



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 05:11 PM UTC

PDB ID : 8TIC / pdb_00008tic
Title : Structure of human beta 1,3-N-acetylglucosaminyltransferase 2 with compound 1
Authors : Sudom, A.; Min, X.
Deposited on : 2023-07-19
Resolution : 2.70 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

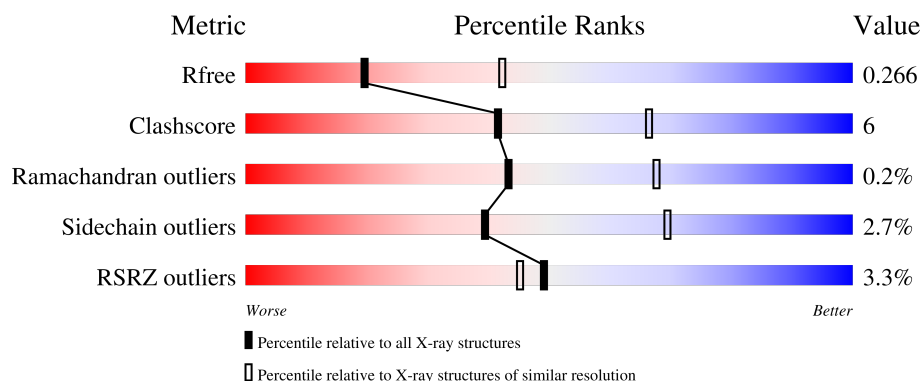
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.70 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	3538 (2.70-2.70)
Clashscore	190562	3843 (2.70-2.70)
Ramachandran outliers	187476	3778 (2.70-2.70)
Sidechain outliers	187428	3778 (2.70-2.70)
RSRZ outliers	180081	3538 (2.70-2.70)

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	397	
1	B	397	
1	C	397	
1	D	397	
2	E	5	

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Mol	Chain	Length	Quality of chain
2	F	5	
2	G	5	
2	H	5	

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

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
4	FKX	A	402	-	X	-	-

2 Entry composition [i](#)

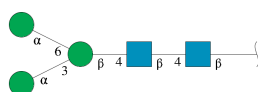
There are 7 unique types of molecules in this entry. The entry contains 11191 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 N-acetylglucosaminide beta-1,3-N-acetylglucosaminyltransferase 2.

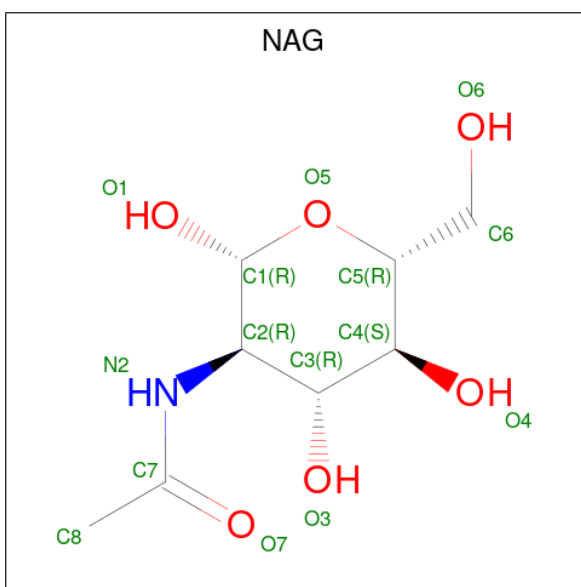
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	328	Total	C	N	O	S	0	0	0
			2710	1748	464	486	12			
1	B	331	Total	C	N	O	S	0	0	0
			2724	1753	467	491	13			
1	C	327	Total	C	N	O	S	0	0	0
			2693	1735	461	484	13			
1	D	320	Total	C	N	O	S	0	0	0
			2639	1698	456	473	12			

- Molecule 2 is an oligosaccharide called 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.



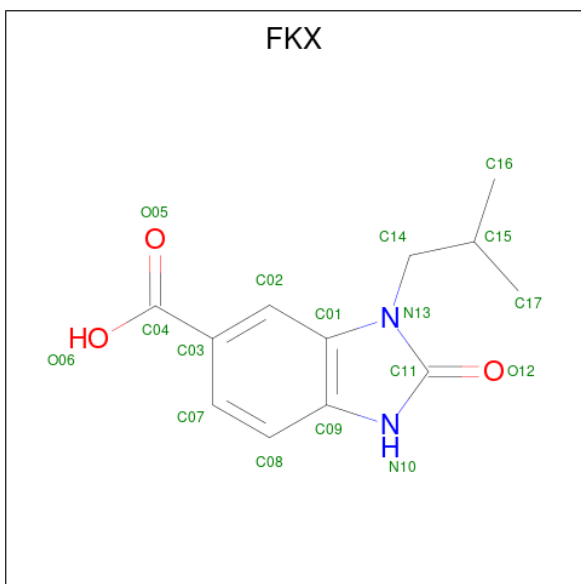
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	E	5	Total	C	N	O	0	0	0
			61	34	2	25			
2	F	5	Total	C	N	O	0	0	0
			61	34	2	25			
2	G	5	Total	C	N	O	0	0	0
			61	34	2	25			
2	H	5	Total	C	N	O	0	0	0
			61	34	2	25			

- Molecule 3 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: C₈H₁₅NO₆).



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

- Molecule 4 is 3-(2-methylpropyl)-2-oxo-2,3-dihydro-1H-benzimidazole-5-carboxylic acid (CCD ID: FKX) (formula: $C_{12}H_{14}N_2O_3$) (labeled as "Ligand of Interest" by depositor).



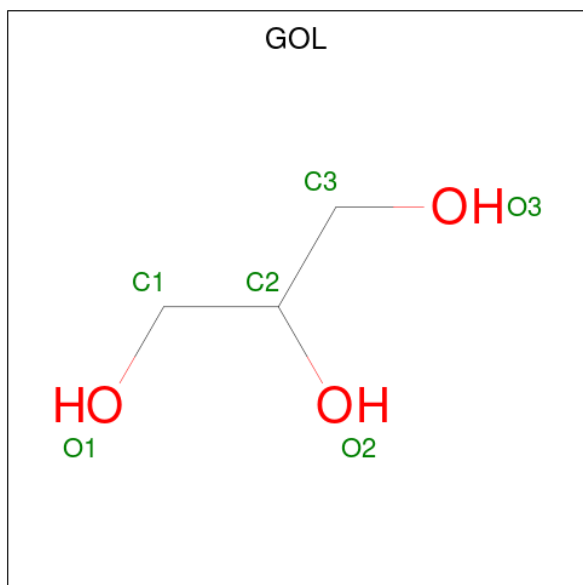
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	N	O	0	0
			17	12	2	3		

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	B	1	Total	C	N	O	0	0
			17	12	2	3		
4	C	1	Total	C	N	O	0	0
			17	12	2	3		
4	D	1	Total	C	N	O	0	0
			17	12	2	3		

- Molecule 5 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



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

- Molecule 6 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

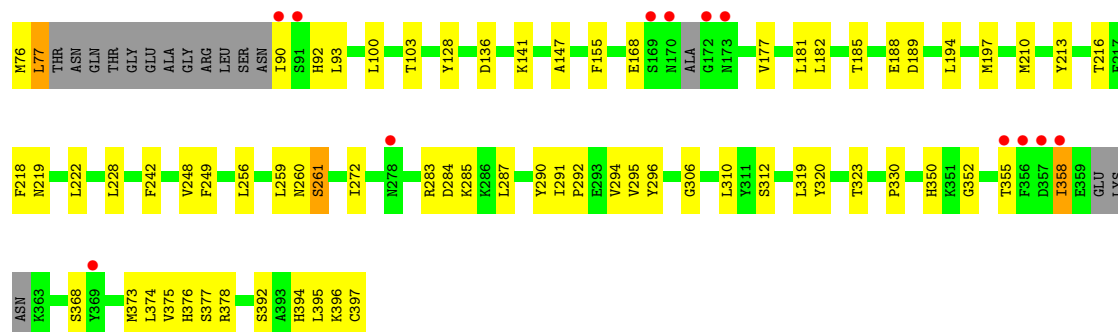
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	C	1	Total	Cl	0	0
			1	1		
6	D	1	Total	Cl	0	0
			1	1		

- Molecule 7 is water.

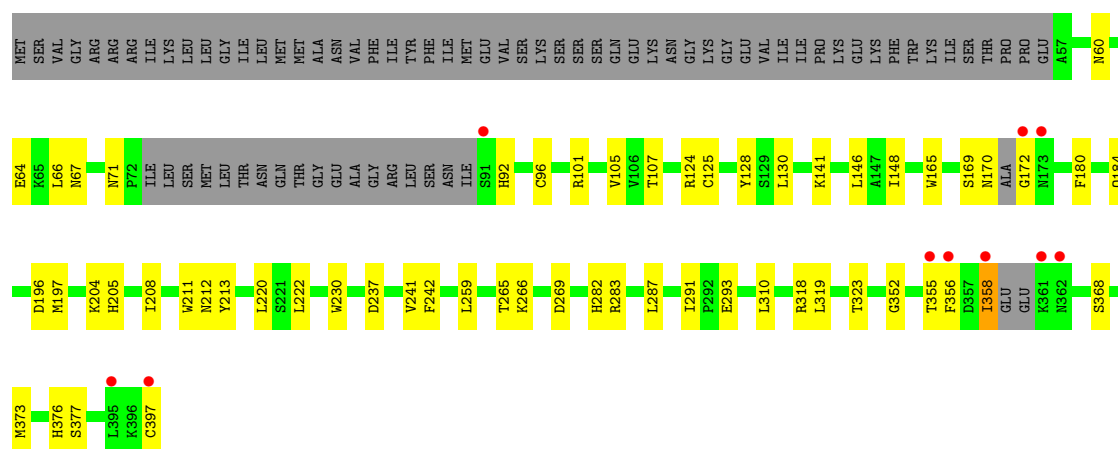
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	14	Total 14	O 14	0	0
7	B	20	Total 20	O 20	0	0
7	C	14	Total 14	O 14	0	0
7	D	11	Total 11	O 11	0	0

- Molecule 1: N-acetyllactosaminide beta-1,3-N-acetylglucosaminyltransferase 2





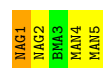
- Molecule 1: N-acetylactosaminide beta-1,3-N-acetylglucosaminyltransferase 2



- Molecule 2: 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



- Molecule 2: 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



- Molecule 2: 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



MAG1	MAG2	BMA3	MAN4	MAN5
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● Molecule 2: α -D-mannopyranose-(1-3)-[α -D-mannopyranose-(1-6)] β -D-mannopyranose-(1-4)-2-acetamido-2-deoxy- β -D-glucopyranose-(1-4)-2-acetamido-2-deoxy- β -D-glucopyranose

Chain H:



MAG1	MAG2	BMA3	MAN4	MAN5
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4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	89.14Å 92.39Å 203.43Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.98 – 2.70 29.98 – 2.70	Depositor EDS
% Data completeness (in resolution range)	100.0 (29.98-2.70) 88.8 (29.98-2.70)	Depositor EDS
R_{merge}	0.39	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	0.84 (at 2.68Å)	Xtriage
Refinement program	PHENIX (1.20.1_4487: ???)	Depositor
R, R_{free}	0.212 , 0.266 0.212 , 0.266	Depositor DCC
R_{free} test set	2004 reflections (4.27%)	wwPDB-VP
Wilson B-factor (Å ²)	35.4	Xtriage
Anisotropy	0.435	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 50.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	0.027 for k,h,-l	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	11191	wwPDB-VP
Average B, all atoms (Å ²)	50.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.23% 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: NAG, BMA, MAN, GOL, CL, FKX

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	0.17	0/2785	0.36	0/3772
1	B	0.18	0/2798	0.37	0/3790
1	C	0.16	0/2765	0.36	0/3743
1	D	0.16	0/2710	0.36	0/3668
All	All	0.17	0/11058	0.36	0/14973

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	2710	0	2651	32	0
1	B	2724	0	2677	36	0
1	C	2693	0	2645	39	0
1	D	2639	0	2586	35	0
2	E	61	0	52	0	0
2	F	61	0	52	6	0
2	G	61	0	52	2	0
2	H	61	0	52	0	0
3	A	14	0	13	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	14	0	13	0	0
4	A	17	0	0	1	0
4	B	17	0	0	0	0
4	C	17	0	0	0	0
4	D	17	0	0	1	0
5	A	6	0	8	0	0
5	B	6	0	8	0	0
5	C	12	0	16	0	0
6	C	1	0	0	0	0
6	D	1	0	0	0	0
7	A	14	0	0	0	0
7	B	20	0	0	0	0
7	C	14	0	0	0	0
7	D	11	0	0	0	0
All	All	11191	0	10825	140	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 6.

