



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 01:15 PM UTC

PDB ID : 9HH0 / pdb_00009hh0
Title : Crystal structure of recombinant soman-aged swine butyrylcholinesterase
Authors : Brazzolotto, X.; Nachon, F.
Deposited on : 2024-11-20
Resolution : 2.66 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

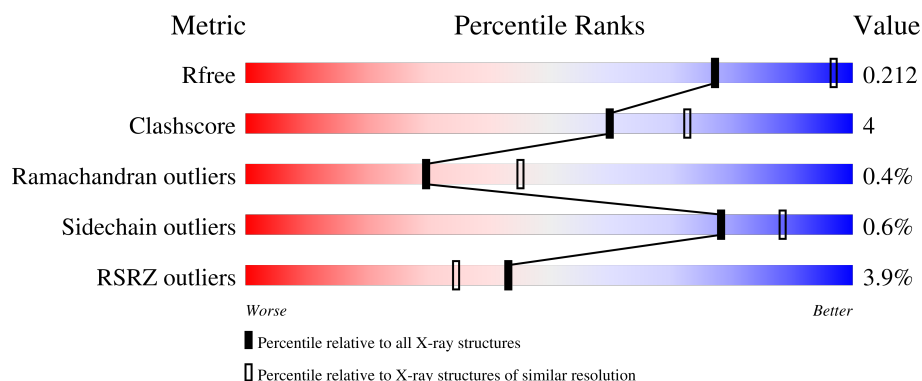
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.66 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	1110 (2.66-2.66)
Clashscore	190562	1141 (2.66-2.66)
Ramachandran outliers	187476	1126 (2.66-2.66)
Sidechain outliers	187428	1126 (2.66-2.66)
RSRZ outliers	180081	1110 (2.66-2.66)

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	529	<div> <div>2%</div> <div> <div></div> <div>89%</div> <div>10%</div> <div>.</div> </div> </div>
1	B	529	<div> <div>2%</div> <div> <div></div> <div>91%</div> <div>7%</div> <div>.</div> </div> </div>
1	C	529	<div> <div>4%</div> <div> <div></div> <div>87%</div> <div>11%</div> <div>.</div> </div> </div>
1	D	529	<div> <div>8%</div> <div> <div></div> <div>78%</div> <div>17%</div> <div>.</div> </div> </div>
2	E	2	<div> <div></div> <div> <div>50%</div> <div>50%</div> </div> </div>

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Mol	Chain	Length	Quality of chain
2	G	2	 100%
3	F	5	 20% 80%
4	H	2	 100%
4	I	2	 100%
4	L	2	 50% 50%
5	J	4	 50% 50%
6	K	3	 67% 33%

2 Entry composition

There are 13 unique types of molecules in this entry. The entry contains 17379 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 Carboxylic ester hydrolase.

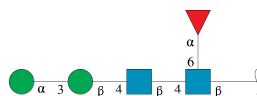
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	523	Total	C	N	O	S	0	1	0
			4188	2702	709	759	18			
1	B	522	Total	C	N	O	S	0	0	0
			4173	2693	704	758	18			
1	C	521	Total	C	N	O	S	0	1	0
			4172	2694	705	755	18			
1	D	508	Total	C	N	O	S	0	0	0
			4057	2617	685	739	16			

- Molecule 2 is an oligosaccharide called alpha-L-fucopyranose-(1-6)-2-acetamido-2-deoxy-beta-D-glucopyranose.



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	E	2	Total	C	N	O	0	0	0
			24	14	1	9			
2	G	2	Total	C	N	O	0	0	0
			24	14	1	9			

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



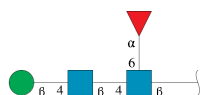
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	F	5	Total	C	N	O	0	0	0
			60	34	2	24			

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



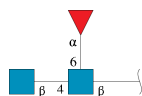
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
4	H	2	Total	C	N	O	0	0	0
			28	16	2	10			
4	I	2	Total	C	N	O	0	0	0
			28	16	2	10			
4	L	2	Total	C	N	O	0	0	0
			28	16	2	10			

- Molecule 5 is an oligosaccharide called beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-[alpha-L-fucopyranose-(1-6)]2-acetamido-2-deoxy-beta-D-glucopyranose.



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
5	J	4	Total	C	N	O	0	0	0
			49	28	2	19			

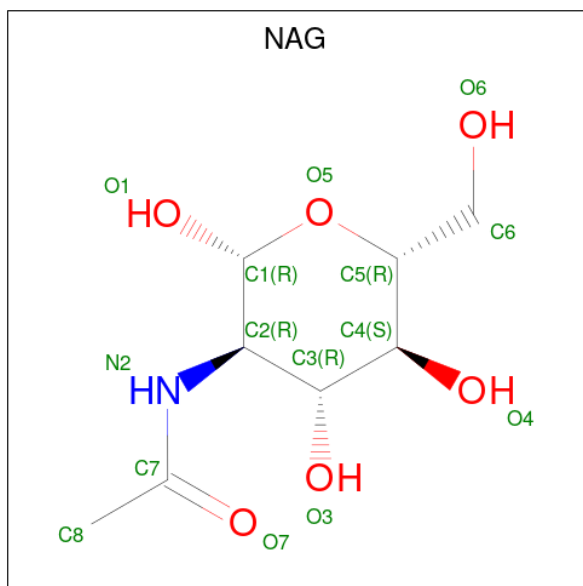
- Molecule 6 is an oligosaccharide called 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-[alpha-L-fucopyranose-(1-6)]2-acetamido-2-deoxy-beta-D-glucopyranose.



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
6	K	3	Total	C	N	O	0	0	0
			38	22	2	14			

- Molecule 7 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula:

C₈H₁₅NO₆).



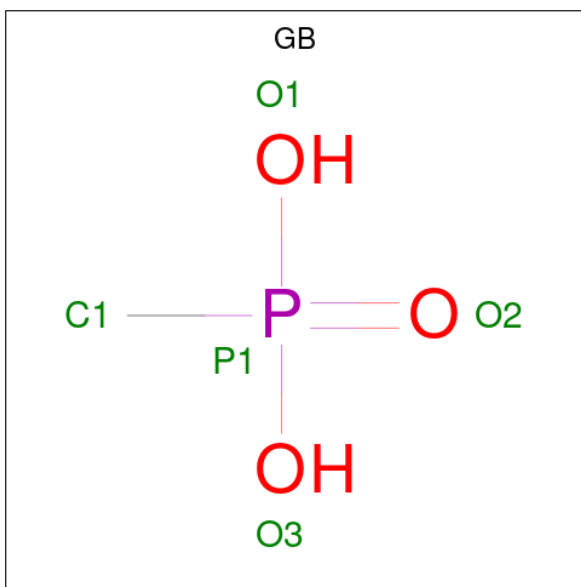
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
7	A	1	Total	C	N	O	0	0
			14	8	1	5		
7	A	1	Total	C	N	O	0	0
			14	8	1	5		
7	A	1	Total	C	N	O	0	0
			14	8	1	5		
7	B	1	Total	C	N	O	0	0
			14	8	1	5		
7	B	1	Total	C	N	O	0	0
			14	8	1	5		

- Molecule 8 is GLYCEROL (CCD ID: GOL) (formula: C₃H₈O₃).



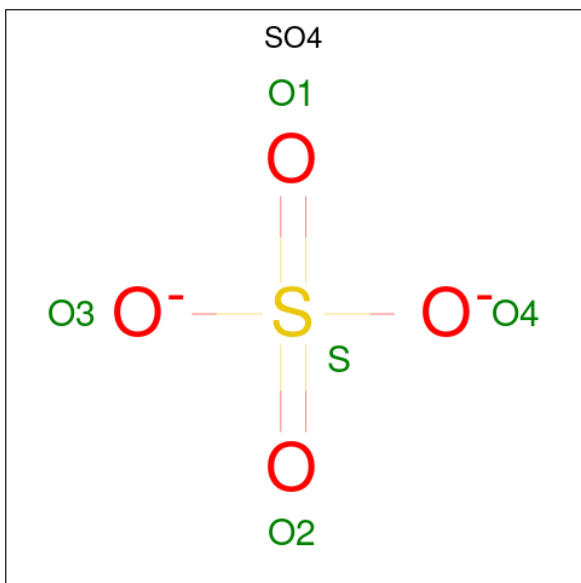
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
8	A	1	Total	C	O	0	0
			6	3	3		
8	A	1	Total	C	O	0	0
			6	3	3		
8	A	1	Total	C	O	0	0
			6	3	3		
8	A	1	Total	C	O	0	0
			6	3	3		
8	B	1	Total	C	O	0	0
			6	3	3		
8	C	1	Total	C	O	0	0
			6	3	3		
8	C	1	Total	C	O	0	0
			6	3	3		

- Molecule 9 is METHYLPHOSPHONIC ACID ESTER GROUP (CCD ID: GB) (formula: $\text{CH}_5\text{O}_3\text{P}$) (labeled as "Ligand of Interest" by depositor).



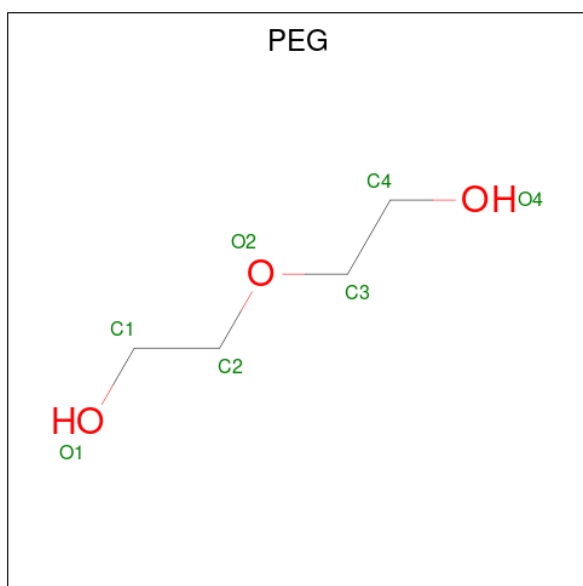
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
9	A	1	Total	C	O	P	0	0
			4	1	2	1		
9	B	1	Total	C	O	P	0	0
			4	1	2	1		
9	C	1	Total	C	O	P	0	0
			4	1	2	1		
9	D	1	Total	C	O	P	0	0
			4	1	2	1		

- Molecule 10 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



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

- Molecule 11 is DI(HYDROXYETHYL)ETHER (CCD ID: PEG) (formula: $C_4H_{10}O_3$).



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

- Molecule 12 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: $C_2H_6O_2$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
12	B	1	Total	C	O	0	0
			4	2	2		
12	B	1	Total	C	O	0	0
			4	2	2		
12	B	1	Total	C	O	0	0
			4	2	2		
12	D	1	Total	C	O	0	0
			4	2	2		

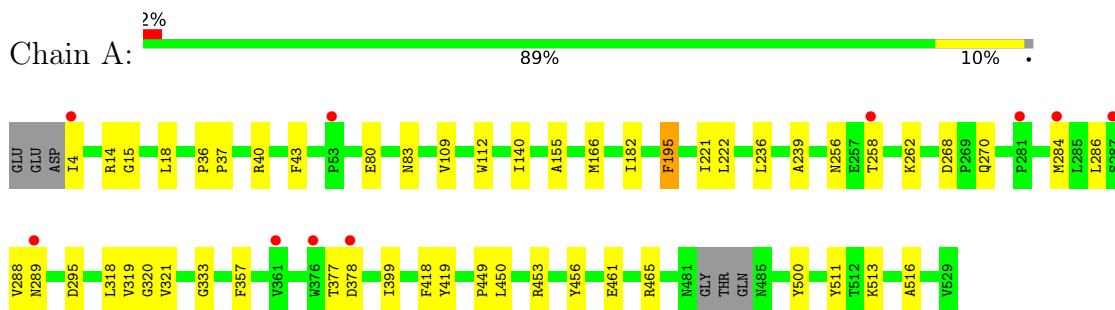
- Molecule 13 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
13	A	122	Total	O	0	0
			122	122		
13	B	110	Total	O	0	0
			110	110		
13	C	73	Total	O	0	0
			73	73		
13	D	19	Total	O	0	0
			19	19		

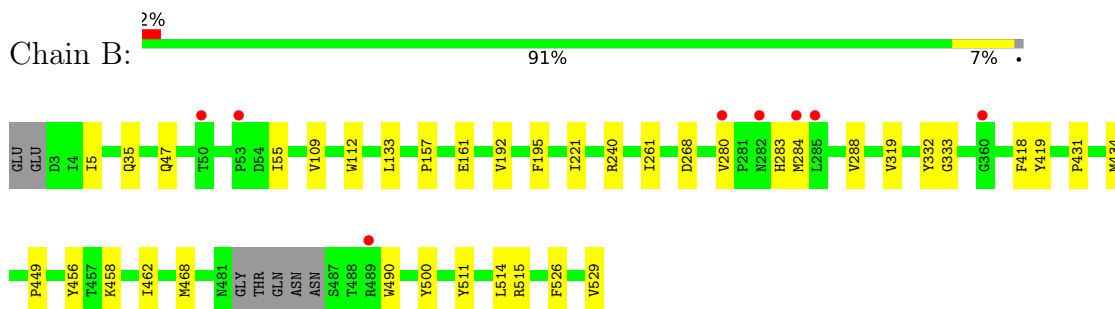
3 Residue-property plots [i](#)

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.

