



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 21, 2020 – 08:49 PM BST

PDB ID : 6VW1
Title : Structure of SARS-CoV-2 chimeric receptor-binding domain complexed with its receptor human ACE2
Authors : Shang, J.; Ye, G.; Shi, K.; Wan, Y.S.; Aihara, H.; Li, F.
Deposited on : 2020-02-18
Resolution : 2.68 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.

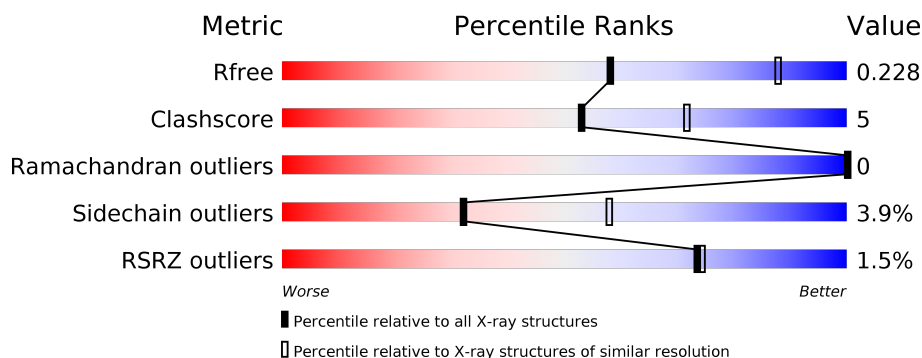
The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.13.1
buster-report : 1.1.7 (2018)
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.13.1

i

X-RAY DIFFRACTION

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.




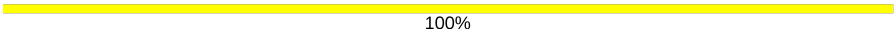

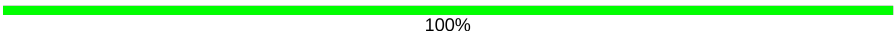

Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R _{free}	130704	3863 (2.70-2.66)
Clashscore	141614	4210 (2.70-2.66)
Ramachandran outliers	138981	4141 (2.70-2.66)
Sidechain outliers	138945	4141 (2.70-2.66)
RSRZ outliers	127900	3780 (2.70-2.66)

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 on 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	597	<div> <div></div> <div>85%</div> <div>14%</div> <div></div> </div>
1	B	597	<div> <div></div> <div>86%</div> <div>13%</div> <div></div> </div>
2	E	217	<div> <div></div> <div>67%</div> <div>20%</div> <div></div> <div>11%</div> </div>
2	F	217	<div> <div></div> <div>73%</div> <div>14%</div> <div></div> <div>11%</div> </div>
3	C	3	<div> <div></div> <div>100%</div> </div>
3	D	3	<div> <div></div> <div>33%</div> <div>67%</div> </div>

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

2 Entry composition [i](#)

There are 10 unique types of molecules in this entry. The entry contains 13180 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 Angiotensin-converting enzyme 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	596	Total	C	N	O	S	0	0	0
			4862	3111	805	917	29			
1	B	596	Total	C	N	O	S	0	0	0
			4862	3111	805	917	29			

- Molecule 2 is a protein called SARS-CoV-2 chimeric RBD.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	E	193	Total	C	N	O	S	0	0	0
			1527	983	247	288	9			
2	F	194	Total	C	N	O	S	0	0	0
			1531	984	248	290	9			

- Molecule 3 is an oligosaccharide called 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
3	C	3	Total	C	N	O	0	0	0
			39	22	2	15			
3	D	3	Total	C	N	O	0	0	0
			39	22	2	15			
3	L	3	Total	C	N	O	0	0	0
			39	22	2	15			

- 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	G	2	Total	C	N	O	0	0	0
			28	16	2	10			
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	K	2	Total	C	N	O	0	0	0
			28	16	2	10			
4	M	2	Total	C	N	O	0	0	0
			28	16	2	10			

- Molecule 5 is an oligosaccharide called 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-3)-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
5	J	4	Total	C	N	O	0	0	0
			53	30	3	20			

- Molecule 6 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	B	1	Total	Zn	0	0
			1	1		
6	A	1	Total	Zn	0	0
			1	1		

- Molecule 7 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	B	1	Total	Cl	0	0
			1	1		
7	A	1	Total	Cl	0	0
			1	1		

- Molecule 8 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: $C_2H_6O_2$).



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

- Molecule 9 is 2-acetamido-2-deoxy-beta-D-glucopyranose (three-letter code: NAG) (formula: $C_8H_{15}NO_6$) (labeled as "Ligand of Interest" by author).