All (140) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:219:ASN:HD21	2:F:1:NAG:C1	1.63	1.10
1:B:219:ASN:HD21	2:F:1:NAG:C2	1.76	0.98
1:B:219:ASN:ND2	2:F:1:NAG:C1	2.29	0.95
1:B:366:ILE:HD13	1:B:389:GLN:HB2	1.67	0.74
1:B:196:ASP:OD2	1:D:204:LYS:NZ	2.22	0.71
1:B:71:ASN:O	1:B:75:SER:OG	2.09	0.69
1:A:90:ILE:HG23	1:A:93:LEU:HD12	1.77	0.67
1:A:272:ILE:HG22	1:A:348:GLU:HB3	1.77	0.67
1:C:306:GLY:HA3	1:C:355:THR:HG21	1.76	0.66
1:C:90:ILE:HG23	1:C:93:LEU:HD12	1.79	0.65
1:C:283:ARG:NH1	1:C:291:ILE:O	2.29	0.63
1:C:358:ILE:O	1:C:378:ARG:NH1	2.32	0.62
1:B:219:ASN:ND2	2:F:1:NAG:C7	2.63	0.61
1:D:67:ASN:O	1:D:71:ASN:ND2	2.33	0.61
1:A:320:TYR:O	1:A:323:THR:OG1	2.18	0.60
1:D:130:LEU:HD11	1:D:230:TRP:HD1	1.65	0.60
1:B:170:ASN:HA	1:B:175:THR:HA	1.84	0.60
1:C:352:GLY:HA3	1:C:373:MET:HG2	1.84	0.59
1:C:256:LEU:O	1:C:260:ASN:ND2	2.36	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:248:VAL:HG11	1:A:307:GLY:HA3	1.85	0.58
1:C:100:LEU:O	1:C:103:THR:OG1	2.21	0.57
1:B:221:SER:OG	1:B:329:TYR:O	2.23	0.57
1:A:128:TYR:O	1:A:229:ARG:NH2	2.29	0.56
1:D:184:GLN:NE2	1:D:212:ASN:OD1	2.38	0.56
1:A:92:HIS:HA	1:A:124:ARG:HB3	1.87	0.56
1:C:66:LEU:HD11	1:C:92:HIS:CD2	2.41	0.56
1:A:62:GLU:OE2	1:A:65:LYS:NZ	2.39	0.56
1:B:137:LYS:HD2	1:B:177:VAL:HG13	1.87	0.55
1:A:269:ASP:OD1	1:A:318:ARG:NH2	2.37	0.55
1:C:248:VAL:HG22	1:C:376:HIS:HB2	1.87	0.55
1:C:147:ALA:HB1	1:C:181:LEU:HD11	1.88	0.54
1:C:128:TYR:HE1	2:G:2:NAG:H82	1.72	0.54
1:C:75:SER:O	1:C:77:LEU:N	2.38	0.54
1:D:101:ARG:O	1:D:105:VAL:HG23	2.07	0.54
1:B:219:ASN:ND2	2:F:1:NAG:C2	2.59	0.54
1:A:259:LEU:HD21	1:A:310:LEU:HD21	1.90	0.54
1:A:199:LYS:O	1:A:203:GLU:HG3	2.09	0.53
1:C:259:LEU:HD21	1:C:310:LEU:HD21	1.90	0.53
1:C:392:SER:HA	1:C:394:HIS:CE1	2.44	0.53
1:A:128:TYR:CE2	1:A:222:LEU:HD21	2.43	0.53
1:C:284:ASP:HB3	1:C:287:LEU:HG	1.90	0.53
1:D:60:ASN:O	1:D:64:GLU:HG3	2.09	0.53
1:A:200:PHE:HD2	1:C:197:MET:HE3	1.73	0.53
1:B:128:TYR:CE2	1:B:222:LEU:HD21	2.44	0.53
1:C:283:ARG:NH2	1:C:296:TYR:O	2.41	0.52
1:B:248:VAL:HG21	1:B:307:GLY:HA3	1.91	0.52
1:D:128:TYR:CE2	1:D:222:LEU:HD21	2.44	0.52
1:B:293:GLU:OE1	1:B:293:GLU:N	2.35	0.51
1:A:289:TYR:HD2	1:A:331:ILE:HD12	1.74	0.51
1:A:224:GLU:HG2	1:A:243:LYS:NZ	2.25	0.51
1:C:128:TYR:CE2	1:C:222:LEU:HD21	2.46	0.51
1:C:272:ILE:HD12	1:C:350:HIS:CD2	2.47	0.50
1:A:223:LYS:NZ	4:A:402:FKX:O12	2.42	0.50
1:B:164:SER:O	1:B:167:GLN:HG3	2.12	0.50
1:A:365:ASN:N	1:A:365:ASN:OD1	2.46	0.49
1:B:354:ARG:NH2	1:B:359:GLU:OE1	2.45	0.49
1:B:362:ASN:HB2	1:B:368:SER:OG	2.12	0.49
1:B:284:ASP:HB3	1:B:287:LEU:HG	1.95	0.49
1:D:130:LEU:HD11	1:D:230:TRP:CD1	2.47	0.49
1:A:292:PRO:HG2	1:A:295:VAL:HG23	1.95	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:142:PRO:HA	1:B:175:THR:OG1	2.13	0.48
1:D:169:SER:OG	1:D:170:ASN:N	2.45	0.48
1:B:73:ILE:HA	1:B:76:MET:HE2	1.95	0.48
1:C:261:SER:O	1:C:261:SER:OG	2.27	0.48
1:A:137:LYS:HD2	1:A:177:VAL:HB	1.95	0.48
1:D:319:LEU:O	1:D:323:THR:HG23	2.14	0.48
1:D:197:MET:HE3	1:D:197:MET:HB3	1.73	0.48
1:B:259:LEU:HD21	1:B:310:LEU:HD21	1.96	0.47
1:D:92:HIS:HA	1:D:124:ARG:HB3	1.96	0.47
1:A:130:LEU:HD13	1:A:211:TRP:CE2	2.49	0.47
1:B:219:ASN:HD21	2:F:1:NAG:C7	2.24	0.47
1:B:365:ASN:OD1	1:B:367:CYS:HB2	2.15	0.47
1:C:319:LEU:O	1:C:323:THR:HG23	2.15	0.47
1:C:396:LYS:HE3	1:C:396:LYS:HB2	1.60	0.47
1:D:204:LYS:HG2	1:D:205:HIS:CD2	2.50	0.47
1:C:242:PHE:HZ	1:C:374:LEU:HD12	1.81	0.46
1:D:211:TRP:HB2	1:D:213:TYR:CE1	2.50	0.46
1:D:269:ASP:OD1	1:D:318:ARG:NH1	2.48	0.46
1:D:180:PHE:HB2	1:D:208:ILE:HD13	1.97	0.45
1:B:393:ALA:O	1:B:397:CYS:N	2.32	0.45
1:C:65:LYS:HE2	1:C:65:LYS:HB3	1.60	0.45
1:C:292:PRO:HG2	1:C:295:VAL:HG23	1.97	0.45
1:C:376:HIS:CD2	1:C:377:SER:HB3	2.52	0.45
1:D:66:LEU:HD11	1:D:92:HIS:CE1	2.52	0.45
1:D:220:LEU:HD21	4:D:501:FKX:C11	2.47	0.45
1:D:293:GLU:H	1:D:293:GLU:CD	2.24	0.45
1:A:110:ASN:N	1:A:110:ASN:OD1	2.49	0.45
1:D:283:ARG:NH2	1:D:293:GLU:HA	2.32	0.45
1:A:141:LYS:HG3	1:A:237:ASP:HB3	1.99	0.45
1:B:249:PHE:HB3	1:B:375:VAL:HG12	1.98	0.44
1:D:96:CYS:HA	1:D:125:CYS:HB2	1.99	0.44
1:D:376:HIS:HA	1:D:377:SER:HA	1.73	0.44
1:A:57:ALA:HB1	1:A:295:VAL:C	2.42	0.44
1:A:281:PRO:HB3	1:A:291:ILE:HB	2.00	0.44
1:D:282:HIS:HB3	1:D:287:LEU:HD12	2.00	0.44
1:A:70:TYR:O	1:A:71:ASN:ND2	2.50	0.44
1:B:286:LYS:HA	1:B:286:LYS:HD3	1.70	0.43
1:B:319:LEU:O	1:B:323:THR:HG23	2.18	0.43
1:D:141:LYS:HG3	1:D:237:ASP:O	2.18	0.43
1:B:366:ILE:HG13	1:B:366:ILE:O	2.18	0.43
1:D:146:LEU:HD23	1:D:242:PHE:HB3	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:103:THR:HG22	1:A:109:PHE:CE2	2.54	0.43
1:D:241:VAL:O	1:D:310:LEU:HD12	2.18	0.43
1:B:76:MET:H	1:B:76:MET:HG2	1.59	0.43
1:B:376:HIS:HA	1:B:377:SER:HA	1.63	0.43
1:C:218:PHE:CD2	1:C:330:PRO:HB3	2.54	0.43
1:D:358:ILE:H	1:D:358:ILE:HG12	1.35	0.43
1:B:170:ASN:OD1	1:B:170:ASN:N	2.52	0.43
1:A:396:LYS:O	1:A:396:LYS:HG3	2.18	0.42
1:C:290:TYR:OH	1:C:292:PRO:HB3	2.18	0.42
1:A:272:ILE:HD11	1:A:310:LEU:HD23	2.01	0.42
1:B:68:ARG:HG2	1:B:74:LEU:HD13	2.02	0.42
1:B:148:ILE:HB	1:B:180:PHE:CD1	2.54	0.42
1:D:265:THR:OG1	1:D:266:LYS:N	2.52	0.42
1:C:188:GLU:CD	1:C:188:GLU:H	2.28	0.42
1:D:211:TRP:HB2	1:D:213:TYR:HE1	1.84	0.42
1:D:259:LEU:HD21	1:D:310:LEU:HD21	2.02	0.42
1:B:359:GLU:OE1	1:B:359:GLU:N	2.27	0.42
1:C:376:HIS:HA	1:C:377:SER:HA	1.81	0.41
1:A:99:ASP:O	1:A:102:VAL:HG22	2.19	0.41
1:A:151:LEU:O	1:A:158:ARG:NH1	2.52	0.41
1:C:216:THR:OG1	1:C:219:ASN:HB2	2.20	0.41
1:C:228:LEU:HB3	1:C:320:TYR:HB2	2.02	0.41
1:C:285:LYS:HG3	1:C:290:TYR:CZ	2.55	0.41
1:D:170:ASN:C	1:D:172:GLY:HA3	2.45	0.41
1:D:352:GLY:HA3	1:D:373:MET:HG2	2.02	0.41
1:C:182:LEU:O	1:C:210:MET:HA	2.21	0.41
1:D:283:ARG:NH2	1:D:291:ILE:O	2.47	0.41
1:C:155:PHE:HE2	1:C:194:LEU:HD13	1.86	0.40
1:C:392:SER:HA	1:C:394:HIS:HE1	1.85	0.40
1:B:272:ILE:HA	1:B:348:GLU:O	2.22	0.40
1:C:168:GLU:HG2	1:C:177:VAL:HA	2.02	0.40
1:C:249:PHE:HB3	1:C:375:VAL:HG12	2.03	0.40
1:D:148:ILE:HD12	1:D:165:TRP:HH2	1.85	0.40
1:A:185:THR:HB	1:A:191:HIS:CG	2.56	0.40
1:A:229:ARG:HB2	1:A:320:TYR:CZ	2.56	0.40
1:A:356:PHE:O	1:A:357:ASP:HB2	2.21	0.40
1:B:66:LEU:HD11	1:B:92:HIS:CD2	2.56	0.40
1:C:213:TYR:HB2	2:G:1:NAG:H62	2.03	0.40
1:D:282:HIS:CD2	1:D:282:HIS:N	2.87	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	320/397 (81%)	303 (95%)	15 (5%)	2 (1%)	21	44
1	B	327/397 (82%)	304 (93%)	23 (7%)	0	100	100
1	C	319/397 (80%)	303 (95%)	15 (5%)	1 (0%)	36	60
1	D	312/397 (79%)	294 (94%)	18 (6%)	0	100	100
All	All	1278/1588 (80%)	1204 (94%)	71 (6%)	3 (0%)	43	68