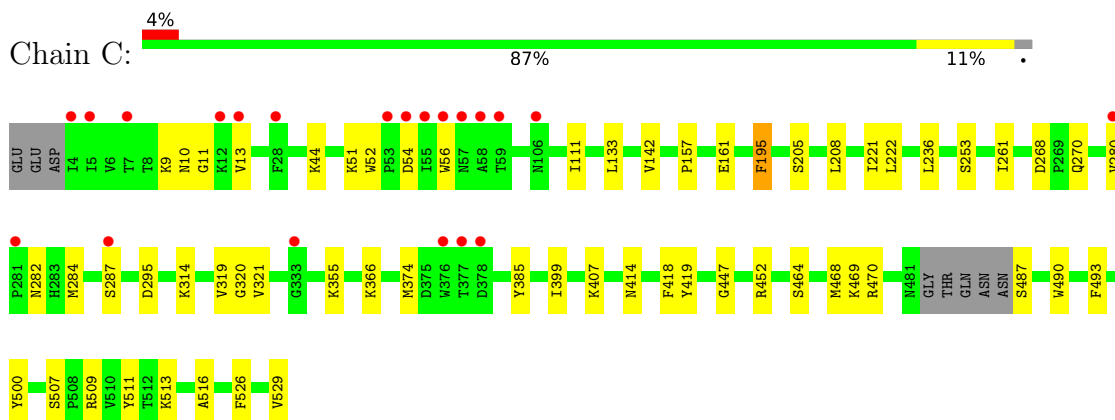
- Molecule 1: Carboxylic ester hydrolase



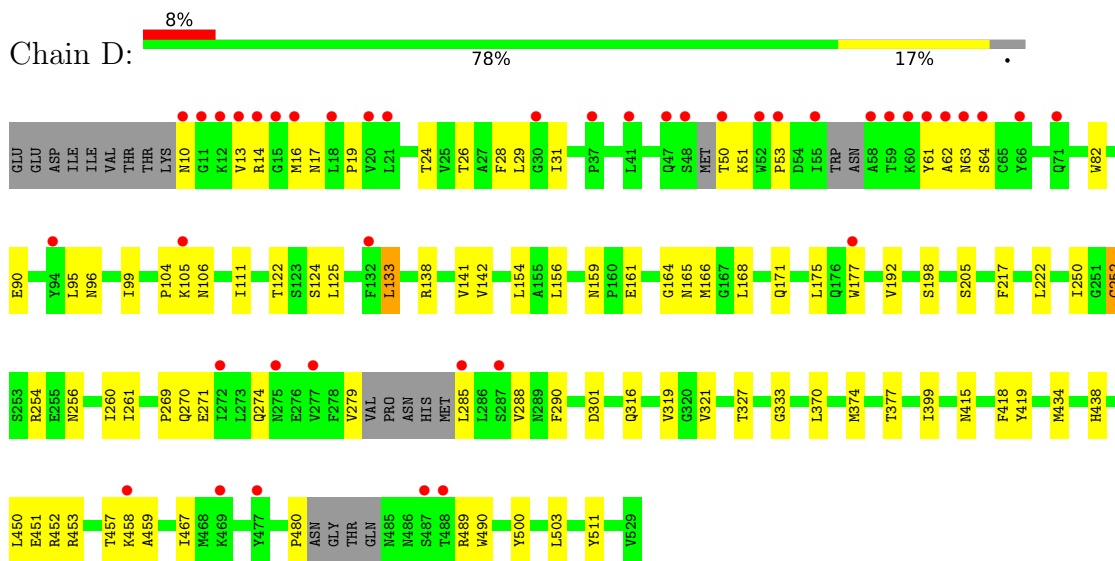
- Molecule 1: Carboxylic ester hydrolase



- Molecule 1: Carboxylic ester hydrolase



- Molecule 1: Carboxylic ester hydrolase



- Molecule 2: alpha-L-fucopyranose-(1-6)-2-acetamido-2-deoxy-beta-D-glucopyranose



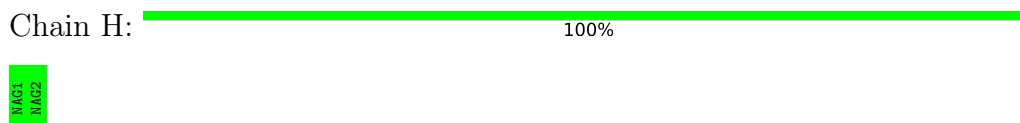
- Molecule 2: alpha-L-fucopyranose-(1-6)-2-acetamido-2-deoxy-beta-D-glucopyranose



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



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



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





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

Chain L:



- Molecule 5: beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-[alpha-L-fucopyranose-(1-6)]2-acetamido-2-deoxy-beta-D-glucopyranose

Chain J:



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

Chain K:



4 Data and refinement statistics

Property	Value	Source
Space group	P 41	Depositor
Cell constants a, b, c, α , β , γ	124.69Å 124.69Å 175.17Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	124.69 – 2.66 124.69 – 2.66	Depositor EDS
% Data completeness (in resolution range)	99.8 (124.69-2.66) 99.8 (124.69-2.66)	Depositor EDS
R_{merge}	0.15	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.54 (at 2.65Å)	Xtriage
Refinement program	PHENIX 1.21.2_5419	Depositor
R, R_{free}	0.195 , 0.236 (Not available) , 0.212	Depositor DCC
R_{free} test set	3953 reflections (5.18%)	wwPDB-VP
Wilson B-factor (Å ²)	41.6	Xtriage
Anisotropy	0.132	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 53.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.033 for h,-k,-l	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	17379	wwPDB-VP
Average B, all atoms (Å ²)	52.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.89% 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: PEG, GOL, EDO, SO4, NAG, GB, FUC, MAN, BMA

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	0.13	0/4309	0.31	0/5847
1	B	0.13	0/4290	0.31	0/5821
1	C	0.12	0/4293	0.30	0/5825
1	D	0.13	0/4167	0.32	0/5647
All	All	0.13	0/17059	0.31	0/23140

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	4188	0	4103	34	0
1	B	4173	0	4089	19	0
1	C	4172	0	4093	35	0
1	D	4057	0	3970	57	0
2	E	24	0	22	0	0
2	G	24	0	22	0	0
3	F	60	0	52	0	0
4	H	28	0	25	0	0
4	I	28	0	25	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	L	28	0	25	0	0
5	J	49	0	43	0	0
6	K	38	0	34	0	0
7	A	42	0	39	0	0
7	B	28	0	26	0	0
8	A	24	0	32	1	0
8	B	6	0	8	0	0
8	C	12	0	16	0	0
9	A	4	0	3	0	0
9	B	4	0	3	0	0
9	C	4	0	3	0	0
9	D	4	0	3	0	0
10	A	20	0	0	0	0
10	B	5	0	0	0	0
10	C	10	0	0	0	0
11	B	7	0	10	3	0
12	B	12	0	18	0	0
12	D	4	0	6	0	0
13	A	122	0	0	0	0
13	B	110	0	0	0	0
13	C	73	0	0	0	0
13	D	19	0	0	0	0
All	All	17379	0	16670	145	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 4.

All (145) 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:268:ASP:HB3	1:C:270:GLN:HE21	1.44	0.81
1:C:507:SER:HB3	1:C:509:ARG:HH21	1.55	0.72
1:D:122:THR:HG22	1:D:124:SER:H	1.55	0.71
1:A:284:MET:HE2	1:A:286:LEU:H	1.55	0.71
1:A:453:ARG:H	1:A:453:ARG:HD2	1.56	0.70
1:C:205:SER:HA	1:C:208:LEU:HD12	1.74	0.69
1:C:513:LYS:HB3	1:C:516:ALA:HB2	1.76	0.68
1:A:80:GLU:HB3	11:B:603:PEG:H41	1.81	0.63
1:D:154:LEU:O	1:D:166:MET:HE1	1.99	0.62
1:C:157:PRO:HB3	1:C:236:LEU:HD13	1.82	0.61
1:A:83:ASN:HD22	11:B:603:PEG:H32	1.65	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:489:ARG:HE	1:D:490:TRP:N	1.99	0.60
1:C:222:LEU:HB2	1:C:319:VAL:HG13	1.83	0.59
1:D:164:GLY:H	1:D:166:MET:HE3	1.67	0.59
1:D:104:PRO:HB3	1:D:138:ARG:HH12	1.66	0.59
1:A:256:ASN:HD21	1:A:258:THR:HG22	1.67	0.59
1:B:133:LEU:HD13	1:B:468:MET:HG2	1.85	0.58
1:D:205:SER:HB3	1:D:222:LEU:HD21	1.84	0.58
1:C:133:LEU:HD23	1:C:468:MET:HE2	1.88	0.56
1:D:26:THR:HG22	1:D:99:ILE:HB	1.86	0.56
1:C:236:LEU:HD21	1:C:295:ASP:HA	1.88	0.56
1:D:161:GLU:HB2	1:D:261:ILE:HD12	1.88	0.56
1:A:513:LYS:HB3	1:A:516:ALA:HB2	1.88	0.55
1:D:192:VAL:HB	1:D:217:PHE:HA	1.88	0.54
1:C:284:MET:H	1:C:287:SER:HB3	1.72	0.54
1:D:111:ILE:HG12	1:D:142:VAL:HB	1.90	0.54
1:C:161:GLU:HB3	1:C:261:ILE:HG21	1.90	0.54
1:C:407:LYS:HG3	1:C:493:PHE:HE1	1.73	0.53
1:A:450:LEU:HD11	1:A:465:ARG:HG3	1.91	0.53
1:D:489:ARG:HE	1:D:490:TRP:H	1.55	0.53
1:A:268:ASP:HB3	1:A:270:GLN:OE1	2.09	0.52
1:B:458:LYS:O	1:B:462:ILE:HG12	2.09	0.52
1:D:319:VAL:O	1:D:418:PHE:HA	2.10	0.52
1:D:28:PHE:HB3	1:D:31:ILE:HD11	1.91	0.52
1:A:256:ASN:ND2	1:A:258:THR:HG22	2.25	0.51
1:A:222:LEU:HB2	1:A:319:VAL:HG13	1.91	0.51
1:B:161:GLU:HB2	1:B:261:ILE:HD13	1.92	0.51
1:D:254:ARG:H	1:D:260:ILE:HG13	1.76	0.50
1:C:10:ASN:HA	1:C:51:LYS:HD3	1.94	0.50
1:C:319:VAL:O	1:C:418:PHE:HA	2.12	0.50
1:D:316:GLN:HG2	1:D:415:ASN:HB2	1.93	0.50
1:C:13:VAL:HG12	1:C:56:TRP:HB3	1.94	0.49
1:D:95:LEU:HD21	1:D:177:TRP:CD2	2.47	0.49
1:D:457:THR:HG22	1:D:459:ALA:H	1.77	0.49
1:D:13:VAL:HG22	1:D:14:ARG:H	1.76	0.49
1:B:109:VAL:HB	1:B:192:VAL:HG22	1.94	0.49
1:C:314:LYS:HD3	1:C:414:ASN:HD21	1.78	0.49
1:A:450:LEU:HA	1:A:461:GLU:HG3	1.95	0.49
1:D:457:THR:HG22	1:D:458:LYS:HD2	1.94	0.49
1:D:480:PRO:C	1:D:489:ARG:HH22	2.22	0.48
1:A:221:ILE:HG12	1:A:318:LEU:HB3	1.95	0.48
1:A:284:MET:HE1	1:A:357:PHE:CE1	2.49	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:19:PRO:HA	1:D:24:THR:HA	1.96	0.48
1:A:4:ILE:HD11	1:A:14:ARG:HG3	1.96	0.48
1:D:489:ARG:HE	1:D:489:ARG:HA	1.79	0.48
1:A:43:PHE:HA	1:A:166:MET:HE3	1.95	0.47
1:A:83:ASN:ND2	11:B:603:PEG:H32	2.27	0.47
1:C:195:PHE:HB3	1:C:221:ILE:HD12	1.96	0.47
1:D:105:LYS:HG2	1:D:106:ASN:H	1.80	0.47
1:D:105:LYS:HG2	1:D:106:ASN:N	2.30	0.47
1:D:500:TYR:CZ	1:D:511:TYR:HB2	2.50	0.47
1:A:18:LEU:HD13	8:A:604:GOL:H11	1.96	0.47
1:A:319:VAL:O	1:A:418:PHE:HA	2.14	0.47
1:B:500:TYR:CZ	1:B:511:TYR:HB2	2.50	0.47
1:D:90:GLU:OE2	1:D:269:PRO:HG3	2.14	0.46
1:D:165:ASN:HB3	1:D:168:LEU:HD12	1.97	0.46
1:C:111:ILE:HG12	1:C:142:VAL:HB	1.96	0.46
1:D:222:LEU:HB2	1:D:319:VAL:HG22	1.97	0.46
1:A:40:ARG:NH1	1:A:262:LYS:HE3	2.31	0.46
1:D:29:LEU:HG	1:D:96:ASN:HD22	1.79	0.46
1:C:419:TYR:HB3	1:C:490:TRP:CZ2	2.50	0.46
1:B:319:VAL:O	1:B:418:PHE:HA	2.16	0.45
1:D:50:THR:OG1	1:D:51:LYS:HD3	2.16	0.45
1:D:159:ASN:HD21	1:D:261:ILE:HD13	1.82	0.45
1:B:514:LEU:HG	1:B:515:ARG:HG2	1.98	0.45
1:D:16:MET:C	1:D:17:ASN:HD22	2.24	0.45
1:B:35:GLN:HB3	1:B:47:GLN:HB2	1.99	0.45
1:D:61:TYR:CE1	1:D:124:SER:HA	2.51	0.45
1:D:104:PRO:HB3	1:D:138:ARG:NH1	2.30	0.45
1:D:419:TYR:HB3	1:D:490:TRP:CZ2	2.52	0.45
1:B:526:PHE:O	1:B:529:VAL:HG22	2.16	0.45
1:D:250:ILE:HG23	1:D:252:CYS:HB2	1.98	0.45
1:D:250:ILE:HD11	1:D:271:GLU:HB3	1.99	0.45
1:A:321:VAL:HG11	1:A:399:ILE:HA	1.99	0.45
1:A:4:ILE:HD13	1:A:15:GLY:H	1.82	0.44
1:B:419:TYR:HB3	1:B:490:TRP:CZ2	2.52	0.44
1:D:270:GLN:O	1:D:274:GLN:HG2	2.18	0.44
1:D:467:ILE:HD12	1:D:503:LEU:HD13	1.99	0.44
1:A:36:PRO:HA	1:A:37:PRO:HD3	1.81	0.44
1:C:407:LYS:HG3	1:C:493:PHE:CE1	2.53	0.44
1:C:447:GLY:HA2	1:C:464:SER:OG	2.18	0.44
1:D:82:TRP:HA	1:D:125:LEU:HD21	1.99	0.44
1:B:284:MET:HA	1:B:284:MET:HE2	2.00	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:453:ARG:HD2	1:A:453:ARG:N	2.29	0.43
1:D:452:ARG:HE	1:D:452:ARG:HB3	1.50	0.43
1:A:112:TRP:HA	1:A:195:PHE:O	2.18	0.43
1:A:320:GLY:HA3	1:A:419:TYR:CD2	2.54	0.43
1:B:284:MET:HE1	1:B:332:TYR:O	2.18	0.43
1:C:500:TYR:CZ	1:C:511:TYR:HB2	2.53	0.43
1:A:500:TYR:CZ	1:A:511:TYR:HB2	2.53	0.43
1:D:327:THR:HB	1:D:434:MET:HG2	2.01	0.43
1:C:470:ARG:HH22	1:C:487:SER:HB3	1.84	0.43
1:C:526:PHE:O	1:C:529:VAL:HG22	2.19	0.43
1:D:451:GLU:HB2	1:D:453:ARG:CZ	2.49	0.43
1:A:320:GLY:HA3	1:A:419:TYR:CE2	2.54	0.42
1:C:452:ARG:HD3	1:C:452:ARG:HA	1.64	0.42
1:D:171:GLN:O	1:D:175:LEU:HG	2.20	0.42
1:A:449:PRO:HA	1:A:456:TYR:CD1	2.54	0.42
1:B:280:VAL:HG12	1:B:283:HIS:H	1.85	0.42
1:D:133:LEU:HG	1:D:141:VAL:HG21	2.01	0.42
1:D:453:ARG:H	1:D:453:ARG:HG2	1.61	0.42
1:A:236:LEU:HD13	1:A:295:ASP:HA	2.00	0.42
1:B:449:PRO:HA	1:B:456:TYR:CD1	2.54	0.42
1:C:11:GLY:HA3	1:C:52:TRP:CE2	2.54	0.42
1:A:270:GLN:H	1:A:270:GLN:CD	2.27	0.41
1:B:195:PHE:CB	1:B:221:ILE:HB	2.50	0.41
1:D:133:LEU:H	1:D:133:LEU:HD23	1.85	0.41
1:D:161:GLU:H	1:D:161:GLU:HG2	1.70	0.41
1:D:321:VAL:HG11	1:D:399:ILE:HA	2.02	0.41
1:A:155:ALA:HB3	1:A:239:ALA:HB1	2.02	0.41
1:C:469:LYS:HE2	1:C:469:LYS:HB2	1.96	0.41
1:C:355:LYS:HG3	1:C:366:LYS:NZ	2.36	0.41
1:B:5:ILE:HD12	1:B:55:ILE:HD13	2.03	0.41
1:C:9:LYS:HD3	1:C:9:LYS:N	2.36	0.41
1:D:489:ARG:HE	1:D:489:ARG:CA	2.33	0.41
1:A:109:VAL:HG22	1:A:140:ILE:HB	2.03	0.41
1:D:10:ASN:CG	1:D:50:THR:HG22	2.45	0.41
1:A:182:ILE:HD12	1:A:182:ILE:HA	1.90	0.41
1:B:112:TRP:HA	1:B:195:PHE:O	2.21	0.41
1:C:195:PHE:CB	1:C:221:ILE:HB	2.51	0.41
1:C:374:MET:HE2	1:C:385:TYR:CE1	2.55	0.41
1:D:279:VAL:HG21	1:D:290:PHE:CE1	2.56	0.41
1:C:133:LEU:HD23	1:C:133:LEU:HA	1.94	0.41
1:C:280:VAL:HG12	1:C:282:ASN:H	1.86	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:377:THR:HG22	1:A:378:ASP:N	2.37	0.40
1:B:431:PRO:HD2	1:B:434:MET:HE3	2.02	0.40
1:D:256:ASN:O	1:D:260:ILE:HD12	2.21	0.40
1:B:157:PRO:HD2	1:B:240:ARG:HD2	2.02	0.40
1:C:321:VAL:HG11	1:C:399:ILE:HA	2.02	0.40
1:D:198:SER:HB2	1:D:438:HIS:CE1	2.57	0.40
1:D:370:LEU:HD23	1:D:374:MET:HB2	2.03	0.40
1:D:63:ASN:HB3	1:D:64:SER:H	1.74	0.40
1:C:44:LYS:HE3	1:C:44:LYS:HB2	1.76	0.40
1:C:320:GLY:HA3	1:C:419:TYR:CE2	2.57	0.40
1:D:450:LEU:HD13	1:D:450:LEU:HA	1.95	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	520/529 (98%)	503 (97%)	15 (3%)	2 (0%)	30	45
1	B	518/529 (98%)	500 (96%)	16 (3%)	2 (0%)	30	45
1	C	518/529 (98%)	490 (95%)	27 (5%)	1 (0%)	43	60
1	D	498/529 (94%)	462 (93%)	32 (6%)	4 (1%)	16	27
All	All	2054/2116 (97%)	1955 (95%)	90 (4%)	9 (0%)	30	45