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

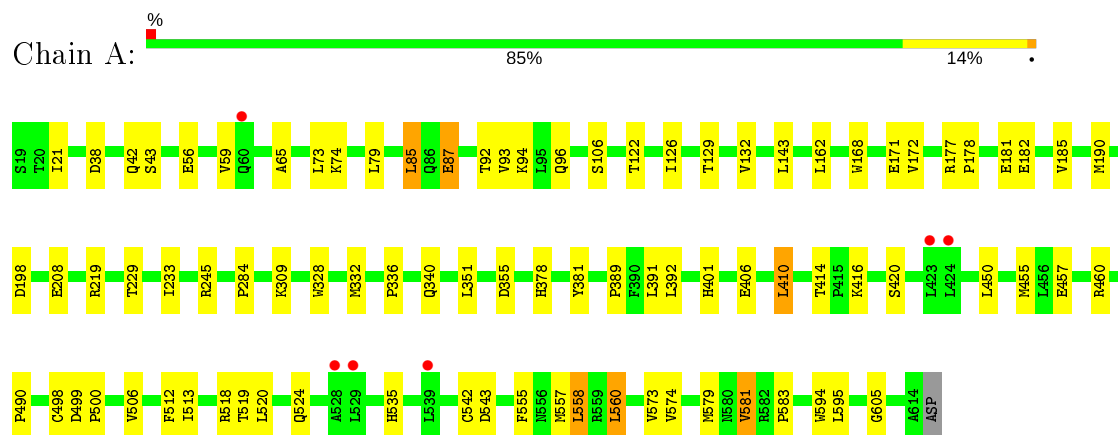
- Molecule 10 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
10	A	13	Total	O	0	0
			13	13		
10	B	8	Total	O	0	0
			8	8		
10	E	4	Total	O	0	0
			4	4		
10	F	1	Total	O	0	0
			1	1		

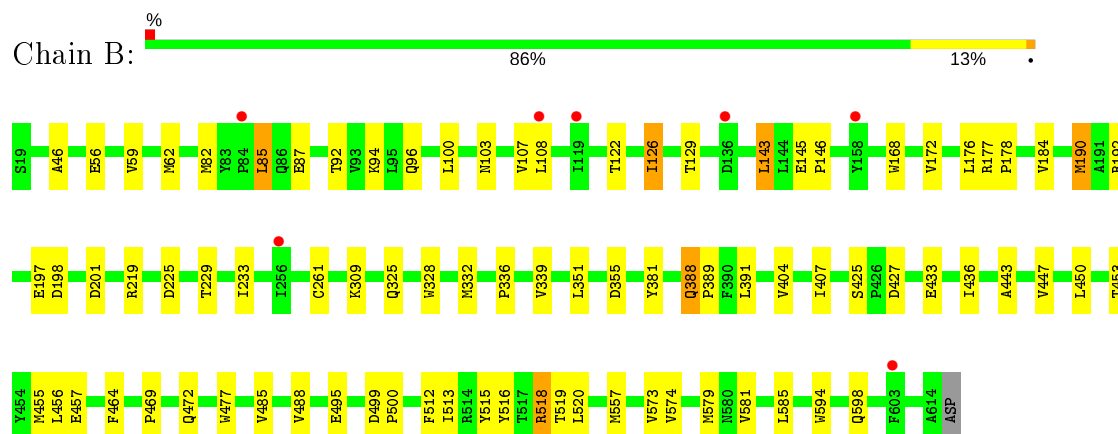
3 Residue-property plots [i](#)

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.

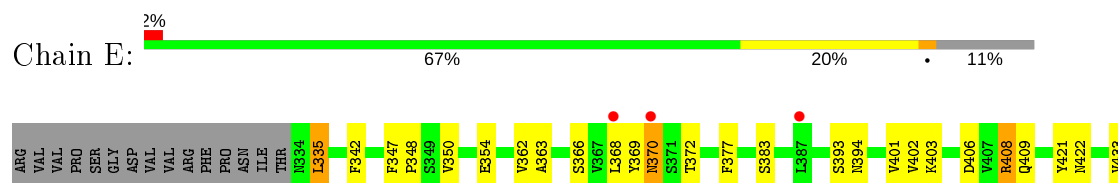
• Molecule 1: Angiotensin-converting enzyme 2

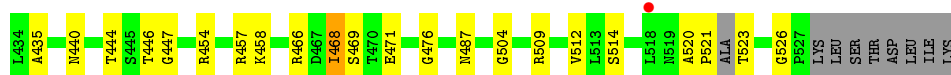


• Molecule 1: Angiotensin-converting enzyme 2

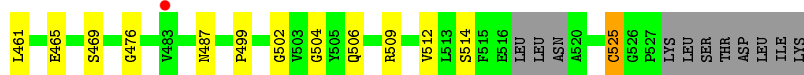


• Molecule 2: SARS-CoV-2 chimeric RBD





- Molecule 2: SARS-CoV-2 chimeric RBD



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



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



- Molecule 3: beta-D-mannopyranose-(1-4)-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