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	357	ASP
1	C	76	MET
1	A	50	LYS

5.3.2 Protein sidechains

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	302/362 (83%)	296 (98%)	6 (2%)	48	76
1	B	304/362 (84%)	297 (98%)	7 (2%)	44	73
1	C	301/362 (83%)	289 (96%)	12 (4%)	28	56
1	D	294/362 (81%)	287 (98%)	7 (2%)	43	72
All	All	1201/1448 (83%)	1169 (97%)	32 (3%)	39	69

All (32) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	91	SER
1	A	175	THR
1	A	255	ILE
1	A	261	SER
1	A	365	ASN
1	A	377	SER
1	B	129	SER
1	B	170	ASN
1	B	248	VAL
1	B	252	THR
1	B	266	LYS
1	B	286	LYS
1	B	392	SER
1	C	77	LEU
1	C	136	ASP
1	C	141	LYS
1	C	185	THR
1	C	189	ASP
1	C	261	SER
1	C	294	VAL
1	C	312	SER
1	C	358	ILE
1	C	368	SER
1	C	395	LEU
1	C	397	CYS
1	D	107	THR
1	D	196	ASP
1	D	355	THR
1	D	356	PHE
1	D	358	ILE
1	D	368	SER
1	D	397	CYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (18) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	71	ASN
1	A	205	HIS
1	A	253	HIS
1	A	391	GLN
1	B	170	ASN
1	B	174	GLN
1	B	219	ASN

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Mol	Chain	Res	Type
1	B	260	ASN
1	C	159	GLN
1	C	184	GLN
1	C	205	HIS
1	C	277	HIS
1	C	381	GLN
1	D	71	ASN
1	D	110	ASN
1	D	190	ASN
1	D	277	HIS
1	D	282	HIS

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 ⓘ

20 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	0.47	0	17,19,21	0.50	0
2	NAG	E	2	2	14,14,15	0.61	0	17,19,21	0.51	0
2	BMA	E	3	2	11,11,12	0.62	0	15,15,17	0.74	0
2	MAN	E	4	2	11,11,12	0.92	1 (9%)	15,15,17	1.19	2 (13%)
2	MAN	E	5	2	11,11,12	0.86	1 (9%)	15,15,17	1.15	2 (13%)
2	NAG	F	1	2	14,14,15	0.76	0	17,19,21	1.59	5 (29%)
2	NAG	F	2	2	14,14,15	0.68	0	17,19,21	1.36	3 (17%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	BMA	F	3	2	11,11,12	0.31	0	15,15,17	0.76	0
2	MAN	F	4	2	11,11,12	0.89	1 (9%)	15,15,17	1.03	1 (6%)
2	MAN	F	5	2	11,11,12	0.77	0	15,15,17	1.09	1 (6%)
2	NAG	G	1	2,1	14,14,15	0.44	0	17,19,21	0.47	0
2	NAG	G	2	2	14,14,15	0.33	0	17,19,21	0.59	0
2	BMA	G	3	2	11,11,12	0.69	0	15,15,17	0.70	0
2	MAN	G	4	2	11,11,12	1.36	2 (18%)	15,15,17	1.49	2 (13%)
2	MAN	G	5	2	11,11,12	0.74	0	15,15,17	1.11	2 (13%)
2	NAG	H	1	2,1	14,14,15	0.26	0	17,19,21	0.50	0
2	NAG	H	2	2	14,14,15	0.52	0	17,19,21	0.61	0
2	BMA	H	3	2	11,11,12	0.50	0	15,15,17	0.76	0
2	MAN	H	4	2	11,11,12	0.66	0	15,15,17	1.06	1 (6%)
2	MAN	H	5	2	11,11,12	0.70	0	15,15,17	0.94	2 (13%)

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	-	2/6/23/26	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	-	2/2/19/22	0/1/1/1
2	MAN	E	5	2	-	2/2/19/22	0/1/1/1
2	NAG	F	1	2	-	0/6/23/26	0/1/1/1
2	NAG	F	2	2	-	0/6/23/26	0/1/1/1
2	BMA	F	3	2	-	2/2/19/22	0/1/1/1
2	MAN	F	4	2	-	0/2/19/22	0/1/1/1
2	MAN	F	5	2	-	0/2/19/22	0/1/1/1
2	NAG	G	1	2,1	-	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	NAG	H	1	2,1	-	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	-	2/2/19/22	0/1/1/1

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	G	4	MAN	C1-C2	3.50	1.60	1.52
2	G	4	MAN	C2-C3	2.57	1.56	1.52
2	E	4	MAN	C1-C2	2.55	1.58	1.52
2	E	5	MAN	C1-C2	2.45	1.58	1.52
2	F	4	MAN	C1-C2	2.44	1.58	1.52

All (21) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	G	5	MAN	C1-O5-C5	3.18	116.45	112.19
2	G	4	MAN	C1-C2-C3	3.01	114.03	109.64
2	F	2	NAG	C8-C7-N2	-2.96	111.20	116.12
2	F	5	MAN	C1-O5-C5	2.95	116.14	112.19
2	E	4	MAN	C1-O5-C5	2.92	116.11	112.19
2	G	4	MAN	C1-O5-C5	2.90	116.07	112.19
2	H	4	MAN	C1-O5-C5	2.88	116.05	112.19
2	F	1	NAG	O5-C1-C2	-2.73	107.06	111.29
2	F	2	NAG	C2-N2-C7	-2.71	119.26	122.90
2	F	1	NAG	C8-C7-N2	-2.63	111.76	116.12
2	E	5	MAN	C1-O5-C5	2.51	115.55	112.19
2	F	4	MAN	C1-O5-C5	2.40	115.40	112.19
2	F	1	NAG	C2-N2-C7	-2.39	119.70	122.90
2	F	1	NAG	C1-O5-C5	-2.36	109.03	112.19
2	F	1	NAG	O7-C7-N2	2.28	126.00	121.98
2	H	5	MAN	C1-O5-C5	2.24	115.19	112.19
2	G	5	MAN	O2-C2-C3	-2.20	105.60	110.15
2	E	4	MAN	O2-C2-C3	-2.12	105.77	110.15
2	E	5	MAN	O2-C2-C3	-2.11	105.79	110.15
2	F	2	NAG	O7-C7-N2	2.08	125.66	121.98
2	H	5	MAN	O2-C2-C3	-2.02	105.97	110.15

There are no chirality outliers.

All (10) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	E	5	MAN	O5-C5-C6-O6
2	E	5	MAN	C4-C5-C6-O6
2	E	4	MAN	O5-C5-C6-O6
2	E	1	NAG	C4-C5-C6-O6
2	E	1	NAG	O5-C5-C6-O6
2	H	5	MAN	C4-C5-C6-O6

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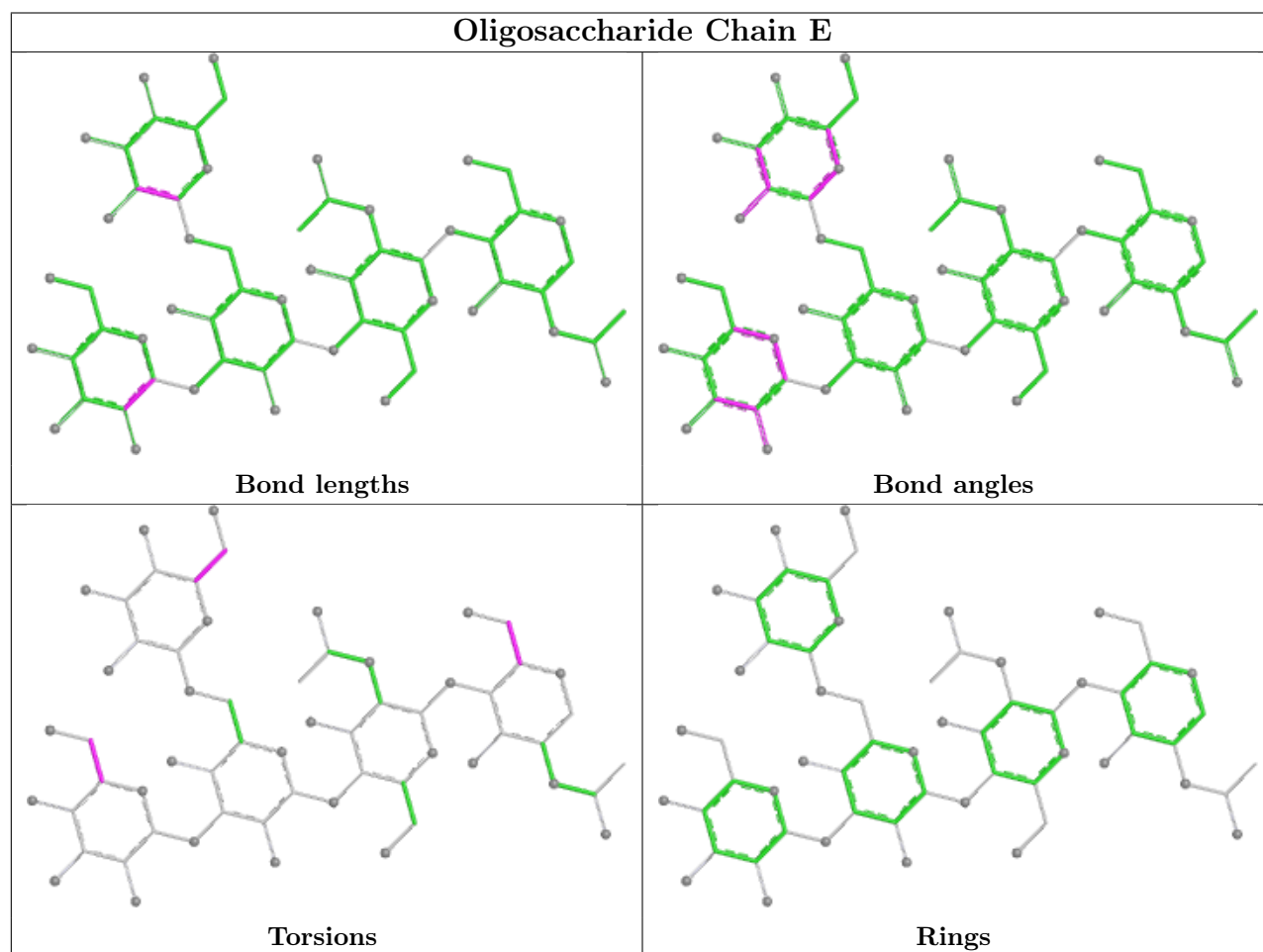
Mol	Chain	Res	Type	Atoms
2	F	3	BMA	C4-C5-C6-O6
2	F	3	BMA	O5-C5-C6-O6
2	H	5	MAN	O5-C5-C6-O6
2	E	4	MAN	C4-C5-C6-O6