All (9) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	53	PRO
1	B	333	GLY
1	C	54	ASP
1	A	333	GLY

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Mol	Chain	Res	Type
1	D	62	ALA
1	D	333	GLY
1	D	288	VAL
1	A	288	VAL
1	B	288	VAL

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	452/456 (99%)	450 (100%)	2 (0%)	84	93
1	B	450/456 (99%)	449 (100%)	1 (0%)	87	95
1	C	450/456 (99%)	448 (100%)	2 (0%)	84	93
1	D	436/456 (96%)	430 (99%)	6 (1%)	59	76
All	All	1788/1824 (98%)	1777 (99%)	11 (1%)	78	88

All (11) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	195	PHE
1	A	289	ASN
1	B	268	ASP
1	C	195	PHE
1	C	253	SER
1	D	133	LEU
1	D	156	LEU
1	D	252	CYS
1	D	285	LEU
1	D	301	ASP
1	D	377	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	172	GLN
1	A	256	ASN
1	B	17	ASN
1	B	172	GLN
1	B	179	GLN
1	C	71	GLN
1	C	270	GLN
1	D	17	ASN
1	D	270	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

22 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	2,1	14,14,15	0.70	0	17,19,21	0.90	1 (5%)
2	FUC	E	2	2	10,10,11	0.76	0	14,14,16	0.95	0
3	NAG	F	1	3,1	14,14,15	0.69	0	17,19,21	1.12	1 (5%)
3	NAG	F	2	3	14,14,15	0.69	0	17,19,21	1.05	1 (5%)
3	BMA	F	3	3	11,11,12	0.89	0	15,15,17	2.40	5 (33%)
3	MAN	F	4	3	11,11,12	0.63	0	15,15,17	1.36	1 (6%)
3	FUC	F	5	3	10,10,11	0.76	0	14,14,16	1.00	0
2	NAG	G	1	2,1	14,14,15	0.70	0	17,19,21	0.83	0
2	FUC	G	2	2	10,10,11	0.76	0	14,14,16	0.99	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	NAG	H	1	1,4	14,14,15	0.69	0	17,19,21	0.95	0
4	NAG	H	2	4	14,14,15	0.70	0	17,19,21	0.82	0
4	NAG	I	1	1,4	14,14,15	0.79	0	17,19,21	2.81	6 (35%)
4	NAG	I	2	4	14,14,15	0.77	0	17,19,21	1.15	1 (5%)
5	NAG	J	1	1,5	14,14,15	0.72	0	17,19,21	1.13	0
5	NAG	J	2	5	14,14,15	0.74	0	17,19,21	1.23	1 (5%)
5	BMA	J	3	5	11,11,12	0.92	0	15,15,17	2.33	6 (40%)
5	FUC	J	4	5	10,10,11	0.75	0	14,14,16	1.00	0
6	NAG	K	1	1,6	14,14,15	0.70	0	17,19,21	0.92	0
6	NAG	K	2	6	14,14,15	0.75	0	17,19,21	0.95	1 (5%)
6	FUC	K	3	6	10,10,11	0.73	0	14,14,16	0.95	0
4	NAG	L	1	1,4	14,14,15	0.67	0	17,19,21	1.01	1 (5%)
4	NAG	L	2	4	14,14,15	0.72	0	17,19,21	0.92	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	2,1	-	0/6/23/26	0/1/1/1
2	FUC	E	2	2	-	-	0/1/1/1
3	NAG	F	1	3,1	-	0/6/23/26	0/1/1/1
3	NAG	F	2	3	-	2/6/23/26	0/1/1/1
3	BMA	F	3	3	-	1/2/19/22	0/1/1/1
3	MAN	F	4	3	-	2/2/19/22	0/1/1/1
3	FUC	F	5	3	-	-	0/1/1/1
2	NAG	G	1	2,1	-	0/6/23/26	0/1/1/1
2	FUC	G	2	2	-	-	0/1/1/1
4	NAG	H	1	1,4	-	1/6/23/26	0/1/1/1
4	NAG	H	2	4	-	0/6/23/26	0/1/1/1
4	NAG	I	1	1,4	-	3/6/23/26	0/1/1/1
4	NAG	I	2	4	-	2/6/23/26	0/1/1/1
5	NAG	J	1	1,5	-	0/6/23/26	0/1/1/1
5	NAG	J	2	5	-	0/6/23/26	0/1/1/1
5	BMA	J	3	5	-	0/2/19/22	0/1/1/1
5	FUC	J	4	5	-	-	0/1/1/1
6	NAG	K	1	1,6	-	2/6/23/26	0/1/1/1
6	NAG	K	2	6	-	0/6/23/26	0/1/1/1
6	FUC	K	3	6	-	-	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	NAG	L	1	1,4	-	2/6/23/26	0/1/1/1
4	NAG	L	2	4	-	0/6/23/26	0/1/1/1

There are no bond length outliers.

All (25) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	I	1	NAG	C2-N2-C7	7.66	133.17	122.90
3	F	3	BMA	C1-O5-C5	6.35	120.69	112.19
5	J	3	BMA	C1-O5-C5	5.55	119.62	112.19
4	I	1	NAG	C1-O5-C5	4.76	118.56	112.19
4	I	1	NAG	C1-C2-N2	4.38	117.33	110.43
3	F	4	MAN	C1-O5-C5	4.31	117.97	112.19
5	J	3	BMA	C3-C4-C5	4.00	117.49	110.23
3	F	3	BMA	C3-C4-C5	3.87	117.25	110.23
5	J	3	BMA	C2-C3-C4	3.62	117.23	110.86
4	I	1	NAG	C4-C3-C2	3.50	116.15	111.02
3	F	3	BMA	C2-C3-C4	2.99	116.12	110.86
4	I	1	NAG	O7-C7-N2	2.56	126.51	121.98
3	F	3	BMA	O5-C5-C4	2.42	116.72	110.83
4	L	1	NAG	C2-N2-C7	2.37	126.08	122.90
3	F	3	BMA	O4-C4-C3	-2.37	104.80	110.38
3	F	2	NAG	O5-C1-C2	-2.36	107.64	111.29
5	J	3	BMA	O4-C4-C3	-2.28	105.00	110.38
5	J	2	NAG	O4-C4-C3	-2.26	105.04	110.38
3	F	1	NAG	C1-O5-C5	2.25	115.20	112.19
4	I	1	NAG	O3-C3-C2	-2.24	104.75	109.40
4	I	2	NAG	C1-O5-C5	2.23	115.17	112.19
2	E	1	NAG	C1-O5-C5	2.18	115.11	112.19
5	J	3	BMA	O3-C3-C2	-2.12	105.72	110.05
5	J	3	BMA	O5-C5-C4	2.10	115.93	110.83
6	K	2	NAG	C1-O5-C5	2.02	114.89	112.19

There are no chirality outliers.

All (15) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	I	1	NAG	C1-C2-N2-C7
4	I	2	NAG	C4-C5-C6-O6
4	I	2	NAG	O5-C5-C6-O6
4	I	1	NAG	O5-C5-C6-O6

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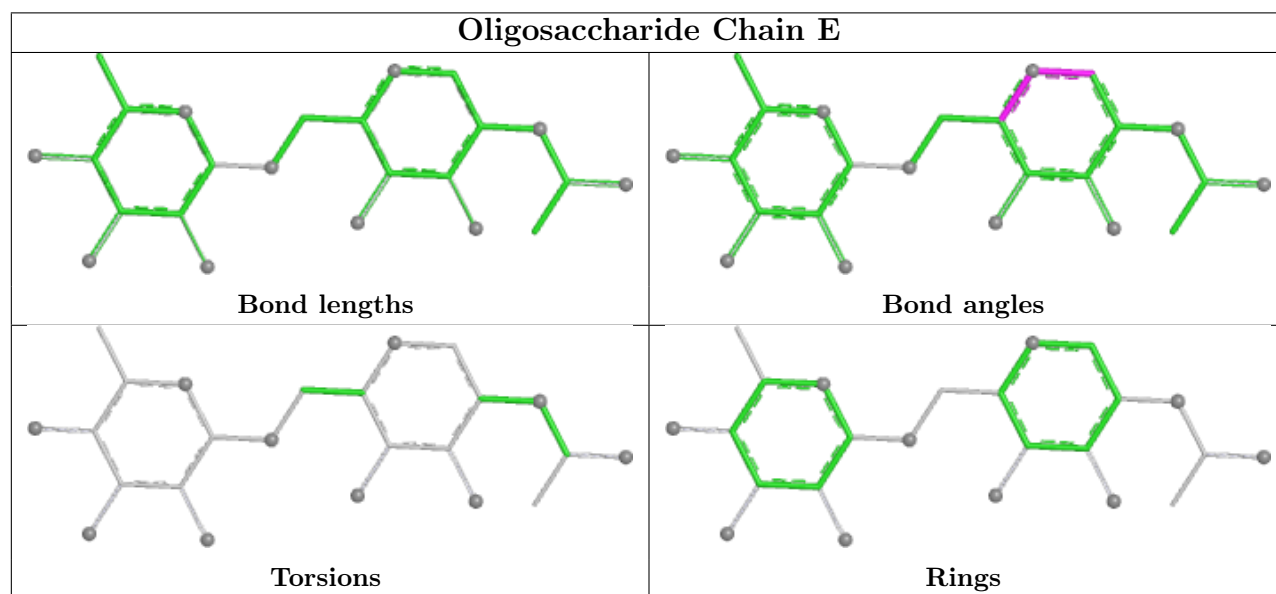
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Mol	Chain	Res	Type	Atoms
4	I	1	NAG	C4-C5-C6-O6
3	F	2	NAG	C8-C7-N2-C2
3	F	2	NAG	O7-C7-N2-C2
4	L	1	NAG	C8-C7-N2-C2
4	L	1	NAG	O7-C7-N2-C2
6	K	1	NAG	C8-C7-N2-C2
6	K	1	NAG	O7-C7-N2-C2
3	F	4	MAN	C4-C5-C6-O6
3	F	3	BMA	C4-C5-C6-O6
3	F	4	MAN	O5-C5-C6-O6
4	H	1	NAG	C1-C2-N2-C7

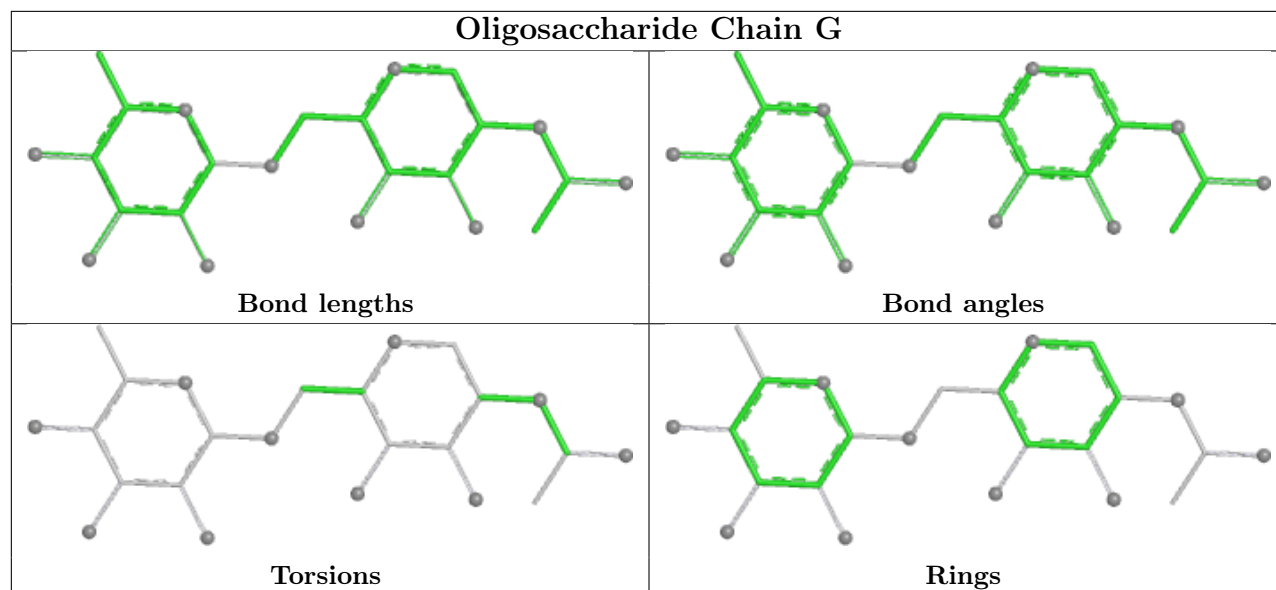
There are no ring outliers.

No monomer is involved in short contacts.

