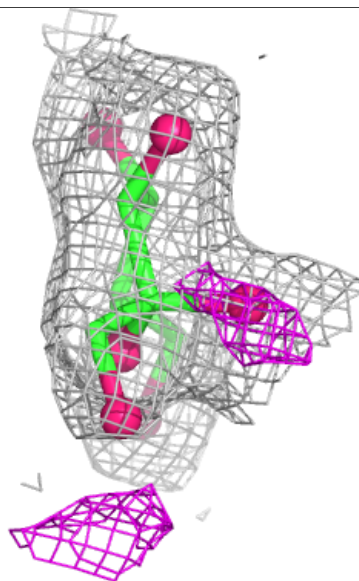
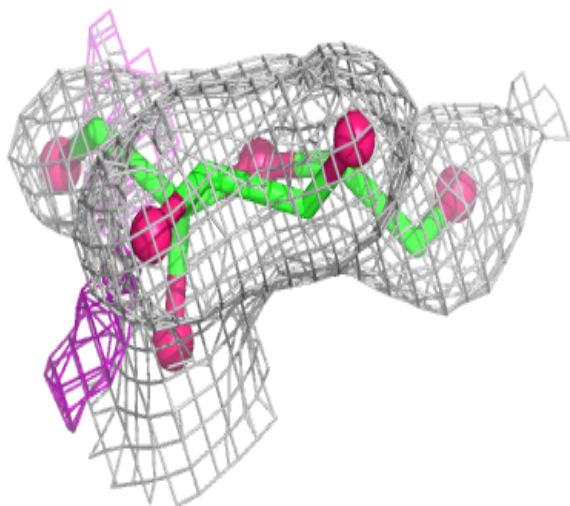
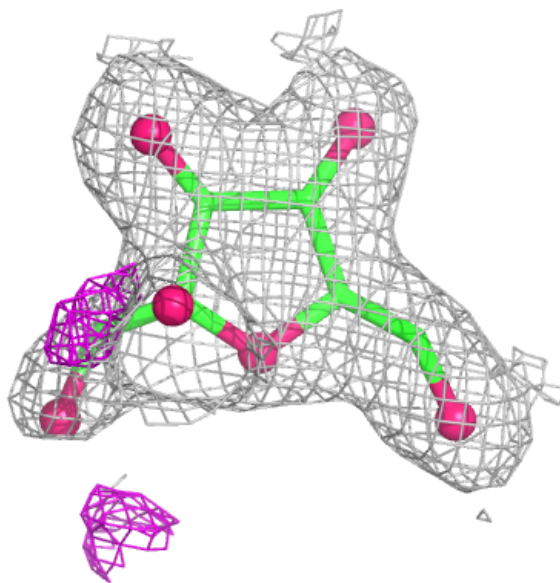
Electron density around SO4 D 604:

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



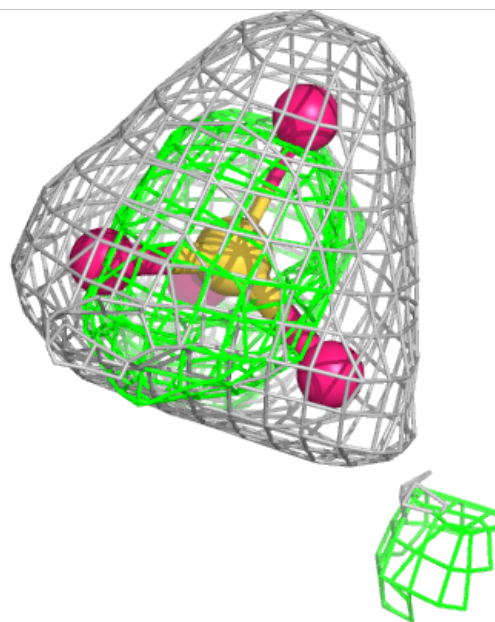
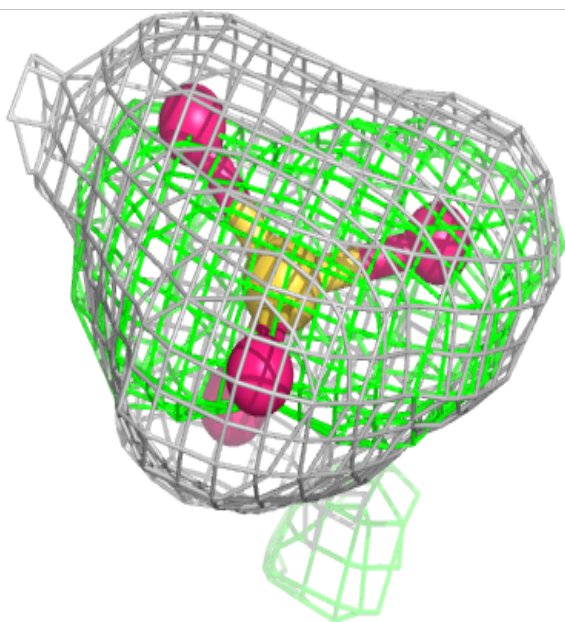
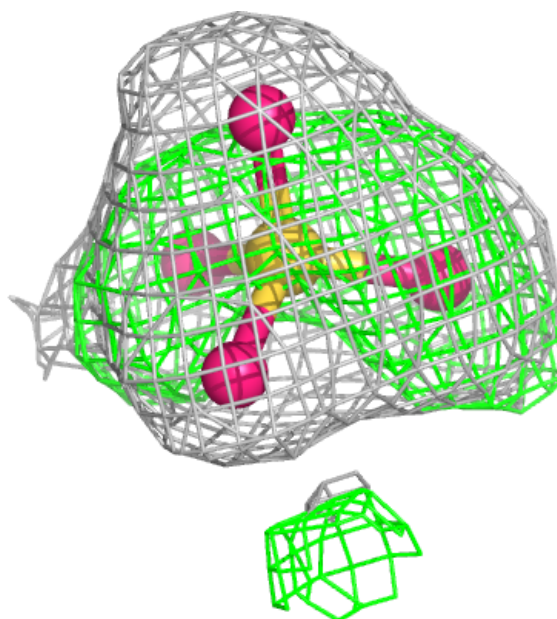
Electron density around FRU C 608:

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



Electron density around SO4 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)



6.5 Other polymers [i](#)

There are no such residues in this entry.