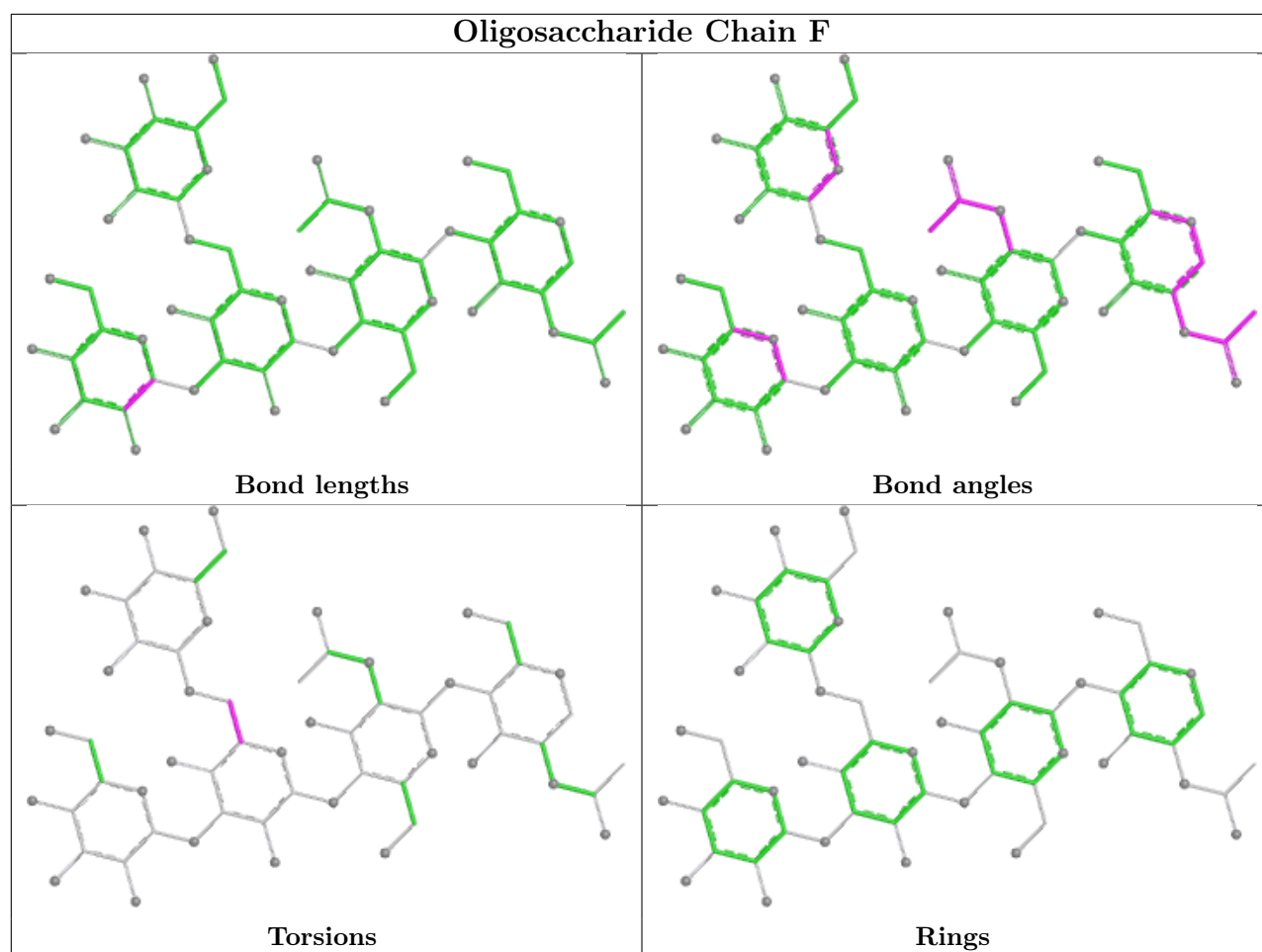
There are no ring outliers.

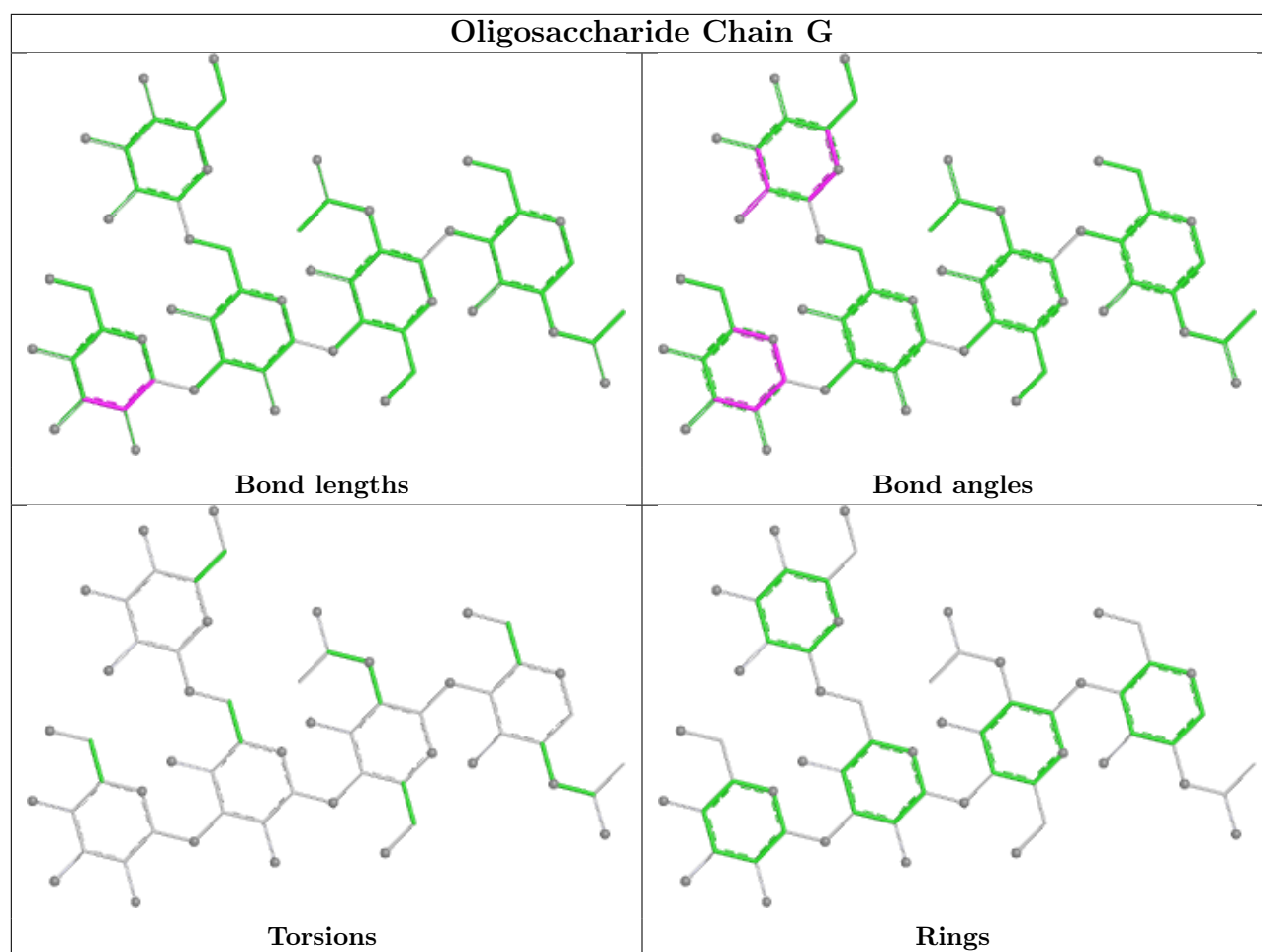
3 monomers are involved in 8 short contacts:

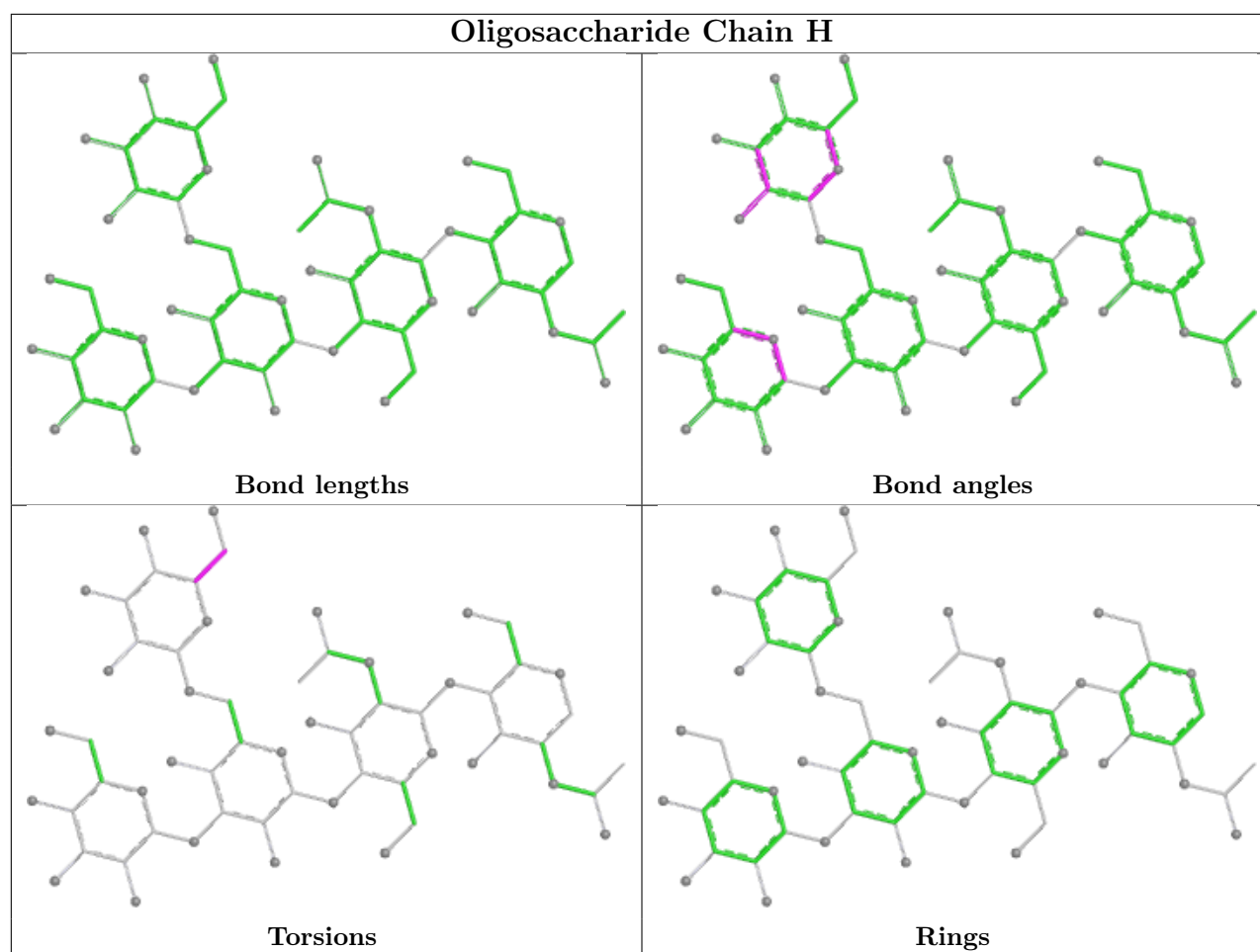
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	G	2	NAG	1	0
2	F	1	NAG	6	0
2	G	1	NAG	1	0