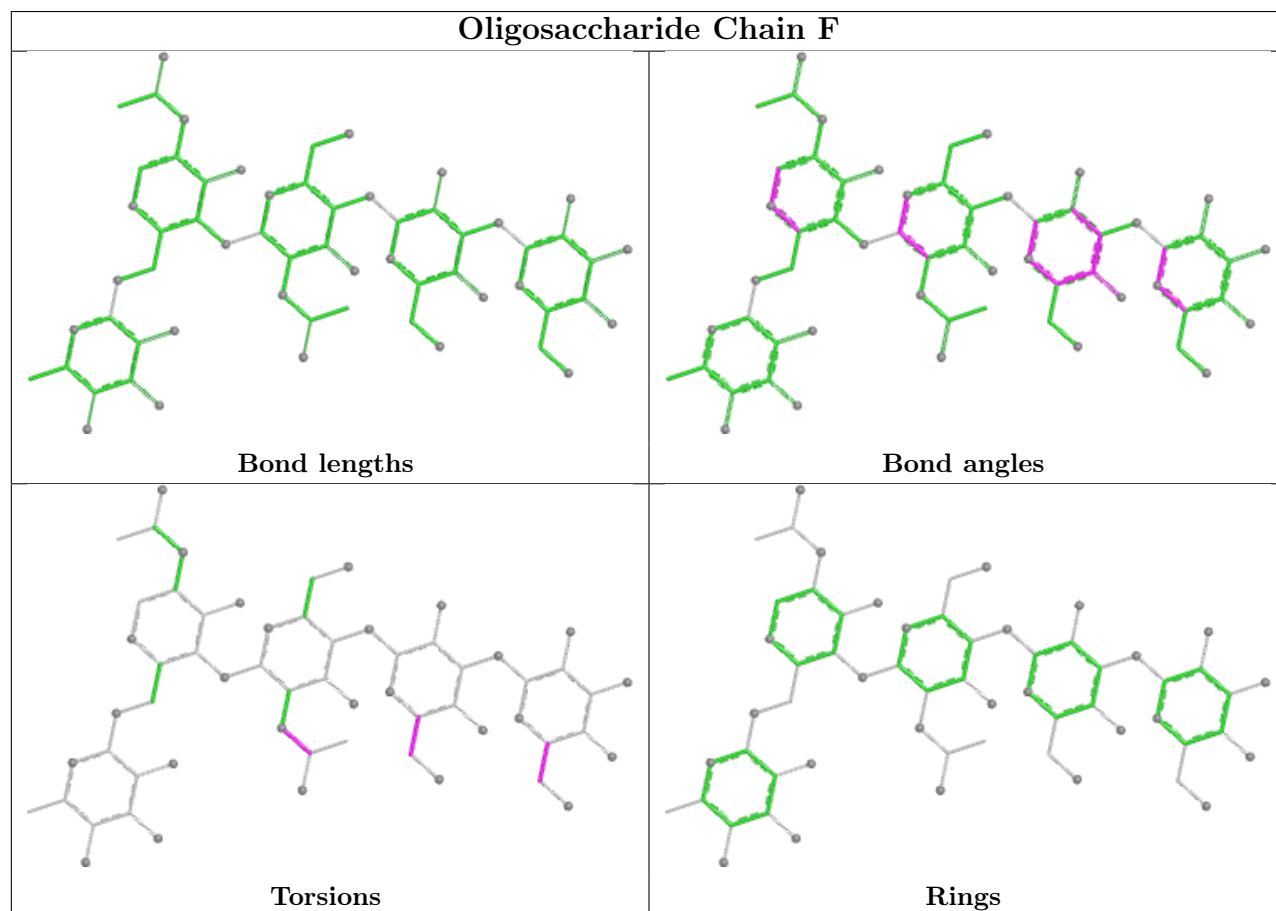
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.