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

Chain I:

100%



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

Chain K:

100%



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

Chain M:

50%

50%



- Molecule 5: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-3)-beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain J:

50%

50%



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	80.44Å 118.03Å 112.08Å 90.00° 93.12° 90.00°	Depositor
Resolution (Å)	59.02 – 2.68 111.91 – 2.68	Depositor EDS
% Data completeness (in resolution range)	99.0 (59.02-2.68) 99.1 (111.91-2.68)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.52 (at 2.69Å)	Xtriage
Refinement program	PHENIX 1.17.1_3660, REFMAC 5.8.0258	Depositor
R, R_{free}	0.197 , 0.229 0.197 , 0.228	Depositor DCC
R_{free} test set	2675 reflections (4.59%)	wwPDB-VP
Wilson B-factor (Å ²)	83.3	Xtriage
Anisotropy	0.179	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 71.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	13180	wwPDB-VP
Average B, all atoms (Å ²)	108.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.08% 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: ZN, BMA, EDO, NAG, CL

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	0.25	0/4999	0.44	0/6792
1	B	0.25	0/4999	0.43	0/6792
2	E	0.27	0/1570	0.47	1/2137 (0.0%)
2	F	0.26	0/1574	0.44	0/2143
All	All	0.26	0/13142	0.44	1/17864 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	E	335	LEU	CA-CB-CG	5.09	127.00	115.30

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	4862	0	4634	49	0
1	B	4862	0	4634	44	0
2	E	1527	0	1437	22	0
2	F	1531	0	1440	17	0
3	C	39	0	34	0	0
3	D	39	0	34	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	L	39	0	34	2	0
4	G	28	0	25	0	0
4	H	28	0	25	4	0
4	I	28	0	25	0	0
4	K	28	0	25	0	0
4	M	28	0	25	1	0
5	J	53	0	46	0	0
6	A	1	0	0	0	0
6	B	1	0	0	0	0
7	A	1	0	0	1	0
7	B	1	0	0	0	0
8	A	4	0	6	0	0
8	B	12	0	18	0	0
9	A	14	0	13	0	0
9	B	28	0	26	1	0
10	A	13	0	0	1	0
10	B	8	0	0	1	0
10	E	4	0	0	0	0
10	F	1	0	0	0	0
All	All	13180	0	12481	138	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 5.