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









5.6 Ligand geometry [i](#)

Of 12 ligands modelled in this entry, 2 are monoatomic - leaving 10 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
3	NAG	A	401	1	14,14,15	0.53	0	17,19,21	0.68	1 (5%)
5	GOL	C	503	-	5,5,5	1.01	0	5,5,5	1.05	0
5	GOL	A	403	-	5,5,5	0.95	0	5,5,5	1.10	0
4	FKX	A	402	-	18,18,18	3.82	8 (44%)	25,26,26	4.30	10 (40%)
4	FKX	C	501	-	18,18,18	3.82	7 (38%)	25,26,26	3.63	9 (36%)
4	FKX	D	501	-	18,18,18	3.80	7 (38%)	25,26,26	4.19	8 (32%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	NAG	B	401	1	14,14,15	0.31	0	17,19,21	0.56	0
5	GOL	B	403	-	5,5,5	0.83	0	5,5,5	1.19	1 (20%)
5	GOL	C	502	-	5,5,5	0.98	0	5,5,5	1.03	0
4	FKX	B	402	-	18,18,18	3.84	8 (44%)	25,26,26	4.50	9 (36%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	NAG	A	401	1	-	1/6/23/26	0/1/1/1
5	GOL	C	503	-	-	2/4/4/4	-
5	GOL	A	403	-	-	0/4/4/4	-
4	FKX	A	402	-	-	4/8/8/8	0/2/2/2
4	FKX	C	501	-	-	0/8/8/8	0/2/2/2
4	FKX	D	501	-	-	0/8/8/8	0/2/2/2
3	NAG	B	401	1	-	2/6/23/26	0/1/1/1
5	GOL	B	403	-	-	4/4/4/4	-
5	GOL	C	502	-	-	4/4/4/4	-
4	FKX	B	402	-	-	0/8/8/8	0/2/2/2

All (30) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	402	FKX	O12-C11	9.72	1.40	1.23
4	A	402	FKX	O12-C11	9.69	1.40	1.23
4	C	501	FKX	O12-C11	9.54	1.40	1.23
4	D	501	FKX	O12-C11	9.49	1.39	1.23
4	C	501	FKX	C02-C01	6.52	1.50	1.39
4	A	402	FKX	C02-C01	6.47	1.50	1.39
4	B	402	FKX	C02-C01	6.47	1.50	1.39
4	D	501	FKX	C02-C01	6.33	1.49	1.39
4	A	402	FKX	C08-C09	5.97	1.49	1.39
4	D	501	FKX	C08-C09	5.97	1.49	1.39
4	C	501	FKX	C08-C09	5.96	1.49	1.39
4	B	402	FKX	C08-C09	5.91	1.49	1.39
4	D	501	FKX	C09-N10	5.62	1.48	1.38
4	A	402	FKX	C09-N10	5.59	1.48	1.38
4	B	402	FKX	C09-N10	5.55	1.48	1.38
4	C	501	FKX	C09-N10	5.48	1.48	1.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	C	501	FKX	C01-N13	4.69	1.48	1.39
4	A	402	FKX	C01-N13	4.63	1.48	1.39
4	B	402	FKX	C01-N13	4.51	1.48	1.39
4	D	501	FKX	C01-N13	4.51	1.48	1.39
4	D	501	FKX	C11-N13	-4.09	1.35	1.38
4	B	402	FKX	C11-N13	-3.85	1.35	1.38
4	C	501	FKX	C11-N13	-3.79	1.35	1.38
4	C	501	FKX	C07-C03	3.62	1.44	1.39
4	B	402	FKX	C07-C03	3.43	1.44	1.39
4	A	402	FKX	C11-N13	-3.41	1.35	1.38
4	D	501	FKX	C07-C03	3.35	1.44	1.39
4	A	402	FKX	C07-C03	3.32	1.44	1.39
4	B	402	FKX	C11-N10	-2.10	1.34	1.37
4	A	402	FKX	C11-N10	-2.06	1.34	1.37

All (38) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	402	FKX	C01-N13-C11	17.86	114.13	109.57
4	D	501	FKX	C01-N13-C11	16.86	113.88	109.57
4	A	402	FKX	C01-N13-C11	16.67	113.83	109.57
4	C	501	FKX	C01-N13-C11	14.15	113.18	109.57
4	B	402	FKX	C09-C01-N13	-8.24	103.78	106.74
4	A	402	FKX	C09-C01-N13	-7.30	104.12	106.74
4	D	501	FKX	C09-C01-N13	-7.20	104.15	106.74
4	C	501	FKX	C09-C01-N13	-6.51	104.40	106.74
4	A	402	FKX	N10-C11-N13	-6.22	103.93	106.76
4	B	402	FKX	N10-C11-N13	-6.15	103.96	106.76
4	D	501	FKX	N10-C11-N13	-5.72	104.16	106.76
4	A	402	FKX	C09-N10-C11	5.63	113.73	110.27
4	B	402	FKX	C09-N10-C11	5.10	113.40	110.27
4	D	501	FKX	C09-N10-C11	4.93	113.30	110.27
4	C	501	FKX	C09-N10-C11	4.73	113.18	110.27
4	A	402	FKX	O12-C11-N13	4.70	129.48	125.82
4	C	501	FKX	N10-C11-N13	-4.47	104.73	106.76
4	B	402	FKX	O12-C11-N13	3.97	128.92	125.82
4	D	501	FKX	O12-C11-N13	3.86	128.83	125.82
4	C	501	FKX	O12-C11-N13	3.63	128.66	125.82
4	B	402	FKX	C15-C14-N13	-3.11	108.19	113.72
4	A	402	FKX	C02-C01-N13	2.74	136.38	130.18
4	B	402	FKX	C02-C01-N13	2.72	136.34	130.18
4	A	402	FKX	C02-C01-C09	-2.52	119.50	121.55

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	D	501	FKX	C08-C09-N10	2.50	135.88	130.83
4	C	501	FKX	C02-C01-N13	2.48	135.81	130.18
4	D	501	FKX	C02-C01-N13	2.33	135.47	130.18
4	C	501	FKX	C08-C09-N10	2.33	135.54	130.83
4	A	402	FKX	C01-C09-N10	-2.30	104.39	106.92
4	A	402	FKX	C08-C09-N10	2.21	135.30	130.83
4	C	501	FKX	C01-C09-N10	-2.20	104.50	106.92
4	D	501	FKX	C01-C09-N10	-2.19	104.51	106.92
4	A	402	FKX	O06-C04-C03	2.18	120.43	114.84
4	C	501	FKX	C02-C01-C09	-2.17	119.78	121.55
3	A	401	NAG	C1-O5-C5	2.13	115.03	112.19
4	B	402	FKX	C08-C09-N10	2.10	135.07	130.83
5	B	403	GOL	C3-C2-C1	-2.05	104.26	111.80
4	B	402	FKX	C02-C01-C09	-2.05	119.88	121.55

There are no chirality outliers.

All (17) torsion outliers are listed below:

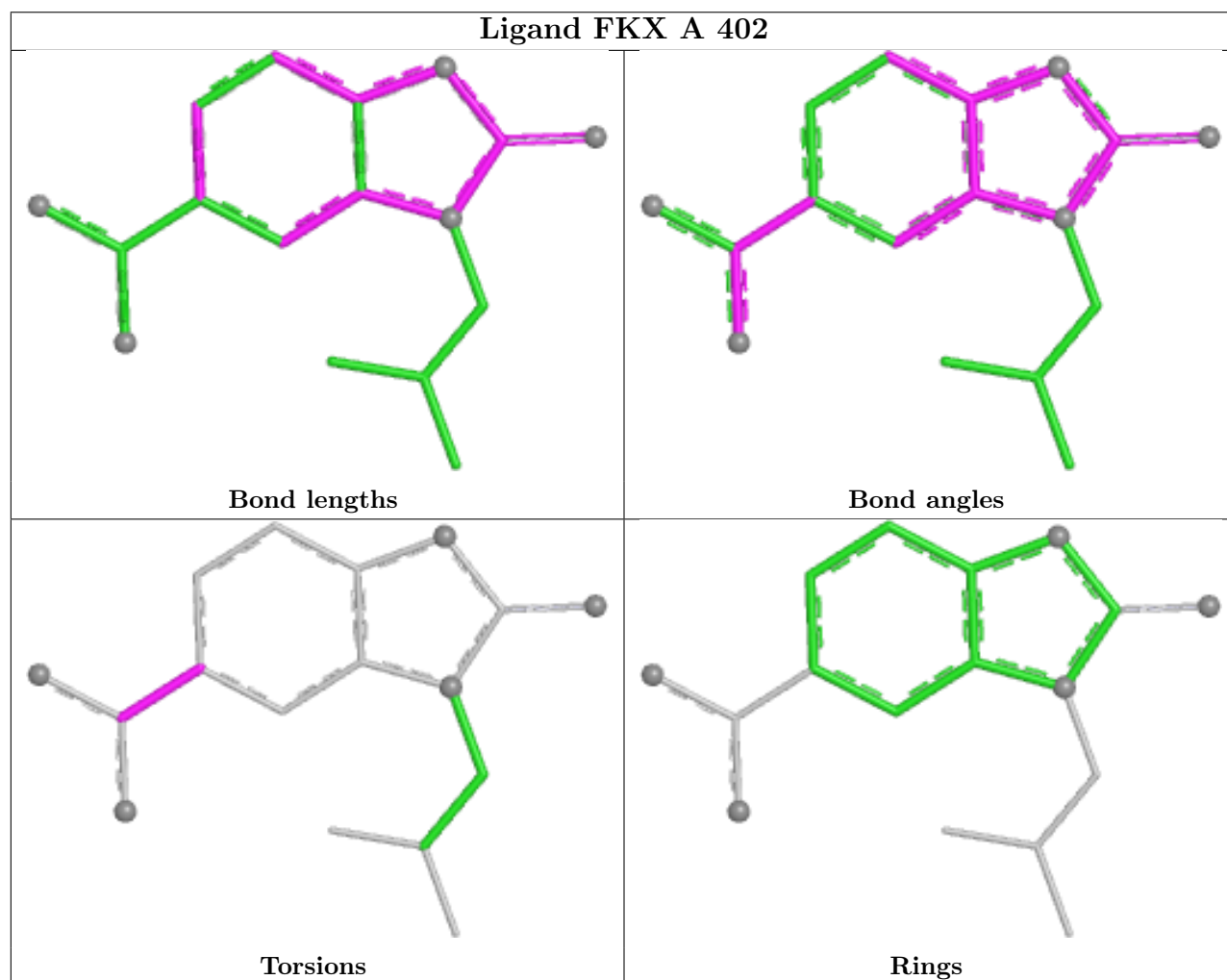
Mol	Chain	Res	Type	Atoms
5	B	403	GOL	O1-C1-C2-C3
5	C	502	GOL	O1-C1-C2-C3
5	C	502	GOL	C1-C2-C3-O3
5	C	503	GOL	O1-C1-C2-O2
5	C	503	GOL	O1-C1-C2-C3
3	B	401	NAG	O5-C5-C6-O6
3	B	401	NAG	C4-C5-C6-O6
5	B	403	GOL	C1-C2-C3-O3
5	C	502	GOL	O1-C1-C2-O2
5	C	502	GOL	O2-C2-C3-O3
4	A	402	FKX	C07-C03-C04-O05
4	A	402	FKX	C02-C03-C04-O05
4	A	402	FKX	C07-C03-C04-O06
4	A	402	FKX	C02-C03-C04-O06
5	B	403	GOL	O2-C2-C3-O3
5	B	403	GOL	O1-C1-C2-O2
3	A	401	NAG	C4-C5-C6-O6