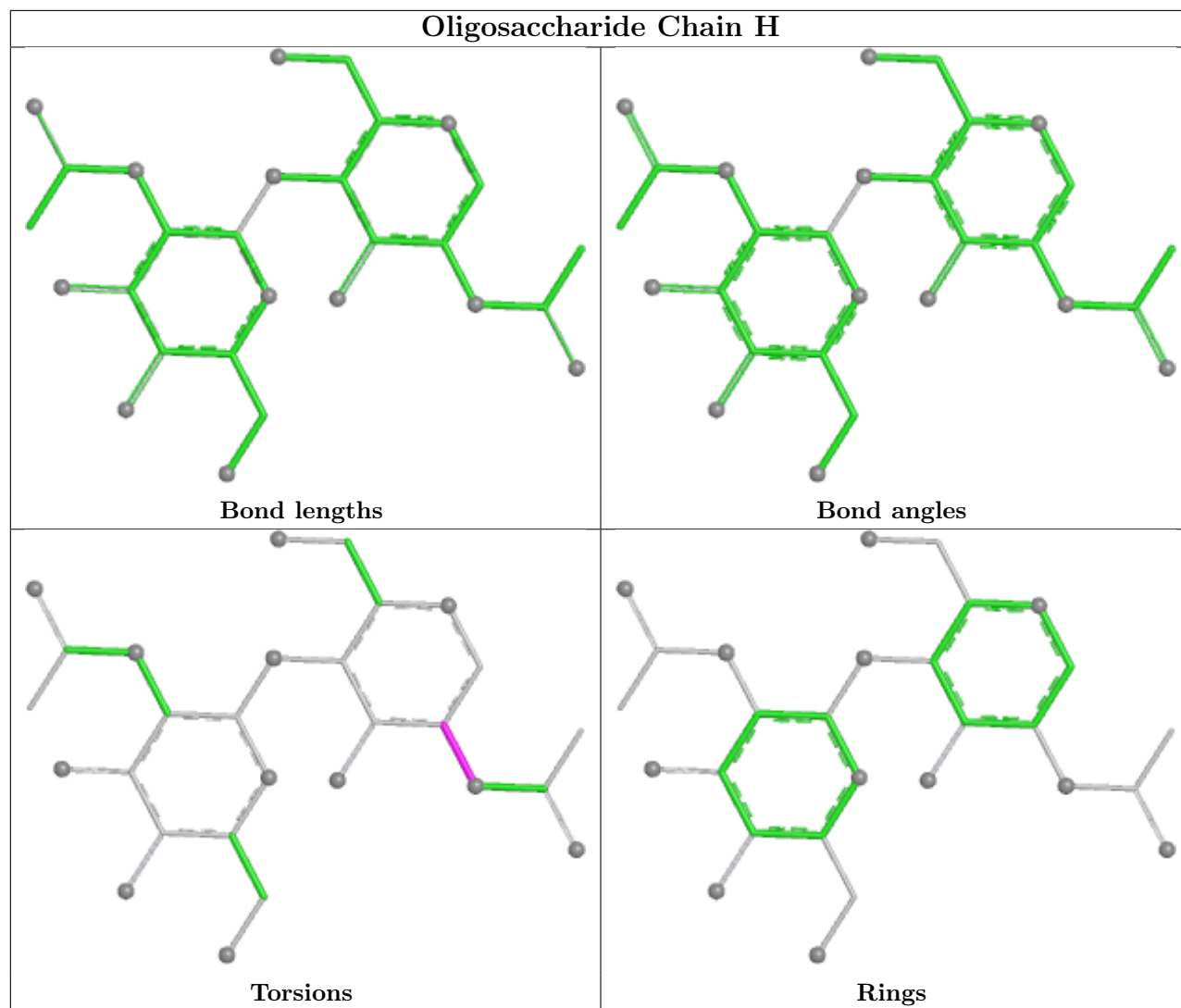


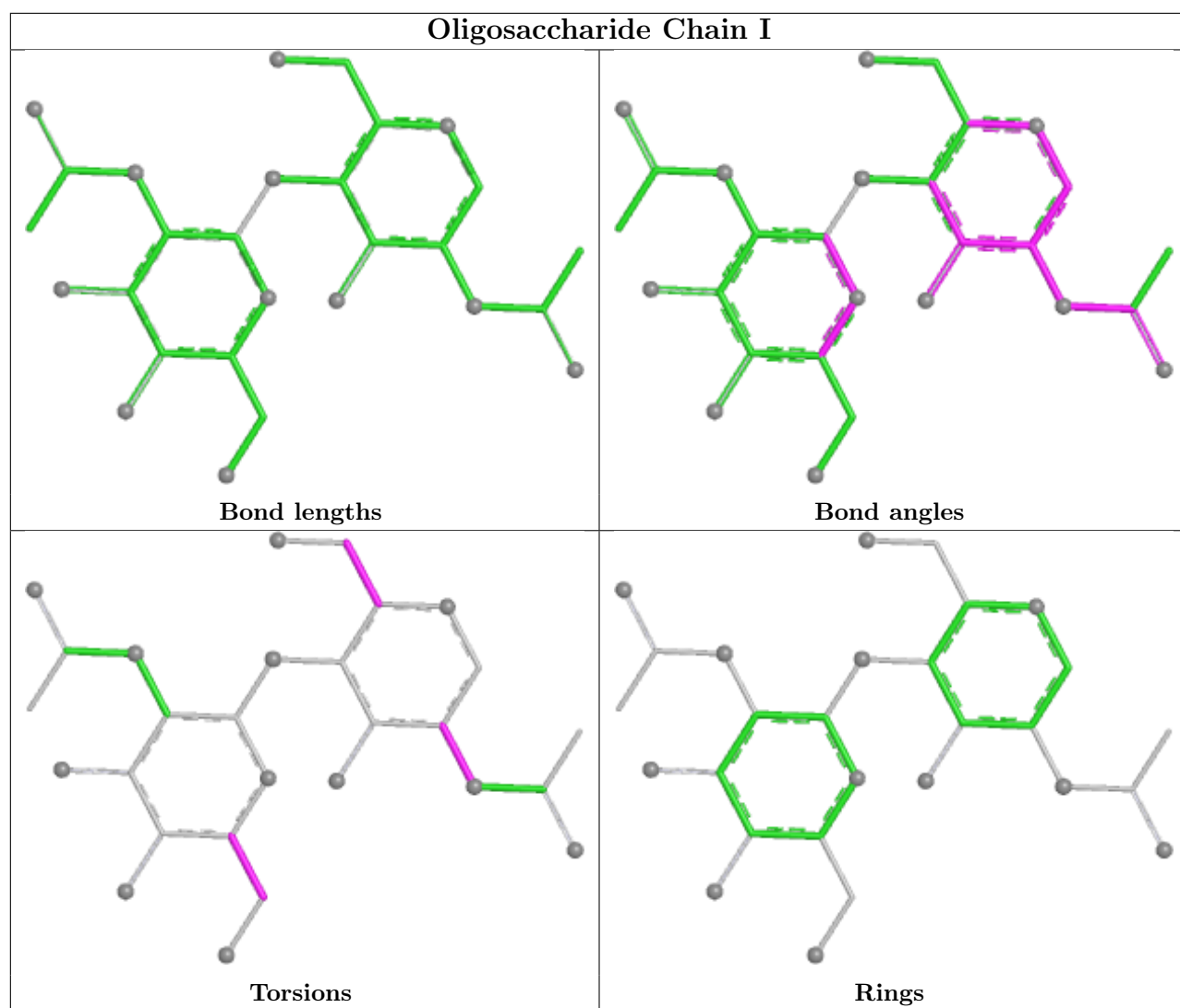
Oligosaccharide Chain G

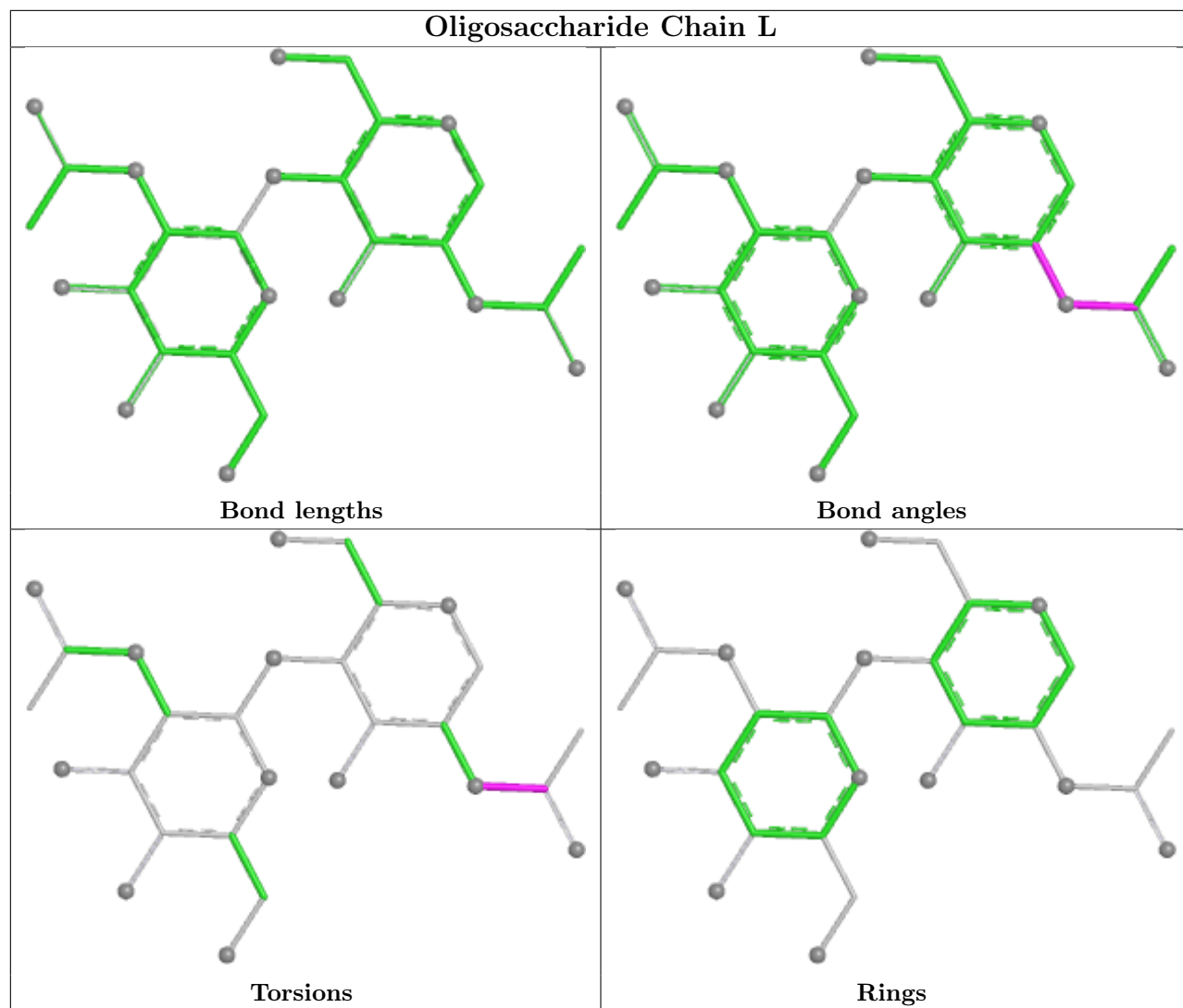


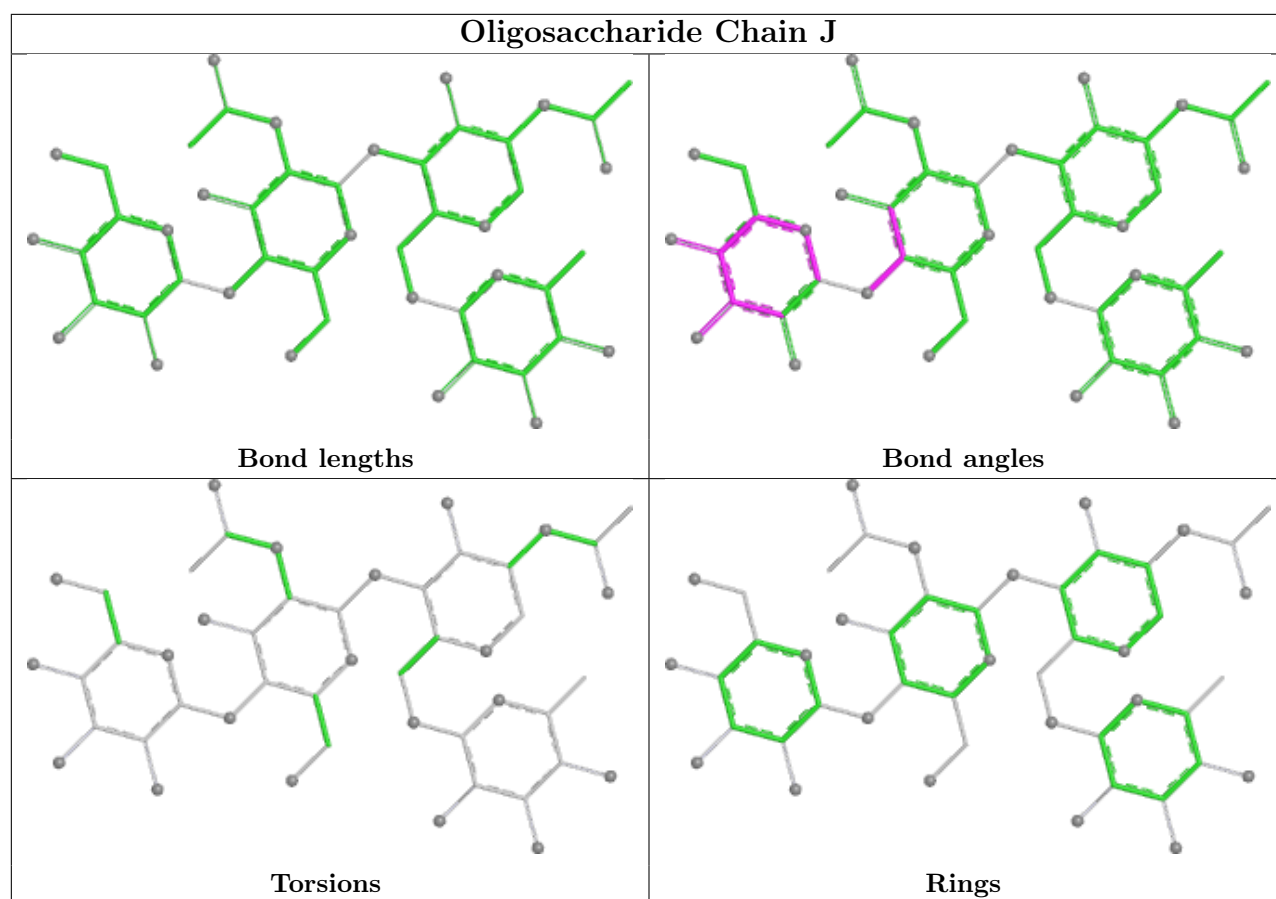
Oligosaccharide Chain F

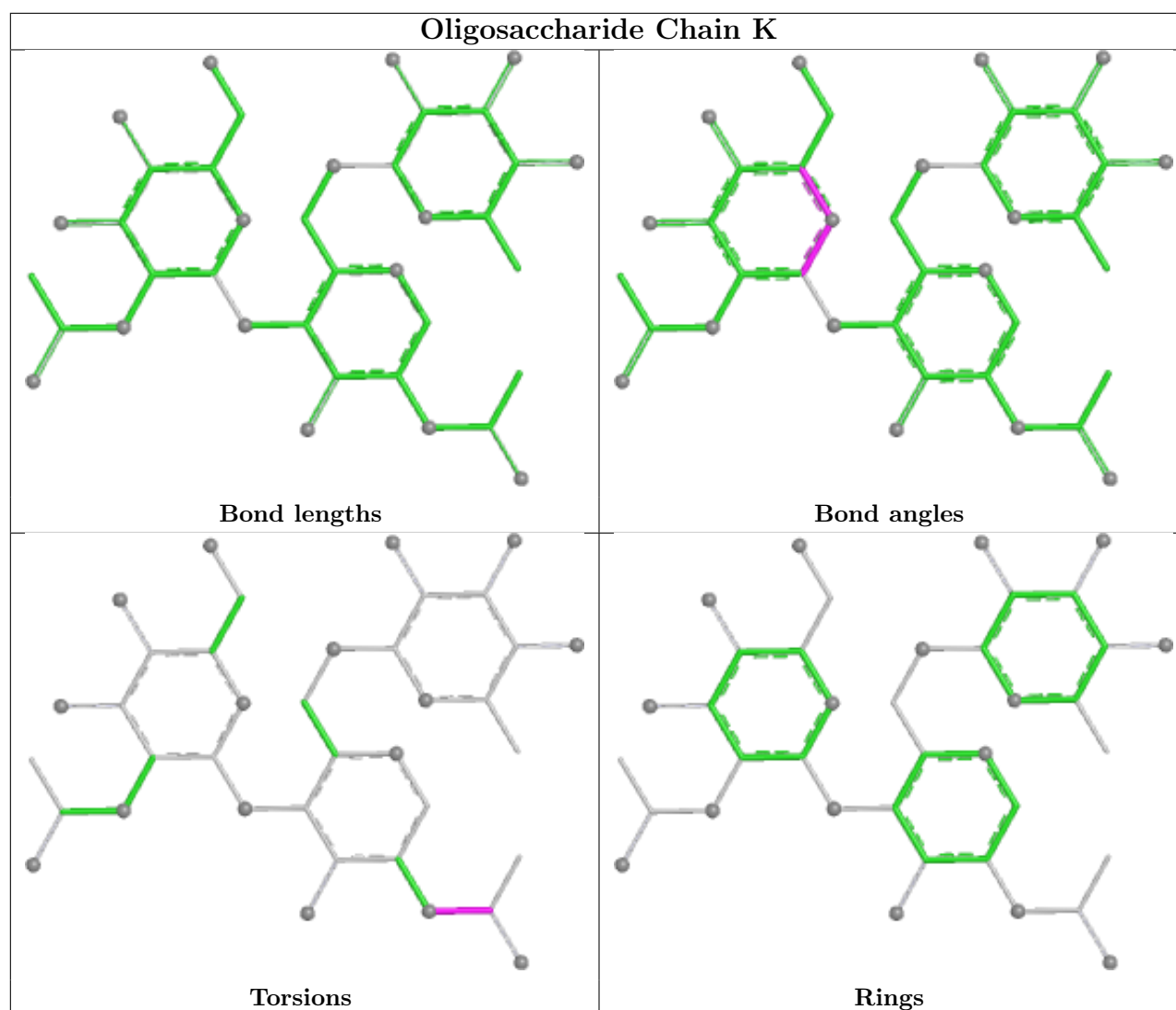












5.6 Ligand geometry [i](#)

28 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	$\# Z > 2$
7	NAG	B	602	1	14,14,15	0.74	0	17,19,21	1.92	3 (17%)
12	EDO	D	602	-	3,3,3	0.24	0	2,2,2	0.43	0
9	GB	A	607	1	0,3,4	-	-	0,3,6	-	-

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
8	GOL	B	605	-	5,5,5	0.38	0	5,5,5	0.47	0
7	NAG	B	601	1	14,14,15	0.70	0	17,19,21	1.13	1 (5%)
8	GOL	A	604	-	5,5,5	0.34	0	5,5,5	0.37	0
8	GOL	C	603	-	5,5,5	0.35	0	5,5,5	0.29	0
7	NAG	A	603	1	14,14,15	0.69	0	17,19,21	1.34	2 (11%)
10	SO4	A	609	-	4,4,4	0.69	0	6,6,6	0.12	0
8	GOL	A	608	-	5,5,5	0.34	0	5,5,5	0.42	0
7	NAG	A	602	1	14,14,15	0.72	0	17,19,21	1.26	2 (11%)
7	NAG	A	601	1	14,14,15	0.68	0	17,19,21	2.48	4 (23%)
10	SO4	A	610	-	4,4,4	0.69	0	6,6,6	0.07	0
12	EDO	B	608	-	3,3,3	0.26	0	2,2,2	0.31	0
10	SO4	B	606	-	4,4,4	0.67	0	6,6,6	0.09	0
10	SO4	C	605	-	4,4,4	0.68	0	6,6,6	0.09	0
10	SO4	A	611	-	4,4,4	0.68	0	6,6,6	0.11	0
9	GB	B	604	1	0,3,4	-	-	0,3,6	-	-
8	GOL	A	605	-	5,5,5	0.33	0	5,5,5	0.52	0
9	GB	D	601	1	0,3,4	-	-	0,3,6	-	-
10	SO4	C	604	-	4,4,4	0.68	0	6,6,6	0.12	0
12	EDO	B	607	-	3,3,3	0.25	0	2,2,2	0.31	0
8	GOL	C	602	-	5,5,5	0.36	0	5,5,5	0.43	0
12	EDO	B	609	-	3,3,3	0.26	0	2,2,2	0.19	0
10	SO4	A	612	-	4,4,4	0.69	0	6,6,6	0.21	0
8	GOL	A	606	-	5,5,5	0.34	0	5,5,5	0.48	0
9	GB	C	601	1	0,3,4	-	-	0,3,6	-	-
11	PEG	B	603	-	6,6,6	0.25	0	5,5,5	0.24	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
7	NAG	A	603	1	-	3/6/23/26	0/1/1/1
12	EDO	B	607	-	-	0/1/1/1	-
8	GOL	C	602	-	-	2/4/4/4	-
7	NAG	B	602	1	-	3/6/23/26	0/1/1/1
12	EDO	B	608	-	-	0/1/1/1	-
12	EDO	B	609	-	-	1/1/1/1	-
8	GOL	A	608	-	-	2/4/4/4	-
7	NAG	A	602	1	-	4/6/23/26	0/1/1/1
12	EDO	D	602	-	-	1/1/1/1	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
8	GOL	A	606	-	-	0/4/4/4	-
8	GOL	B	605	-	-	2/4/4/4	-
7	NAG	B	601	1	-	3/6/23/26	0/1/1/1
7	NAG	A	601	1	-	3/6/23/26	0/1/1/1
11	PEG	B	603	-	-	3/4/4/4	-
8	GOL	A	604	-	-	2/4/4/4	-
8	GOL	A	605	-	-	2/4/4/4	-
8	GOL	C	603	-	-	2/4/4/4	-

There are no bond length outliers.

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	A	601	NAG	C2-N2-C7	8.55	134.35	122.90
7	B	602	NAG	C2-N2-C7	5.92	130.83	122.90
7	A	603	NAG	C2-N2-C7	3.64	127.77	122.90
7	A	602	NAG	C1-O5-C5	3.34	116.66	112.19
7	B	602	NAG	C1-C2-N2	3.29	115.62	110.43
7	B	601	NAG	C1-O5-C5	3.08	116.32	112.19
7	A	601	NAG	C8-C7-N2	2.65	120.51	116.12
7	A	601	NAG	C1-O5-C5	2.49	115.53	112.19
7	A	602	NAG	C2-N2-C7	2.35	126.05	122.90
7	B	602	NAG	O7-C7-N2	2.32	126.08	121.98
7	A	601	NAG	O5-C1-C2	-2.17	107.93	111.29
7	A	603	NAG	O5-C1-C2	-2.06	108.11	111.29

There are no chirality outliers.

All (33) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
7	B	602	NAG	C1-C2-N2-C7
8	A	605	GOL	C1-C2-C3-O3
8	B	605	GOL	C1-C2-C3-O3
7	A	602	NAG	C4-C5-C6-O6
7	A	602	NAG	O5-C5-C6-O6
7	A	601	NAG	C8-C7-N2-C2
7	A	601	NAG	O7-C7-N2-C2
7	A	602	NAG	C8-C7-N2-C2
7	A	602	NAG	O7-C7-N2-C2

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Mol	Chain	Res	Type	Atoms
7	B	601	NAG	C8-C7-N2-C2
7	B	601	NAG	O7-C7-N2-C2
7	B	602	NAG	C4-C5-C6-O6
8	A	604	GOL	C1-C2-C3-O3
8	C	602	GOL	C1-C2-C3-O3
8	C	603	GOL	C1-C2-C3-O3
8	A	605	GOL	O2-C2-C3-O3
7	B	602	NAG	O5-C5-C6-O6
12	B	609	EDO	O1-C1-C2-O2
7	A	603	NAG	O5-C5-C6-O6
7	B	601	NAG	O5-C5-C6-O6
8	B	605	GOL	O2-C2-C3-O3
8	A	608	GOL	C1-C2-C3-O3
12	D	602	EDO	O1-C1-C2-O2
11	B	603	PEG	O1-C1-C2-O2
7	A	601	NAG	C3-C2-N2-C7
7	A	603	NAG	C3-C2-N2-C7
11	B	603	PEG	O2-C3-C4-O4
8	A	604	GOL	O2-C2-C3-O3
7	A	603	NAG	C1-C2-N2-C7
8	A	608	GOL	O2-C2-C3-O3
8	C	603	GOL	O2-C2-C3-O3
8	C	602	GOL	O2-C2-C3-O3
11	B	603	PEG	C1-C2-O2-C3

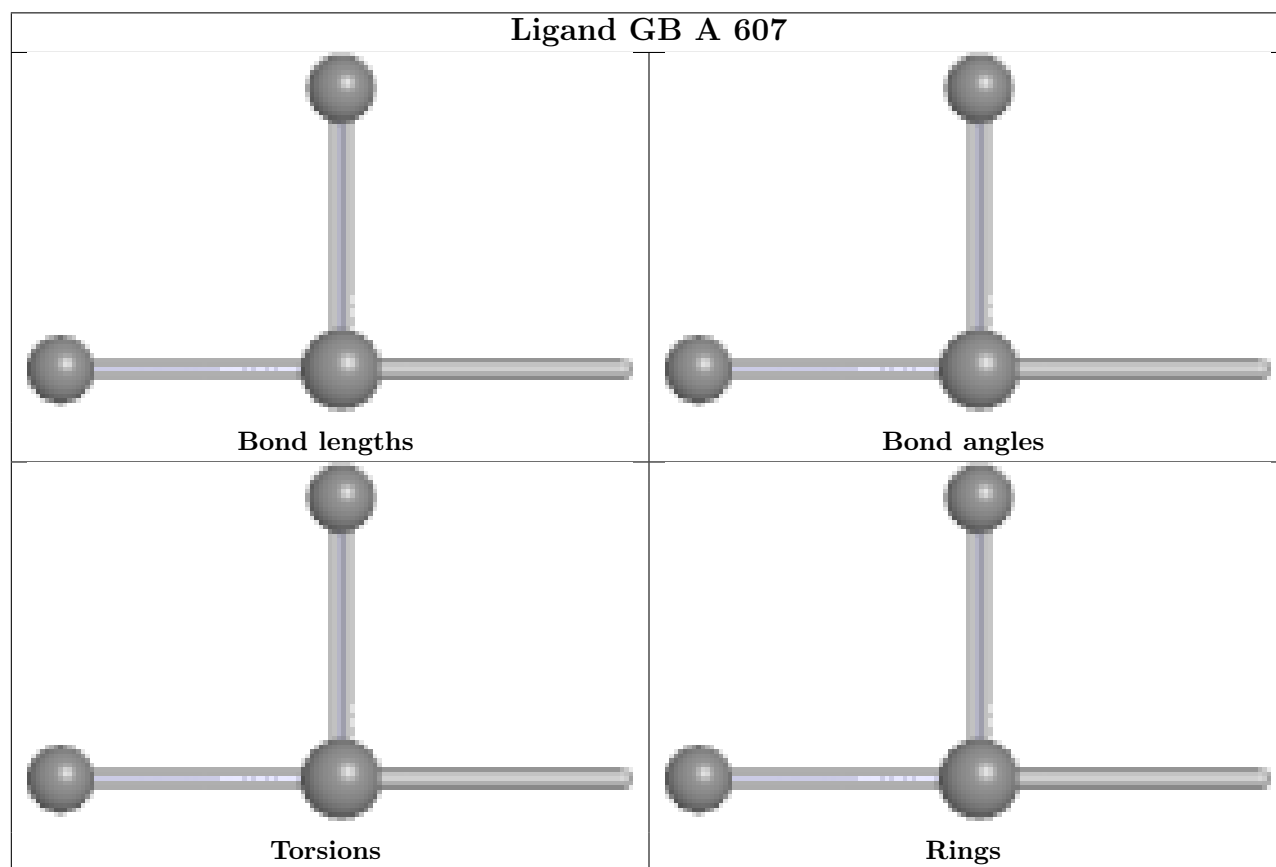
There are no ring outliers.