All (138) 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)
2:E:454:ARG:NH2	2:E:469:SER:O	2.28	0.66
1:B:85:LEU:HD11	1:B:94:LYS:HG3	1.78	0.64
1:B:126:ILE:HD11	1:B:176:LEU:HG	1.80	0.64
1:B:433:GLU:HG2	9:B:714:NAG:H82	1.79	0.63
1:A:177:ARG:HD3	1:A:498:CYS:HB2	1.79	0.63
1:B:388:GLN:HG3	1:B:389:PRO:HD2	1.83	0.61
2:E:366:SER:HB2	2:E:370:ASN:HD21	1.65	0.61
4:M:2:NAG:H3	4:M:2:NAG:H83	1.84	0.60
1:A:74:LYS:HE3	1:A:106:SER:HB3	1.83	0.60
2:E:366:SER:O	2:E:370:ASN:ND2	2.32	0.60
2:E:393:SER:O	2:E:523:THR:OG1	2.18	0.60
2:E:401:VAL:HG22	2:E:509:ARG:HG2	1.86	0.58
1:A:208:GLU:OE1	1:A:219:ARG:NH1	2.38	0.56
1:A:457:GLU:HG2	1:A:512:PHE:HB3	1.87	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:132:VAL:HG12	1:A:171:GLU:HG3	1.88	0.56
1:A:557:MET:HA	1:A:560:LEU:HD22	1.86	0.56
1:B:457:GLU:HG2	1:B:513:ILE:HB	1.88	0.56
2:F:401:VAL:HG22	2:F:509:ARG:HG2	1.88	0.55
2:F:387:LEU:HA	2:F:390:LEU:HD12	1.88	0.55
1:B:229:THR:HG23	1:B:516:TYR:OH	2.06	0.55
1:B:573:VAL:HG13	1:B:574:VAL:HG13	1.87	0.55
1:B:425:SER:OG	1:B:427:ASP:OD1	2.20	0.54
1:A:245:ARG:NH1	1:A:605:GLY:O	2.29	0.54
1:A:573:VAL:HG13	1:A:574:VAL:HG13	1.90	0.54
1:B:450:LEU:HD21	1:B:519:THR:HG21	1.89	0.54
1:B:520:LEU:HD22	1:B:579:MET:HE2	1.88	0.53
2:E:363:ALA:H	2:E:526:GLY:HA3	1.73	0.53
2:F:433:VAL:HG22	2:F:512:VAL:HG13	1.91	0.53
1:B:122:THR:O	1:B:126:ILE:HG23	2.09	0.53
2:E:444:THR:HG23	2:E:447:GLY:H	1.73	0.53
1:B:201:ASP:OD2	1:B:219:ARG:NH1	2.42	0.52
1:A:332:MET:HE3	1:A:336:PRO:HD3	1.91	0.52
1:A:85:LEU:HD11	1:A:94:LYS:HG3	1.92	0.51
1:B:92:THR:HG22	1:B:96:GLN:HE21	1.76	0.51
2:F:350:VAL:HG22	2:F:422:ASN:HB3	1.91	0.51
2:F:454:ARG:NH2	2:F:469:SER:O	2.44	0.51
2:E:421:TYR:HB3	2:E:454:ARG:HG2	1.93	0.51
1:B:233:ILE:HD13	1:B:450:LEU:HD13	1.93	0.51
1:A:336:PRO:HG2	1:A:340:GLN:O	2.11	0.50
1:A:460:ARG:HH21	1:A:506:VAL:HA	1.76	0.50
4:H:1:NAG:H4	4:H:2:NAG:HN2	1.76	0.50
1:A:351:LEU:HB2	1:A:355:ASP:HB3	1.94	0.50
2:E:348:PRO:HG3	2:E:354:GLU:HB2	1.94	0.50
1:B:177:ARG:NH1	1:B:495:GLU:OE2	2.45	0.50
1:A:177:ARG:HB3	1:A:178:PRO:HD3	1.94	0.49
1:A:92:THR:HG22	1:A:96:GLN:HE21	1.77	0.49
1:B:100:LEU:HG	1:B:391:LEU:HD21	1.94	0.49
1:B:225:ASP:O	1:B:229:THR:HG22	2.12	0.49
2:E:335:LEU:HD22	2:E:362:VAL:O	2.12	0.49
1:A:85:LEU:HD12	1:A:94:LYS:HE3	1.94	0.49
1:A:520:LEU:HD21	1:A:581:VAL:HG13	1.95	0.49
1:B:515:TYR:O	1:B:518:ARG:HG3	2.13	0.49
1:A:378:HIS:CE1	1:A:401:HIS:HB3	2.47	0.49
1:B:192:ARG:NH2	1:B:197:GLU:O	2.46	0.48
2:E:476:GLY:H	2:E:487:ASN:HB3	1.79	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:454:ARG:HD3	2:E:457:ARG:HB2	1.94	0.48
1:B:453:THR:HG23	1:B:512:PHE:CD2	2.49	0.48
1:B:177:ARG:HB3	1:B:178:PRO:HD3	1.95	0.48
2:E:520:ALA:HB1	2:E:521:PRO:HD2	1.95	0.48
1:A:94:LYS:NZ	10:A:802:HOH:O	2.47	0.47
2:F:391:CYS:SG	2:F:525:CYS:HB3	2.55	0.47
1:B:168:TRP:O	1:B:172:VAL:HG22	2.15	0.47
2:F:461:LEU:HD22	2:F:465:GLU:HB3	1.97	0.47
1:A:56:GLU:HA	1:A:59:VAL:HG12	1.96	0.46
1:A:168:TRP:O	1:A:172:VAL:HG22	2.15	0.46
1:B:443:ALA:HA	1:B:447:VAL:HG13	1.98	0.46
1:A:229:THR:O	1:A:233:ILE:HG13	2.15	0.46
1:A:460:ARG:NH2	1:A:506:VAL:HA	2.30	0.46
1:B:456:LEU:HD13	1:B:477:TRP:HH2	1.81	0.46
1:A:416:LYS:HE3	1:A:543:ASP:HB3	1.97	0.46
1:B:46:ALA:HB1	1:B:62:MET:HA	1.97	0.46
1:A:499:ASP:HB3	7:A:705:CL:CL	2.53	0.45