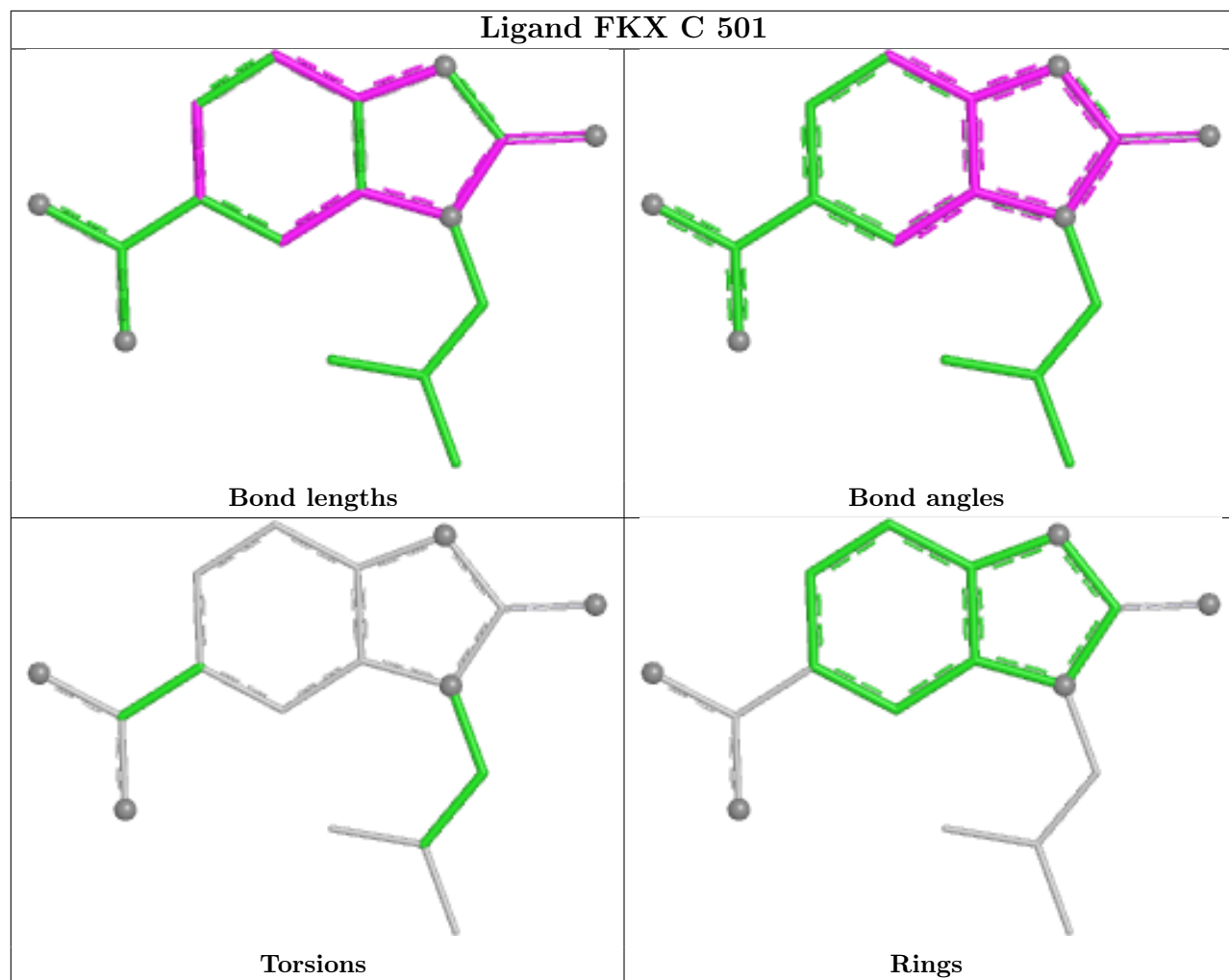
There are no ring outliers.

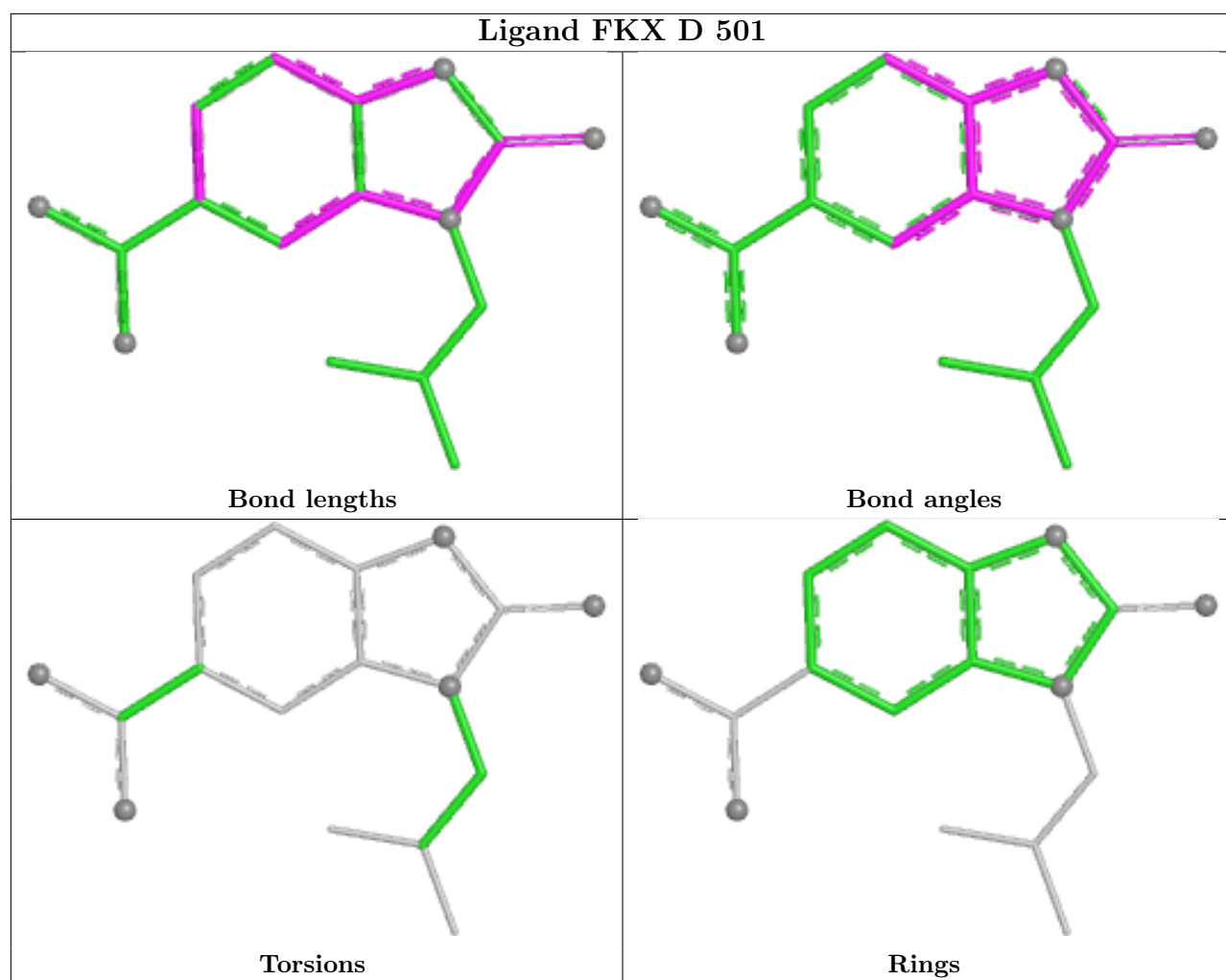
2 monomers are involved in 2 short contacts:

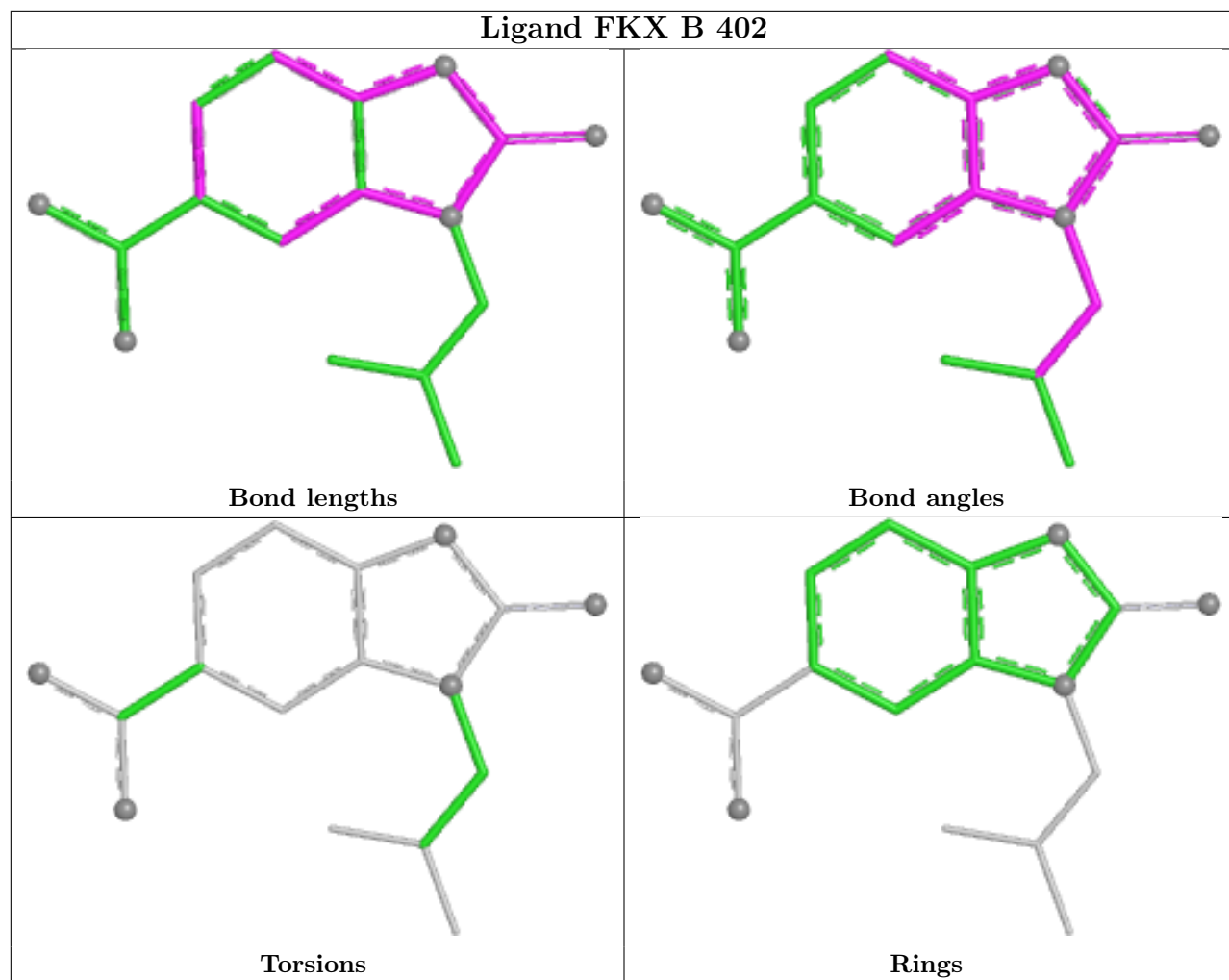
Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	402	FKX	1	0
4	D	501	FKX	1	0

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









5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 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	328/397 (82%)	0.18	10 (3%)	52	49	24, 44, 77, 103	0
1	B	331/397 (83%)	0.23	10 (3%)	52	49	25, 45, 75, 102	0
1	C	327/397 (82%)	0.41	13 (3%)	42	38	28, 50, 78, 109	0
1	D	320/397 (80%)	0.36	10 (3%)	51	48	30, 51, 78, 96	0
All	All	1306/1588 (82%)	0.29	43 (3%)	49	45	24, 48, 77, 109	0

All (43) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	90	ILE	4.5
1	B	55	PRO	4.1
1	B	171	ALA	4.0
1	C	358	ILE	4.0
1	C	170	ASN	3.6
1	B	397	CYS	3.6
1	C	90	ILE	3.5
1	C	356	PHE	3.3
1	C	169	SER	3.1
1	C	173	ASN	3.1
1	A	358	ILE	3.1
1	A	356	PHE	3.1
1	C	55	PRO	3.0
1	D	358	ILE	3.0
1	D	172	GLY	2.9
1	B	173	ASN	2.9
1	C	91	SER	2.7
1	D	362	ASN	2.7
1	C	357	ASP	2.7
1	B	265	THR	2.6
1	D	355	THR	2.6

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Mol	Chain	Res	Type	RSRZ
1	A	173	ASN	2.5
1	D	91	SER	2.5
1	A	355	THR	2.4
1	A	397	CYS	2.4
1	B	172	GLY	2.4
1	A	95	TYR	2.4
1	B	253	HIS	2.4
1	C	355	THR	2.3
1	D	173	ASN	2.3
1	A	395	LEU	2.3
1	D	395	LEU	2.3
1	A	90	ILE	2.3
1	C	369	TYR	2.3
1	B	170	ASN	2.3
1	C	172	GLY	2.3
1	C	278	ASN	2.3
1	D	397	CYS	2.2
1	D	356	PHE	2.2
1	B	321	HIS	2.1
1	A	351	LYS	2.1
1	A	172	GLY	2.1
1	D	361	LYS	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.