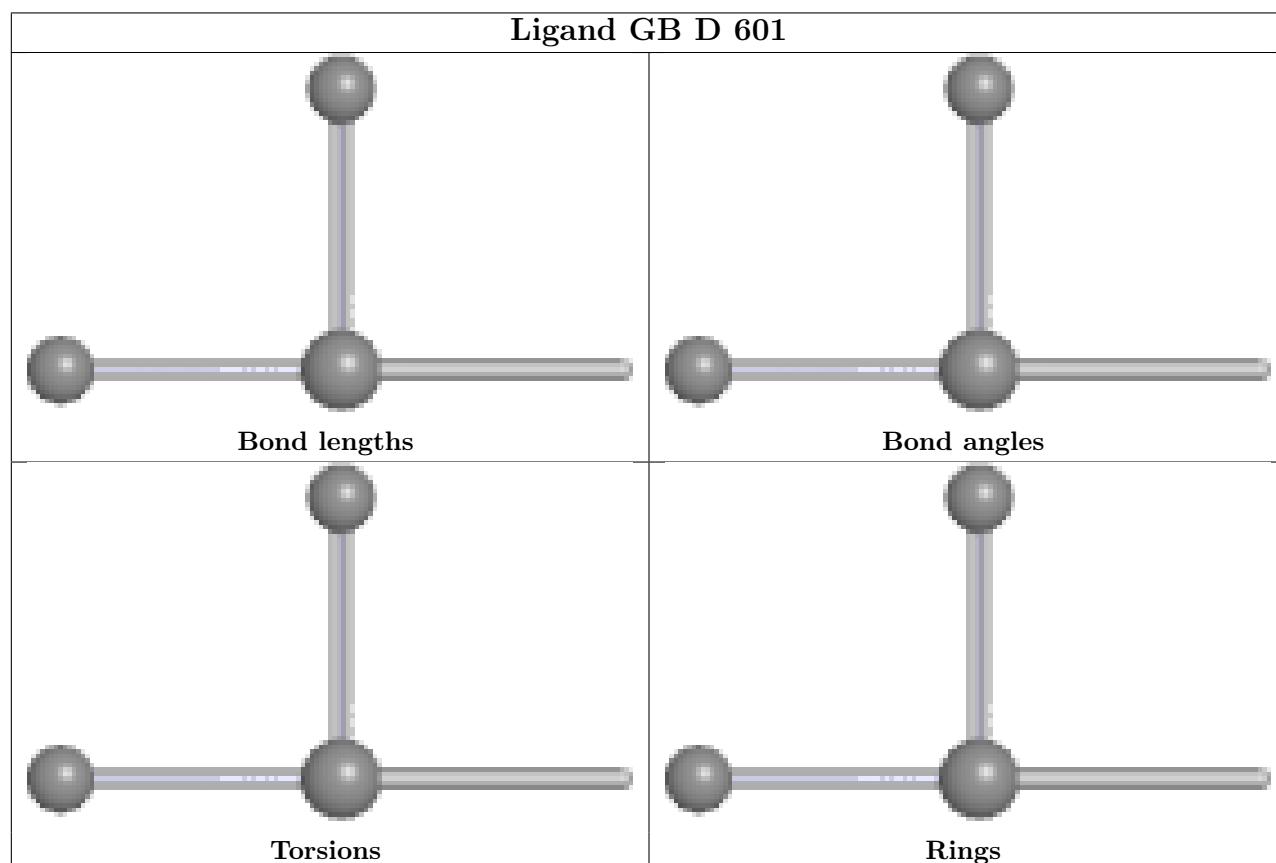
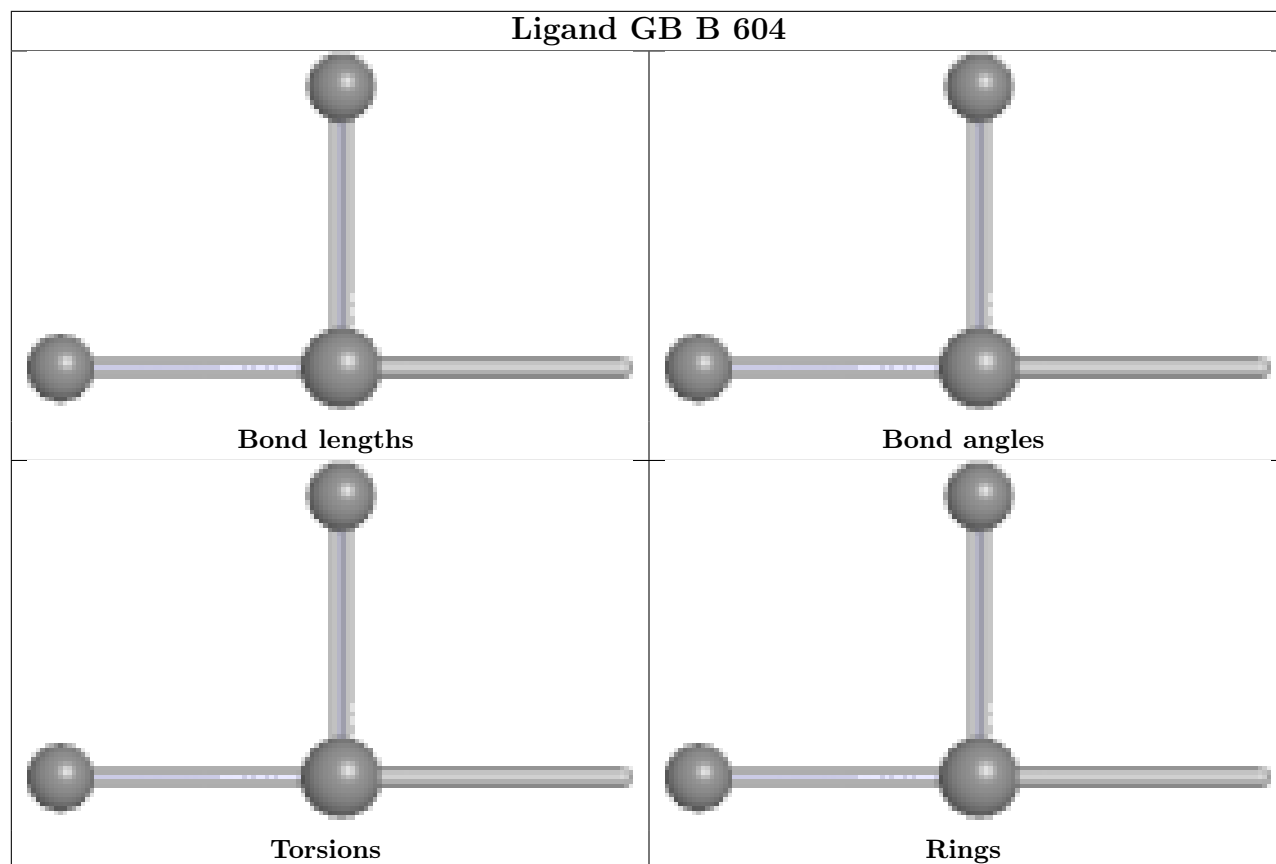
2 monomers are involved in 4 short contacts:

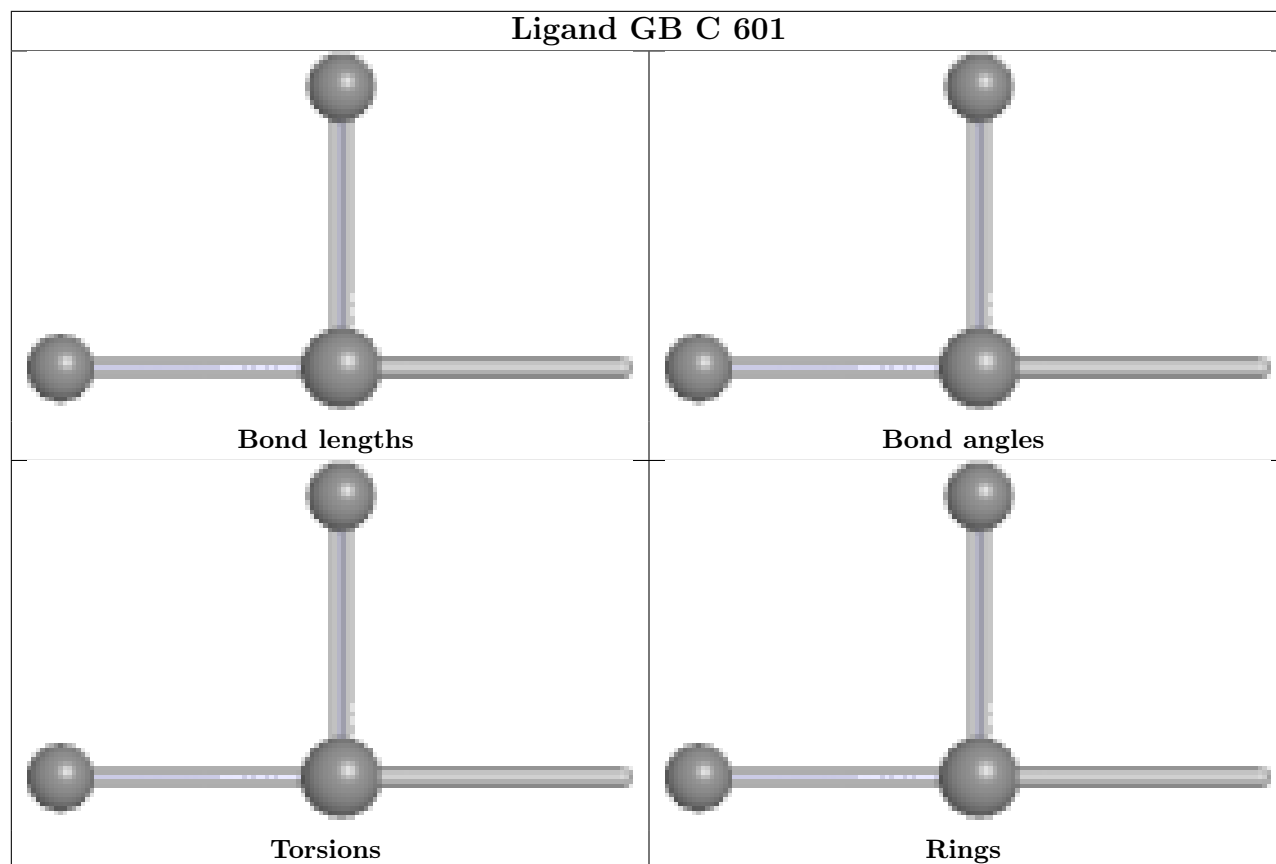
Mol	Chain	Res	Type	Clashes	Symm-Clashes
8	A	604	GOL	1	0
11	B	603	PEG	3	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier.

The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.







5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 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	523/529 (98%)	-0.10	10 (1%) 66 61	20, 39, 76, 120	1 (0%)
1	B	522/529 (98%)	-0.14	8 (1%) 72 67	24, 39, 69, 138	0
1	C	521/529 (98%)	0.28	21 (4%) 42 34	24, 48, 94, 163	1 (0%)
1	D	508/529 (96%)	0.70	42 (8%) 17 12	32, 64, 107, 151	0
All	All	2074/2116 (98%)	0.18	81 (3%) 43 35	20, 46, 96, 163	2 (0%)

All (81) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	63	ASN	5.2
1	D	21	LEU	4.6
1	C	377	THR	4.4
1	D	58	ALA	4.4
1	D	477	TYR	4.3
1	B	282	ASN	4.0
1	D	13	VAL	3.9
1	D	41	LEU	3.8
1	C	4	ILE	3.7
1	D	61	TYR	3.7
1	C	54	ASP	3.6
1	B	360	GLY	3.6
1	D	59	THR	3.5
1	D	62	ALA	3.5
1	D	53	PRO	3.5
1	A	376	TRP	3.5
1	C	333	GLY	3.4
1	C	53	PRO	3.3
1	D	285	LEU	3.1
1	D	55	ILE	3.1
1	B	285	LEU	3.0

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Mol	Chain	Res	Type	RSRZ
1	A	4	ILE	3.0
1	D	277	VAL	3.0
1	D	66	TYR	2.9
1	D	287	SER	2.9
1	D	10	ASN	2.9
1	C	280	VAL	2.9
1	D	11	GLY	2.9
1	C	55	ILE	2.9
1	D	52	TRP	2.7
1	D	275	ASN	2.7
1	D	30	GLY	2.7
1	C	376	TRP	2.7
1	C	12	LYS	2.7
1	A	281	PRO	2.7
1	C	13	VAL	2.6
1	D	50	THR	2.6
1	A	378	ASP	2.5
1	C	281	PRO	2.5
1	D	272	ILE	2.5
1	B	284	MET	2.5
1	D	60	LYS	2.4
1	A	284	MET	2.4
1	C	378	ASP	2.4
1	A	361	VAL	2.4
1	D	94	TYR	2.3
1	B	50	THR	2.3
1	D	458	LYS	2.3
1	A	258	THR	2.3
1	C	287	SER	2.3
1	D	64	SER	2.3
1	D	487	SER	2.3
1	D	177	TRP	2.3
1	D	18	LEU	2.3
1	A	53	PRO	2.3
1	D	47	GLN	2.3
1	D	20	VAL	2.3
1	D	14	ARG	2.2
1	C	5	ILE	2.2
1	C	58	ALA	2.2
1	D	488	THR	2.2
1	B	280	VAL	2.2
1	D	12	LYS	2.2