1:A:450:LEU:HD21	1:A:519:THR:HG21	1.97	0.45
1:B:351:LEU:HB2	1:B:355:ASP:HB3	1.99	0.45
1:B:332:MET:HE3	1:B:336:PRO:HD3	1.98	0.45
1:B:103:ASN:HB3	1:B:107:VAL:HG13	1.98	0.45
1:B:339:VAL:HG22	10:B:804:HOH:O	2.17	0.45
1:B:184:VAL:HG22	1:B:464:PHE:HE1	1.81	0.45
1:A:420:SER:HA	3:D:2:NAG:H82	1.99	0.45
2:F:394:ASN:N	2:F:394:ASN:OD1	2.49	0.45
1:B:391:LEU:H	1:B:391:LEU:HD12	1.81	0.45
1:A:555:PHE:HA	1:A:558:LEU:HD22	1.98	0.45
1:B:229:THR:OG1	1:B:581:VAL:HB	2.17	0.45
2:E:435:ALA:HA	2:E:509:ARG:O	2.17	0.45
2:F:439:ARG:HD3	2:F:499:PRO:HA	1.99	0.45
1:A:389:PRO:HG2	1:A:392:LEU:HD22	1.98	0.44
1:A:520:LEU:HD22	1:A:579:MET:HE2	1.99	0.44
2:F:383:SER:HG	2:F:386:LYS:HZ2	1.62	0.44
4:H:1:NAG:H61	4:H:2:NAG:C7	2.47	0.44
4:H:1:NAG:H4	4:H:2:NAG:N2	2.32	0.44
1:A:182:GLU:HA	1:A:185:VAL:HG12	2.00	0.44
1:A:309:LYS:HD2	1:A:328:TRP:CH2	2.53	0.44
1:B:108:LEU:HD11	1:B:190:MET:HB2	2.00	0.44
1:A:43:SER:HA	1:A:65:ALA:HB1	1.99	0.44
1:B:469:PRO:HD2	1:B:472:GLN:HE21	1.83	0.44
2:E:402:VAL:HG12	2:E:406:ASP:HB2	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:21:ILE:H	1:A:21:ILE:HD12	1.82	0.44
2:F:335:LEU:HA	2:F:362:VAL:HG23	2.00	0.44
2:E:350:VAL:HG22	2:E:422:ASN:HB3	1.98	0.43
1:B:455:MET:HE2	1:B:485:VAL:HG21	2.00	0.43
1:A:162:LEU:HD13	1:A:490:PRO:HB2	1.99	0.43
2:E:342:PHE:HB2	3:L:1:NAG:H82	1.99	0.43
3:L:2:NAG:H4	3:L:3:BMA:O2	2.18	0.43
1:A:38:ASP:O	1:A:42:GLN:HG2	2.19	0.43
1:B:145:GLU:HA	1:B:146:PRO:HA	1.80	0.43
1:A:177:ARG:O	1:A:181:GLU:HG3	2.18	0.43
1:A:524:GLN:HG3	1:A:583:PRO:HG2	2.00	0.43
2:F:502:GLY:O	2:F:506:GLN:HG3	2.19	0.43
1:B:404:VAL:O	1:B:407:ILE:HG12	2.18	0.43
1:B:499:ASP:N	1:B:500:PRO:HD2	2.34	0.42
2:E:440:ASN:N	2:E:440:ASN:OD1	2.51	0.42
2:E:433:VAL:HG12	2:E:512:VAL:HG22	2.00	0.42
2:F:440:ASN:N	2:F:440:ASN:OD1	2.51	0.42
1:A:391:LEU:HD12	1:A:391:LEU:H	1.83	0.42
1:A:21:ILE:HD13	1:A:87:GLU:HG2	2.02	0.42
2:F:421:TYR:HB3	2:F:454:ARG:HG2	2.01	0.42
2:F:476:GLY:N	2:F:487:ASN:HB3	2.35	0.42
2:E:408:ARG:HD3	2:E:409:GLN:HG3	2.01	0.42
2:F:438:THR:O	2:F:442:ASP:HB2	2.20	0.42
2:F:403:LYS:HG3	2:F:504:GLY:O	2.19	0.42
1:B:56:GLU:HA	1:B:59:VAL:HG12	2.01	0.42
2:E:466:ARG:NH1	2:E:468:ILE:HD11	2.35	0.42
1:A:177:ARG:NH2	1:A:181:GLU:OE2	2.53	0.41
1:A:535:HIS:CD2	1:A:542:CYS:HB2	2.55	0.41
1:A:406:GLU:O	1:A:410:LEU:HD13	2.19	0.41
2:E:403:LYS:HG3	2:E:504:GLY:O	2.19	0.41
1:A:122:THR:O	1:A:126:ILE:HG12	2.20	0.41
1:A:499:ASP:N	1:A:500:PRO:HD2	2.35	0.41
1:A:284:PRO:HB3	1:A:594:TRP:CH2	2.55	0.41
1:B:143:LEU:H	1:B:143:LEU:HD23	1.86	0.41
1:B:261:CYS:HB2	1:B:488:VAL:HB	2.02	0.41
1:B:594:TRP:CZ2	1:B:598:GLN:HG3	2.56	0.41
1:A:85:LEU:CD1	1:A:94:LYS:HG3	2.50	0.41
1:B:309:LYS:HD2	1:B:328:TRP:CH2	2.56	0.41
1:A:457:GLU:HG3	1:A:513:ILE:N	2.36	0.40
1:A:414:THR:HG21	1:A:542:CYS:O	2.22	0.40
1:B:126:ILE:HA	1:B:129:THR:HG22	2.02	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:H:1:NAG:H61	4:H:2:NAG:N2	2.36	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	594/597 (100%)	578 (97%)	16 (3%)	0	100	100
1	B	594/597 (100%)	579 (98%)	15 (2%)	0	100	100
2	E	189/217 (87%)	172 (91%)	17 (9%)	0	100	100
2	F	190/217 (88%)	168 (88%)	22 (12%)	0	100	100
All	All	1567/1628 (96%)	1497 (96%)	70 (4%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains [i](#)