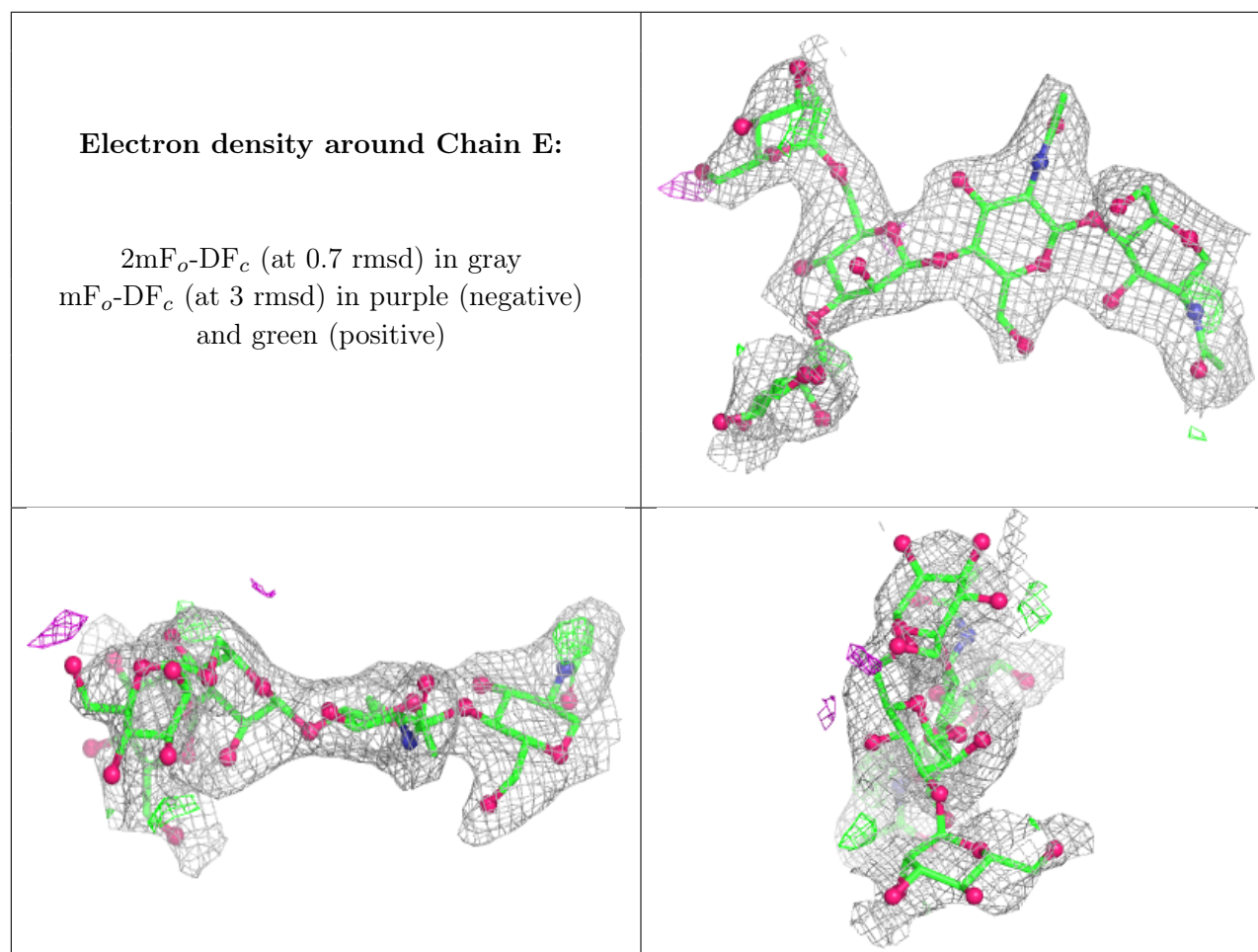
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	MAN	E	4	11/12	0.35	0.19	89,104,109,113	0
2	MAN	G	4	11/12	0.36	0.18	83,92,104,105	0
2	MAN	H	4	11/12	0.47	0.17	102,113,130,130	0
2	MAN	F	5	11/12	0.53	0.17	74,94,98,99	0
2	MAN	H	5	11/12	0.53	0.19	85,108,121,126	0
2	MAN	E	5	11/12	0.56	0.17	88,99,104,105	0

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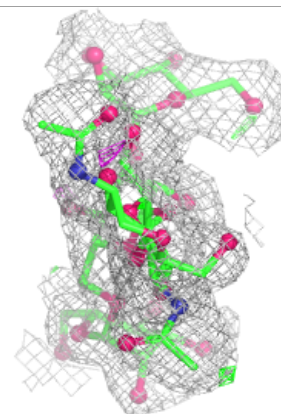
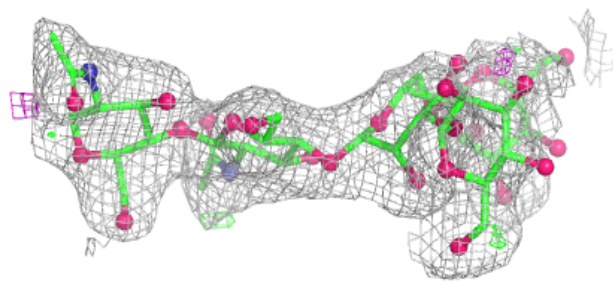
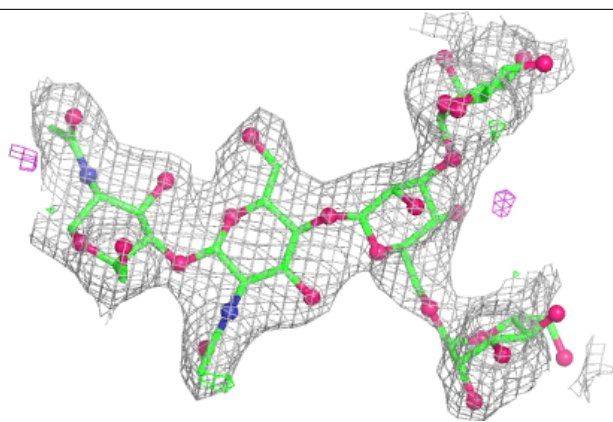
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	MAN	G	5	11/12	0.63	0.15	81,92,102,104	0
2	MAN	F	4	11/12	0.69	0.14	60,89,101,102	0
2	BMA	H	3	11/12	0.75	0.14	70,93,99,106	0
2	BMA	G	3	11/12	0.80	0.13	57,67,83,89	0
2	BMA	E	3	11/12	0.84	0.11	59,75,85,95	0
2	BMA	F	3	11/12	0.84	0.12	56,72,88,93	0
2	NAG	H	2	14/15	0.89	0.11	44,55,65,66	0
2	NAG	F	2	14/15	0.89	0.11	35,42,51,52	0
2	NAG	G	2	14/15	0.92	0.09	37,43,58,59	0
2	NAG	F	1	14/15	0.92	0.10	25,35,40,42	0
2	NAG	G	1	14/15	0.93	0.09	27,31,35,36	0
2	NAG	E	1	14/15	0.93	0.09	23,36,42,44	0
2	NAG	E	2	14/15	0.93	0.08	30,42,53,55	0
2	NAG	H	1	14/15	0.94	0.09	33,39,44,51	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.

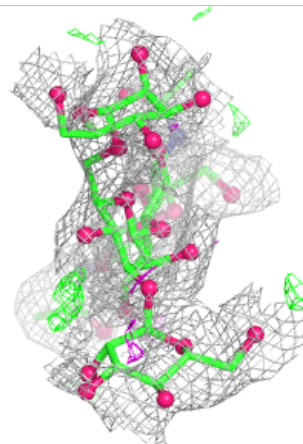
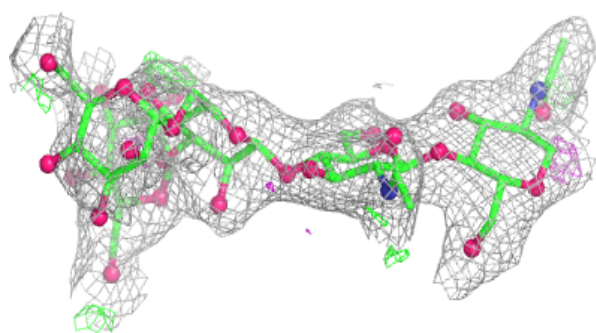
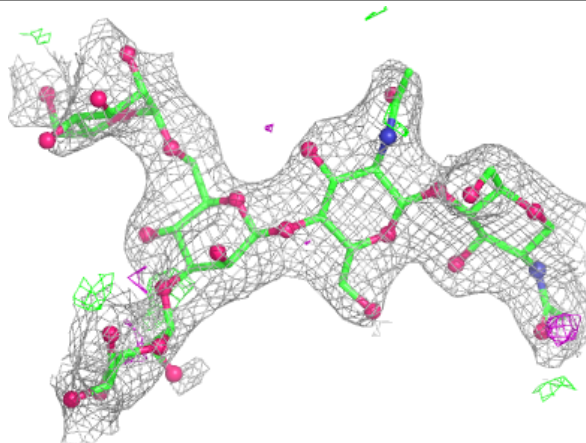