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Mol	Chain	Res	Type	RSRZ
1	C	57	ASN	2.2
1	B	53	PRO	2.2
1	D	16	MET	2.2
1	C	59	THR	2.2
1	D	37	PRO	2.2
1	A	287	SER	2.2
1	D	15	GLY	2.1
1	A	289	ASN	2.1
1	B	489	ARG	2.1
1	C	7	THR	2.1
1	C	106	ASN	2.1
1	C	28	PHE	2.1
1	D	48	SER	2.0
1	D	105	LYS	2.0
1	C	56	TRP	2.0
1	D	132	PHE	2.0
1	D	469	LYS	2.0
1	D	71	GLN	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
5	BMA	J	3	11/12	0.19	0.21	108,112,118,121	0
6	FUC	K	3	10/11	0.23	0.21	124,137,142,145	0
6	NAG	K	2	14/15	0.34	0.18	110,117,123,128	0
4	NAG	L	2	14/15	0.61	0.18	109,118,128,128	0
6	NAG	K	1	14/15	0.61	0.15	97,107,122,128	0
3	BMA	F	3	11/12	0.63	0.18	75,82,86,91	0
2	NAG	E	1	14/15	0.68	0.14	103,111,117,122	0
2	FUC	E	2	10/11	0.70	0.16	109,127,133,134	0
4	NAG	I	1	14/15	0.70	0.16	74,79,85,88	0
5	FUC	J	4	10/11	0.70	0.19	96,112,114,116	0

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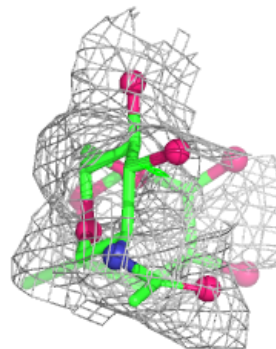
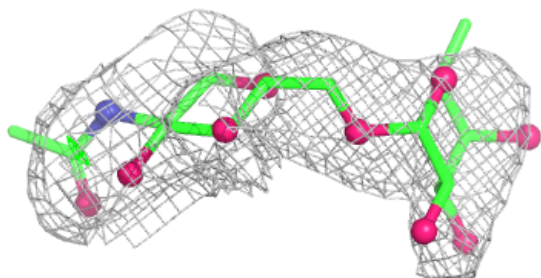
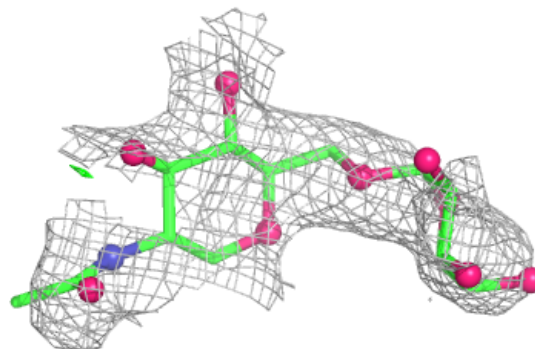
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
4	NAG	H	2	14/15	0.71	0.15	74,94,105,108	0
3	FUC	F	5	10/11	0.71	0.17	89,98,101,104	0
4	NAG	I	2	14/15	0.75	0.14	69,81,89,89	0
3	NAG	F	2	14/15	0.81	0.14	57,68,76,81	0
2	NAG	G	1	14/15	0.81	0.12	73,84,91,98	0
3	MAN	F	4	11/12	0.82	0.13	68,79,83,84	0
5	NAG	J	2	14/15	0.84	0.12	55,84,91,103	0
2	FUC	G	2	10/11	0.84	0.17	98,111,114,116	0
4	NAG	L	1	14/15	0.85	0.12	67,75,91,100	0
4	NAG	H	1	14/15	0.87	0.10	49,57,69,76	0
3	NAG	F	1	14/15	0.88	0.10	42,53,68,84	0
5	NAG	J	1	14/15	0.90	0.11	52,63,80,94	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.

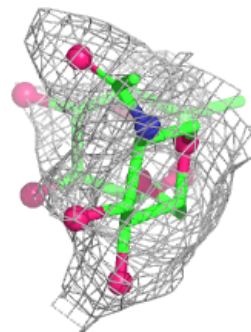
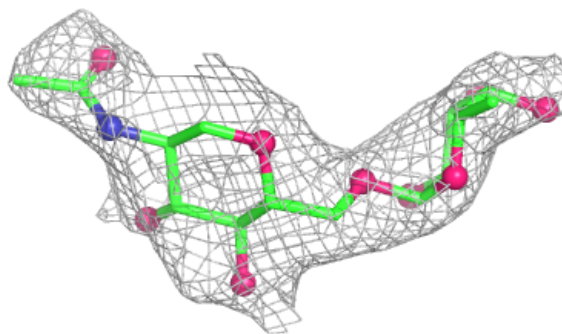
Electron density around Chain E:

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

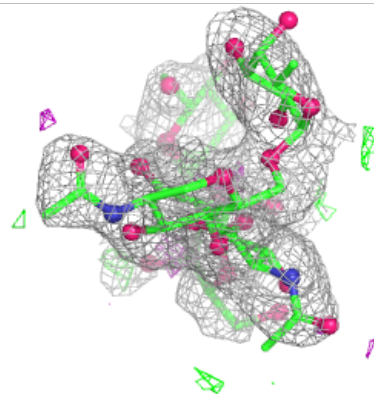
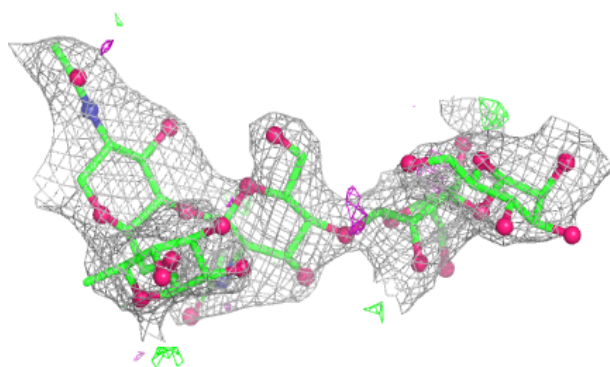
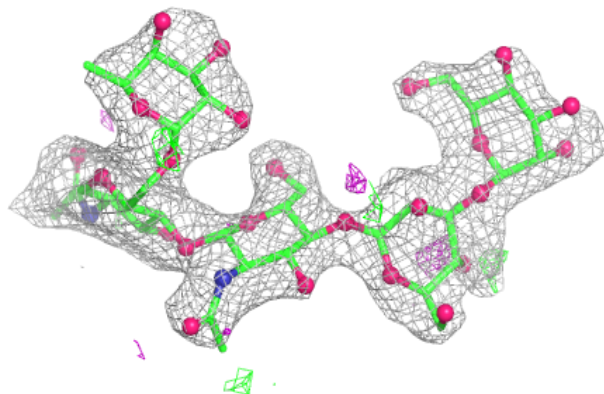


Electron density around Chain G:

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

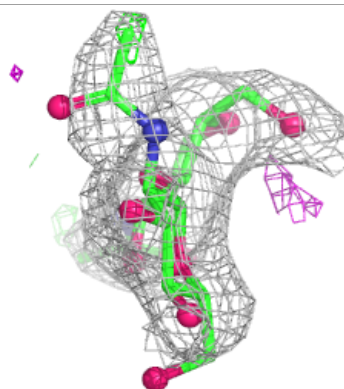
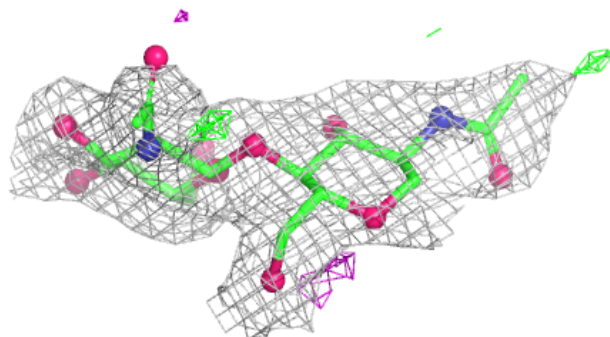
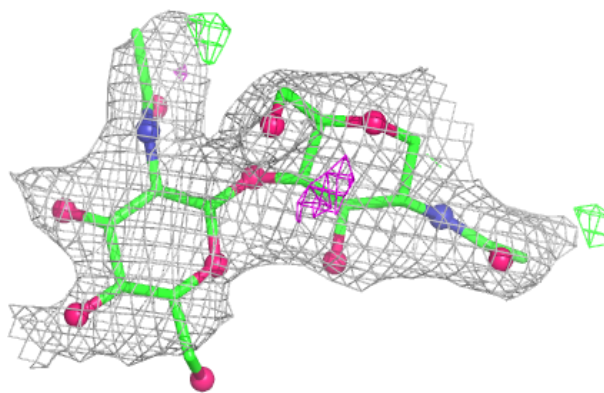
**Electron density around Chain F:**

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



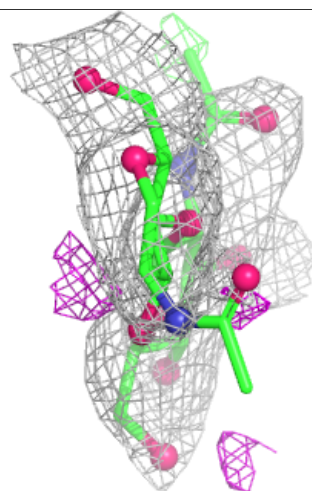
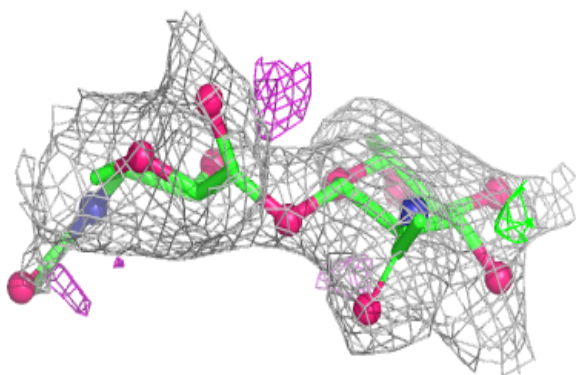
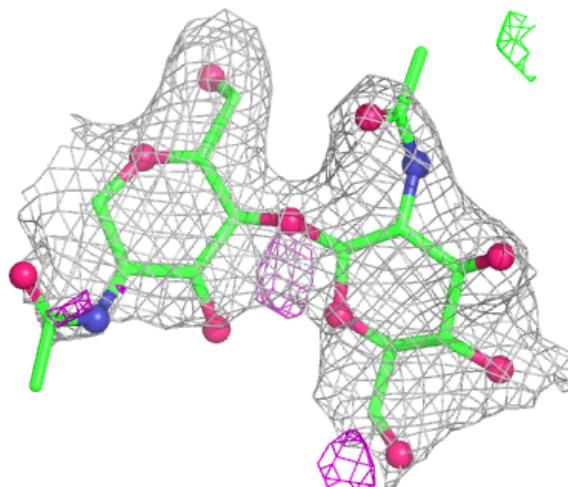
Electron density around Chain H:

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



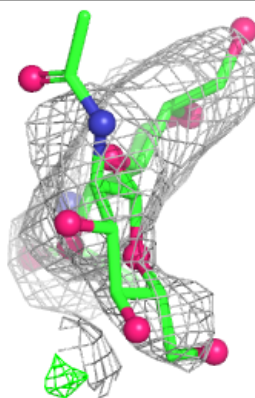
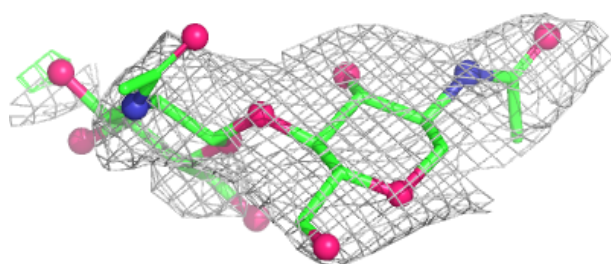
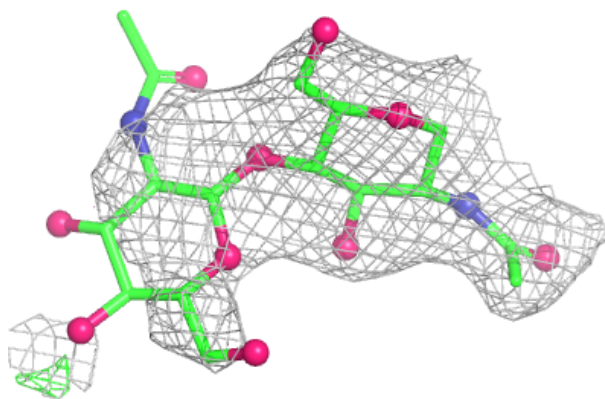
Electron density around Chain I:

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

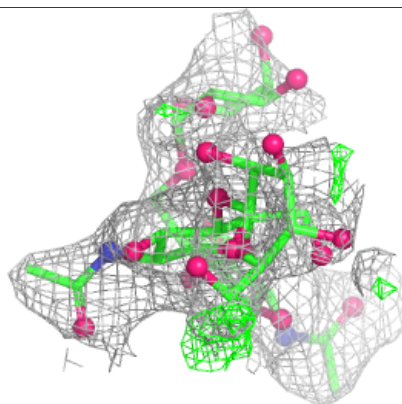
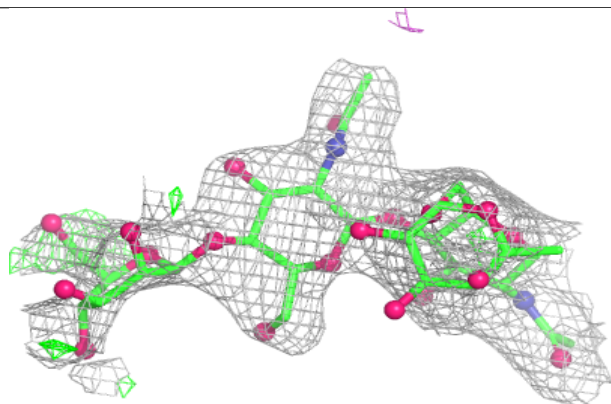
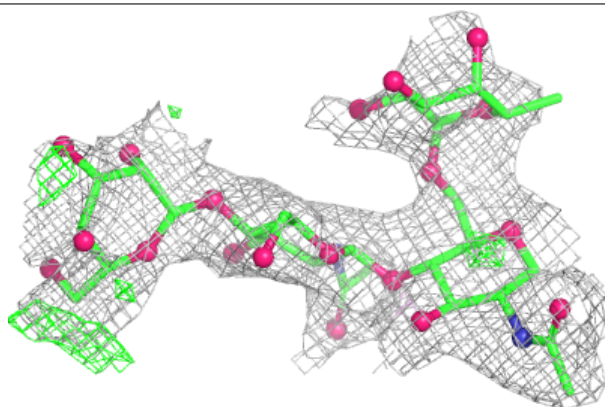