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	526/527 (100%)	509 (97%)	17 (3%)	39	65
1	B	526/527 (100%)	512 (97%)	14 (3%)	44	71
2	E	167/190 (88%)	153 (92%)	14 (8%)	11	22
2	F	167/190 (88%)	158 (95%)	9 (5%)	22	44

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
All	All	1386/1434 (97%)	1332 (96%)	54 (4%)	32 58

All (54) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	73	LEU
1	A	79	LEU
1	A	85	LEU
1	A	87	GLU
1	A	93	VAL
1	A	129	THR
1	A	143	LEU
1	A	190	MET
1	A	198	ASP
1	A	381	TYR
1	A	410	LEU
1	A	455	MET
1	A	518	ARG
1	A	558	LEU
1	A	560	LEU
1	A	581	VAL
1	A	595	LEU
1	B	82	MET
1	B	85	LEU
1	B	87	GLU
1	B	126	ILE
1	B	143	LEU
1	B	190	MET
1	B	198	ASP
1	B	325	GLN
1	B	381	TYR
1	B	388	GLN
1	B	436	ILE
1	B	518	ARG
1	B	557	MET
1	B	585	LEU
2	E	347	PHE
2	E	368	LEU
2	E	369	TYR
2	E	370	ASN
2	E	372	THR
2	E	377	PHE

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Mol	Chain	Res	Type
2	E	383	SER
2	E	394	ASN
2	E	408	ARG
2	E	446	THR
2	E	458	LYS
2	E	468	ILE
2	E	471	GLU
2	E	514	SER
2	F	335	LEU
2	F	372	THR
2	F	377	PHE
2	F	383	SER
2	F	387	LEU
2	F	394	ASN
2	F	446	THR
2	F	514	SER
2	F	525	CYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	60	GLN
1	A	81	GLN
1	B	42	GLN
1	B	81	GLN
1	B	96	GLN
1	B	388	GLN
2	E	370	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

23 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
3	NAG	C	1	1,3	14,14,15	0.32	0	17,19,21	0.53	0
3	NAG	C	2	3	14,14,15	0.27	0	17,19,21	0.38	0
3	BMA	C	3	3	11,11,12	0.60	0	15,15,17	0.73	0
3	NAG	D	1	1,3	14,14,15	0.75	1 (7%)	17,19,21	0.88	0
3	NAG	D	2	3	14,14,15	0.24	0	17,19,21	0.51	0
3	BMA	D	3	3	11,11,12	0.80	0	15,15,17	0.98	0
4	NAG	G	1	1,4	14,14,15	0.57	0	17,19,21	1.04	1 (5%)
4	NAG	G	2	4	14,14,15	0.93	1 (7%)	17,19,21	1.17	1 (5%)
4	NAG	H	1	1,4	14,14,15	0.31	0	17,19,21	0.47	0
4	NAG	H	2	4	14,14,15	0.42	0	17,19,21	1.26	2 (11%)
4	NAG	I	1	1,4	14,14,15	0.49	0	17,19,21	0.52	0
4	NAG	I	2	4	14,14,15	0.25	0	17,19,21	0.35	0
5	NAG	J	1	1,5	14,14,15	0.32	0	17,19,21	0.50	0
5	NAG	J	2	5	14,14,15	0.22	0	17,19,21	0.40	0
5	BMA	J	3	5	11,11,12	0.82	0	15,15,17	1.07	1 (6%)
5	NAG	J	4	5	14,14,15	1.46	2 (14%)	17,19,21	1.09	1 (5%)
4	NAG	K	1	1,4	14,14,15	0.25	0	17,19,21	0.48	0
4	NAG	K	2	4	14,14,15	0.28	0	17,19,21	0.42	0
3	NAG	L	1	3,2	14,14,15	0.24	0	17,19,21	0.47	0
3	NAG	L	2	3	14,14,15	0.21	0	17,19,21	0.43	0
3	BMA	L	3	3	11,11,12	1.22	1 (9%)	15,15,17	1.20	2 (13%)
4	NAG	M	1	2,4	14,14,15	0.50	0	17,19,21	0.72	0
4	NAG	M	2	4	14,14,15	0.75	1 (7%)	17,19,21	1.24	1 (5%)