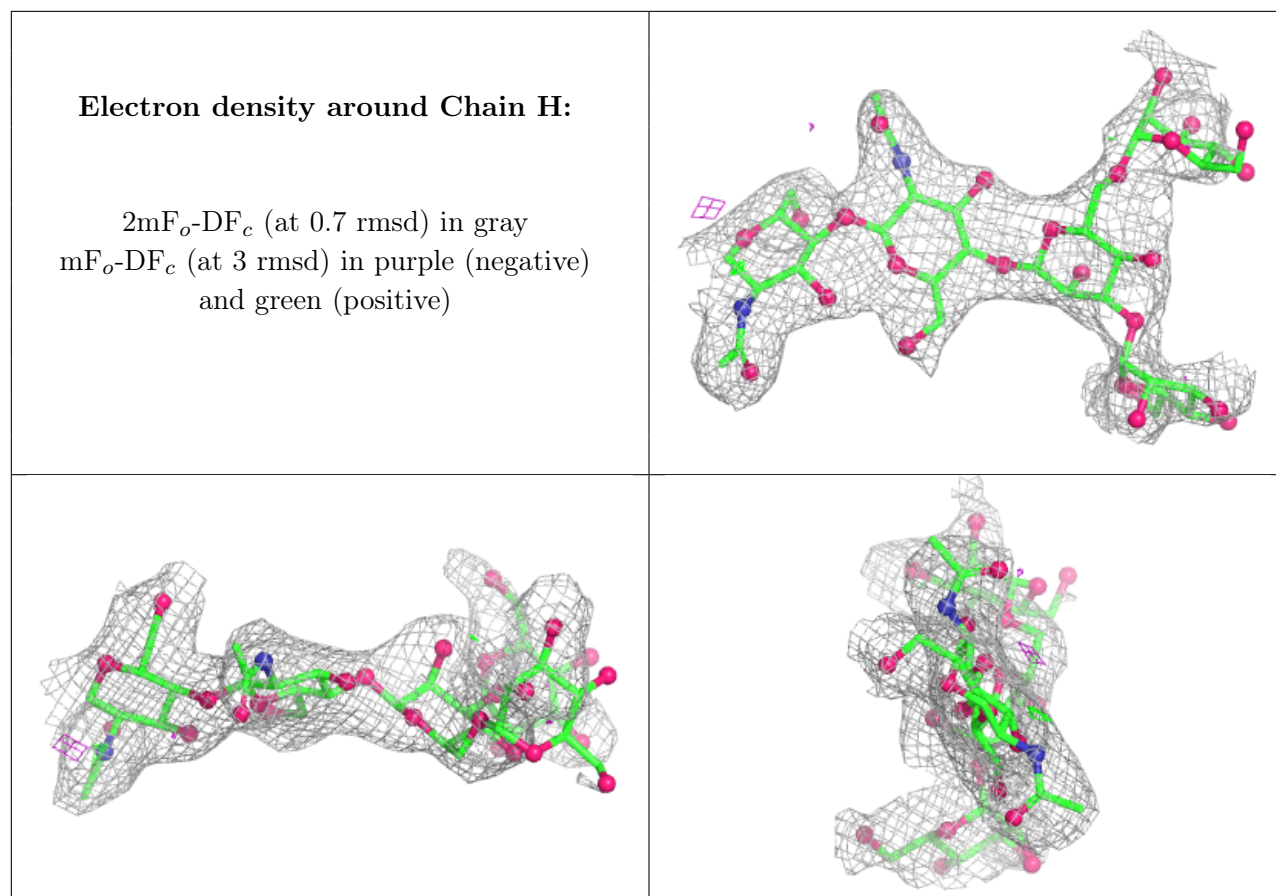
Electron density around Chain F:

$2mF_o-DF_c$ (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)





6.4 Ligands ⓘ

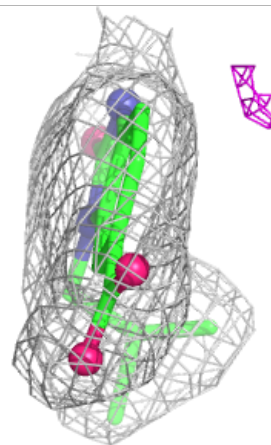
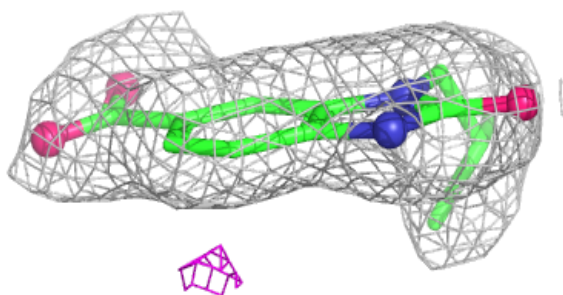
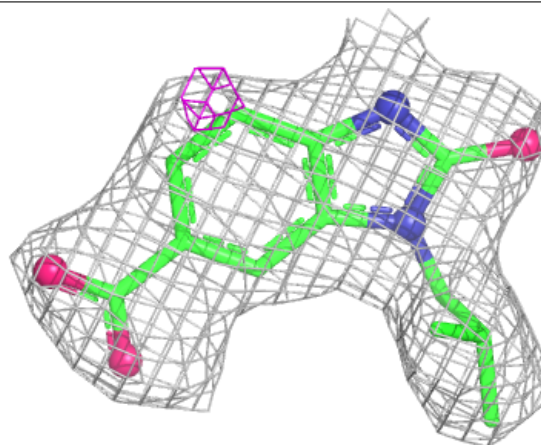
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
5	GOL	C	503	6/6	0.79	0.17	45,60,67,68	0
5	GOL	C	502	6/6	0.81	0.16	45,57,69,69	0
3	NAG	A	401	14/15	0.83	0.11	41,50,58,59	0
3	NAG	B	401	14/15	0.84	0.10	52,64,68,69	0
5	GOL	B	403	6/6	0.85	0.17	37,46,49,55	0
5	GOL	A	403	6/6	0.91	0.09	39,44,49,60	0
4	FKX	D	501	17/17	0.92	0.10	26,34,49,51	0
4	FKX	B	402	17/17	0.94	0.09	19,30,39,42	0
4	FKX	A	402	17/17	0.94	0.09	30,34,42,51	0
4	FKX	C	501	17/17	0.95	0.09	27,32,38,44	0
6	CL	D	502	1/1	0.96	0.07	43,43,43,43	0
6	CL	C	504	1/1	0.97	0.05	40,40,40,40	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.

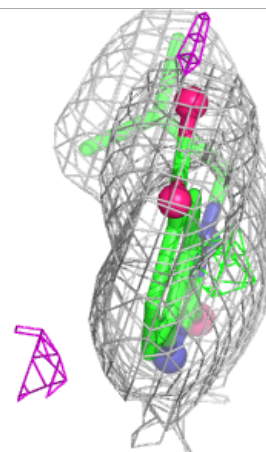
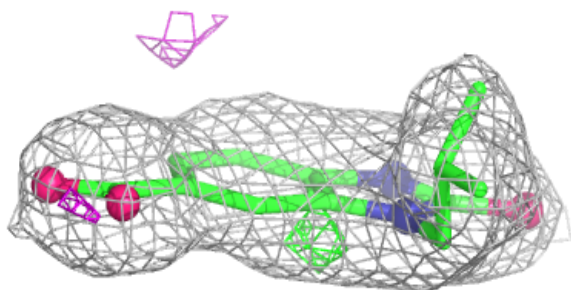
Electron density around FKX D 501:

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



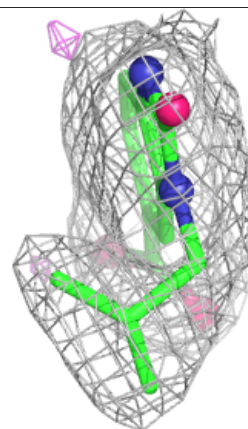
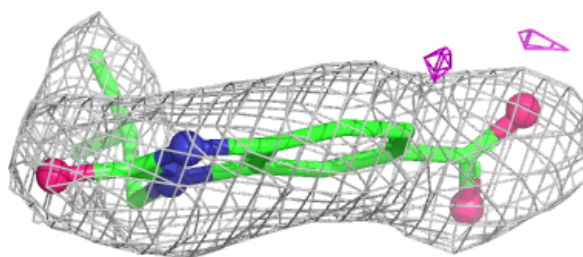
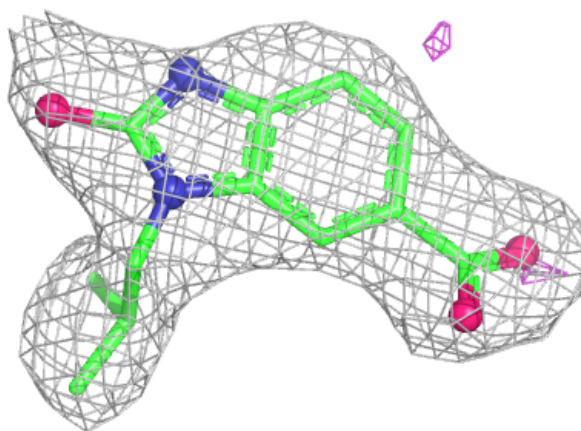
Electron density around FKX B 402:

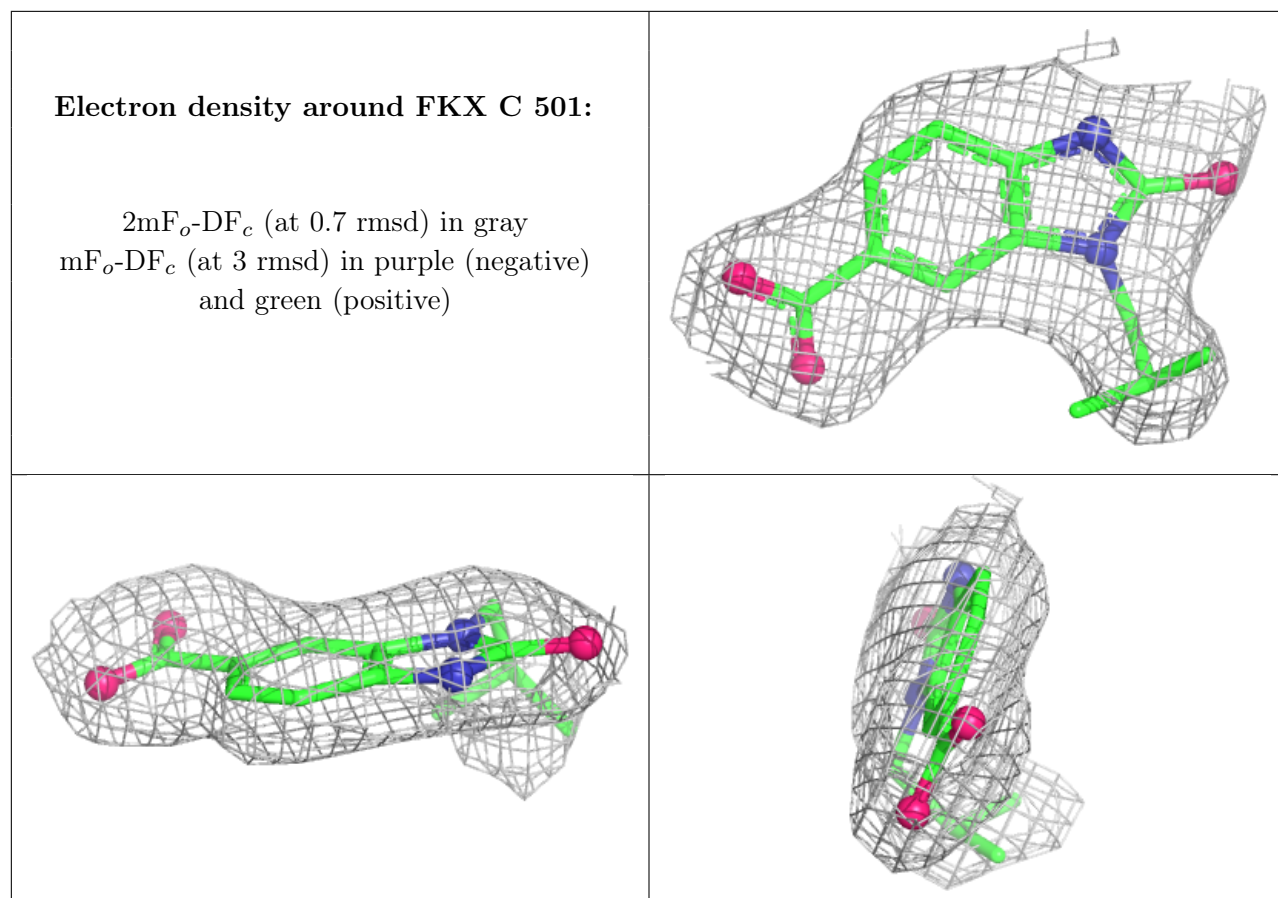
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around FKX A 402:

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





6.5 Other polymers ⓘ

There are no such residues in this entry.