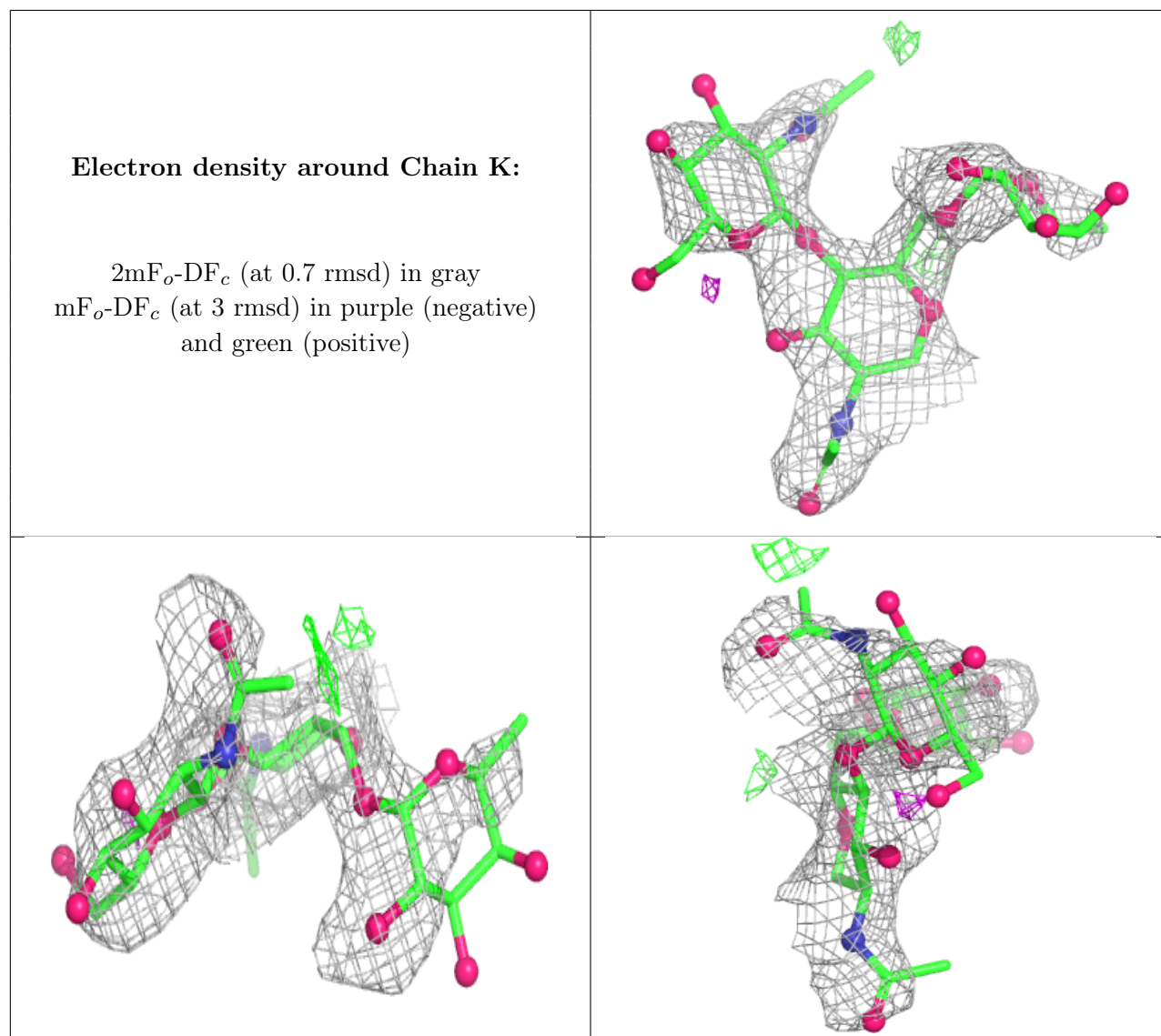
Electron density around Chain L:

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

**Electron density around Chain J:**

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





6.4 Ligands [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
7	NAG	A	603	14/15	0.44	0.18	97,113,121,121	0
7	NAG	B	601	14/15	0.57	0.24	91,104,118,122	0
7	NAG	B	602	14/15	0.70	0.15	112,119,127,127	0
7	NAG	A	602	14/15	0.71	0.16	75,89,95,97	0
10	SO4	B	606	5/5	0.72	0.15	81,84,96,96	0
7	NAG	A	601	14/15	0.73	0.14	66,69,74,86	0
10	SO4	A	611	5/5	0.82	0.13	84,87,101,106	0

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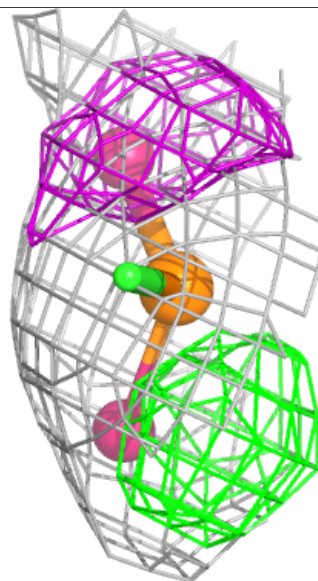
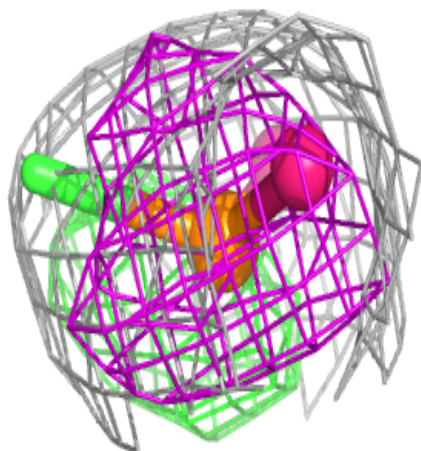
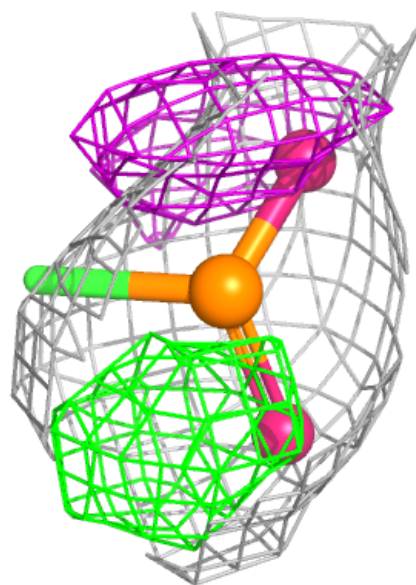
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
12	EDO	D	602	4/4	0.83	0.21	54,55,56,60	0
8	GOL	A	605	6/6	0.85	0.15	44,51,55,56	0
10	SO4	C	605	5/5	0.88	0.16	72,73,85,90	0
12	EDO	B	608	4/4	0.88	0.11	43,44,47,47	0
10	SO4	A	612	5/5	0.88	0.20	61,69,76,82	0
12	EDO	B	607	4/4	0.89	0.12	41,44,47,48	0
10	SO4	A	609	5/5	0.90	0.11	52,56,59,67	0
8	GOL	A	606	6/6	0.90	0.15	62,64,68,69	0
8	GOL	C	603	6/6	0.90	0.14	50,54,57,58	0
9	GB	D	601	4/5	0.90	0.20	60,61,67,73	0
8	GOL	A	608	6/6	0.91	0.13	46,48,49,60	0
8	GOL	B	605	6/6	0.91	0.14	52,52,57,58	0
12	EDO	B	609	4/4	0.92	0.12	36,42,42,43	0
11	PEG	B	603	7/7	0.92	0.13	46,46,49,52	0
10	SO4	A	610	5/5	0.93	0.15	54,55,64,69	0
10	SO4	C	604	5/5	0.93	0.11	64,64,69,77	0
8	GOL	C	602	6/6	0.93	0.14	52,56,57,60	0
8	GOL	A	604	6/6	0.93	0.11	36,44,45,46	0
9	GB	C	601	4/5	0.98	0.07	21,23,28,34	0
9	GB	A	607	4/5	0.99	0.07	13,20,21,23	0
9	GB	B	604	4/5	0.99	0.05	20,25,25,28	0

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

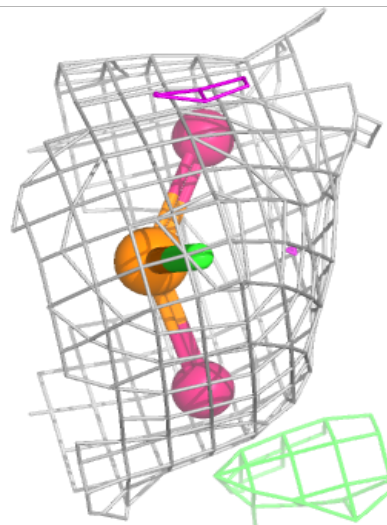
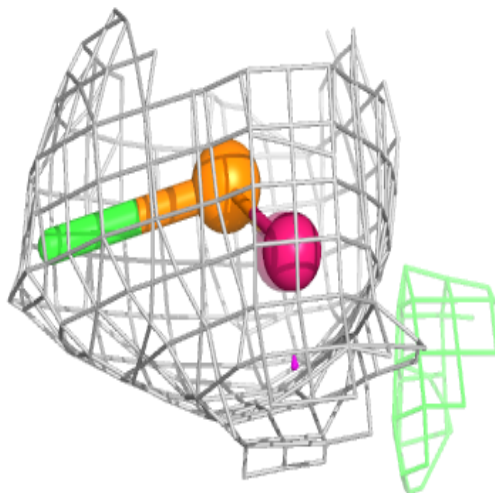
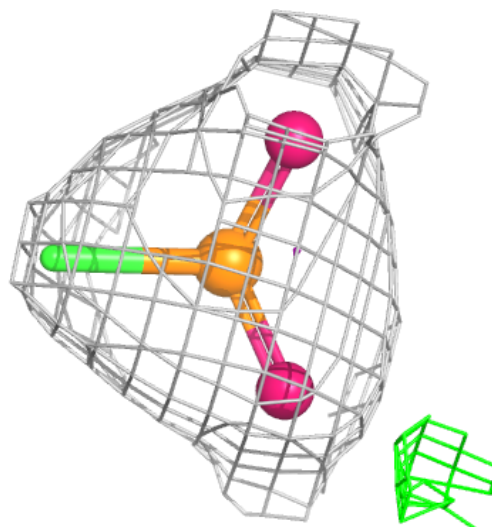
Electron density around GB D 601:

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



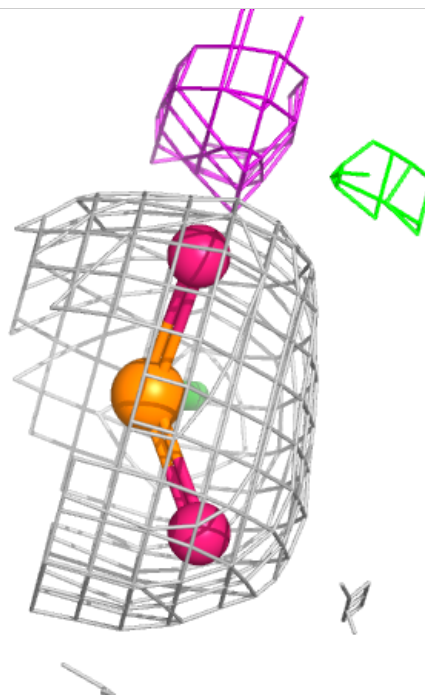
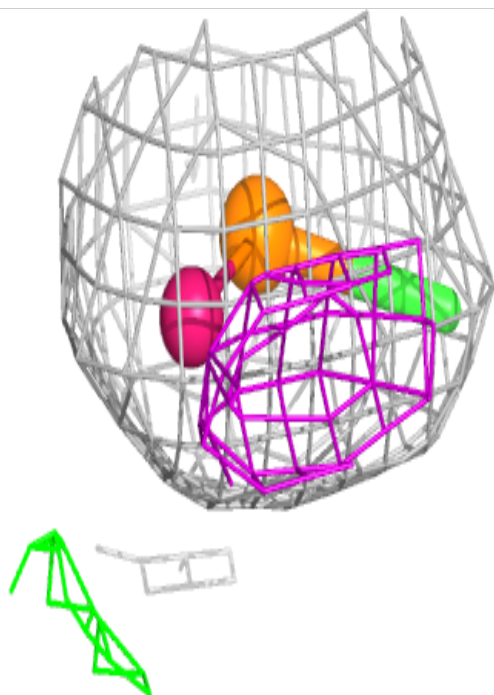
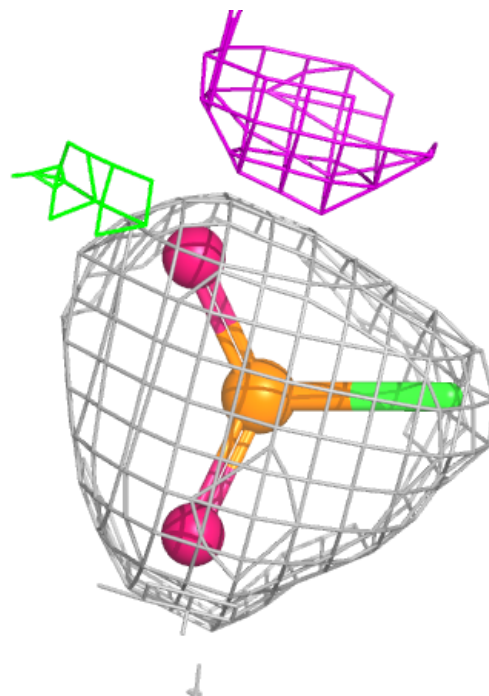
Electron density around GB C 601:

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



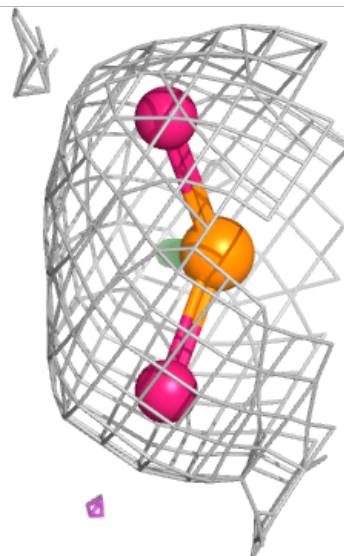
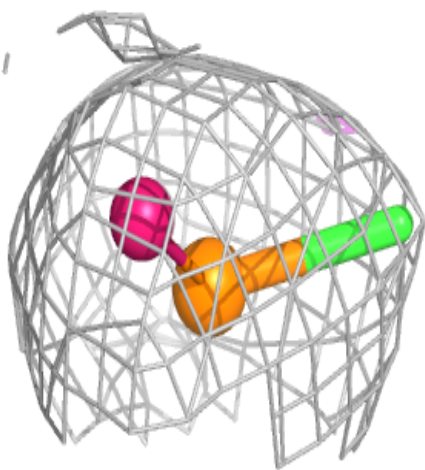
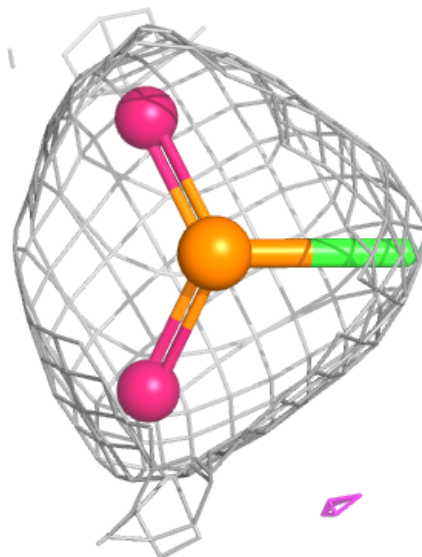
Electron density around GB A 607:

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



Electron density around GB 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)



6.5 Other polymers [i](#)

There are no such residues in this entry.