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	C	1	1,3	-	2/6/23/26	0/1/1/1
3	NAG	C	2	3	-	2/6/23/26	0/1/1/1
3	BMA	C	3	3	-	1/2/19/22	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	NAG	D	1	1,3	-	4/6/23/26	0/1/1/1
3	NAG	D	2	3	-	2/6/23/26	0/1/1/1
3	BMA	D	3	3	-	0/2/19/22	0/1/1/1
4	NAG	G	1	1,4	-	2/6/23/26	0/1/1/1
4	NAG	G	2	4	-	3/6/23/26	0/1/1/1
4	NAG	H	1	1,4	-	2/6/23/26	0/1/1/1
4	NAG	H	2	4	-	2/6/23/26	0/1/1/1
4	NAG	I	1	1,4	-	2/6/23/26	0/1/1/1
4	NAG	I	2	4	-	2/6/23/26	0/1/1/1
5	NAG	J	1	1,5	-	2/6/23/26	0/1/1/1
5	NAG	J	2	5	-	2/6/23/26	0/1/1/1
5	BMA	J	3	5	-	0/2/19/22	0/1/1/1
5	NAG	J	4	5	-	4/6/23/26	0/1/1/1
4	NAG	K	1	1,4	-	0/6/23/26	0/1/1/1
4	NAG	K	2	4	-	2/6/23/26	0/1/1/1
3	NAG	L	1	3,2	-	2/6/23/26	0/1/1/1
3	NAG	L	2	3	-	2/6/23/26	0/1/1/1
3	BMA	L	3	3	-	1/2/19/22	0/1/1/1
4	NAG	M	1	2,4	-	0/6/23/26	0/1/1/1
4	NAG	M	2	4	-	3/6/23/26	0/1/1/1

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	J	4	NAG	O5-C1	4.60	1.51	1.43
3	L	3	BMA	C1-C2	3.37	1.59	1.52
4	G	2	NAG	O5-C1	3.15	1.48	1.43
5	J	4	NAG	C1-C2	2.81	1.56	1.52
3	D	1	NAG	O5-C1	-2.66	1.39	1.43
4	M	2	NAG	C1-C2	2.44	1.56	1.52

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	G	2	NAG	C1-O5-C5	4.39	118.14	112.19
4	H	2	NAG	C1-O5-C5	4.31	118.03	112.19
4	M	2	NAG	C2-N2-C7	4.26	128.96	122.90
5	J	4	NAG	C1-O5-C5	4.25	117.95	112.19
4	G	1	NAG	C1-O5-C5	3.49	116.92	112.19

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	L	3	BMA	O2-C2-C3	-2.64	104.86	110.14
5	J	3	BMA	C1-C2-C3	2.51	112.75	109.67
3	L	3	BMA	O2-C2-C1	2.05	113.34	109.15
4	H	2	NAG	C3-C4-C5	2.02	113.83	110.24

There are no chirality outliers.

All (42) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	D	1	NAG	C4-C5-C6-O6
5	J	2	NAG	O5-C5-C6-O6
3	D	1	NAG	O5-C5-C6-O6
4	H	2	NAG	O5-C5-C6-O6
4	I	2	NAG	O5-C5-C6-O6
3	D	2	NAG	O5-C5-C6-O6
4	G	2	NAG	O5-C5-C6-O6
4	H	1	NAG	O5-C5-C6-O6
4	I	1	NAG	O5-C5-C6-O6
5	J	2	NAG	C4-C5-C6-O6
3	L	1	NAG	O5-C5-C6-O6
4	H	2	NAG	C4-C5-C6-O6
5	J	4	NAG	O5-C5-C6-O6
3	C	1	NAG	C4-C5-C6-O6
4	H	1	NAG	C4-C5-C6-O6
4	G	1	NAG	O5-C5-C6-O6
4	M	2	NAG	C8-C7-N2-C2
4	M	2	NAG	O7-C7-N2-C2
5	J	4	NAG	C8-C7-N2-C2
5	J	4	NAG	O7-C7-N2-C2
3	D	1	NAG	C8-C7-N2-C2
3	D	1	NAG	O7-C7-N2-C2
4	K	2	NAG	O5-C5-C6-O6
3	C	1	NAG	O5-C5-C6-O6
3	D	2	NAG	C4-C5-C6-O6
5	J	4	NAG	C4-C5-C6-O6
3	L	2	NAG	O5-C5-C6-O6
3	L	1	NAG	C4-C5-C6-O6
3	L	2	NAG	C4-C5-C6-O6
4	G	1	NAG	C4-C5-C6-O6
4	I	2	NAG	C4-C5-C6-O6
4	I	1	NAG	C4-C5-C6-O6
4	G	2	NAG	C4-C5-C6-O6

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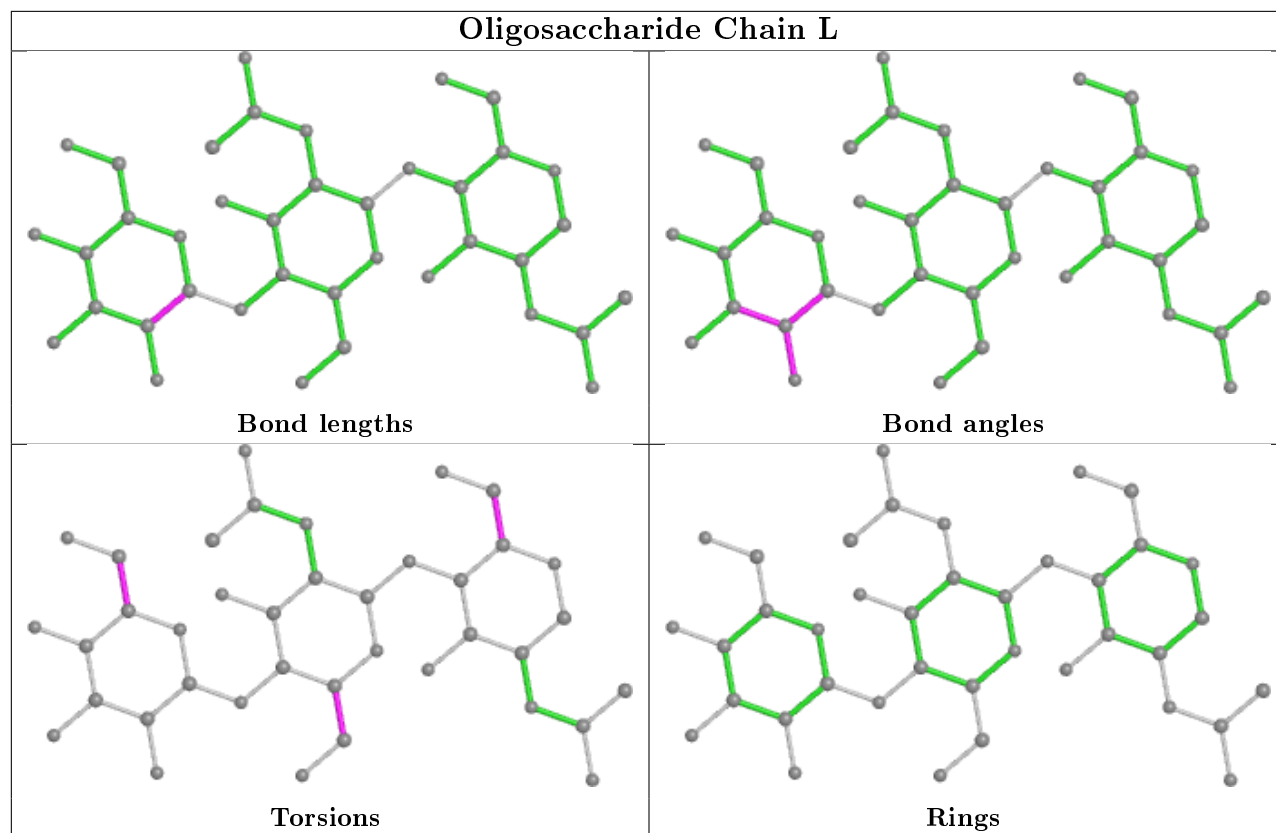
Mol	Chain	Res	Type	Atoms
3	L	3	BMA	O5-C5-C6-O6
3	C	3	BMA	O5-C5-C6-O6
3	C	2	NAG	C4-C5-C6-O6
3	C	2	NAG	O5-C5-C6-O6
4	G	2	NAG	C3-C2-N2-C7
4	M	2	NAG	C3-C2-N2-C7
5	J	1	NAG	C4-C5-C6-O6
4	K	2	NAG	C4-C5-C6-O6
5	J	1	NAG	O5-C5-C6-O6

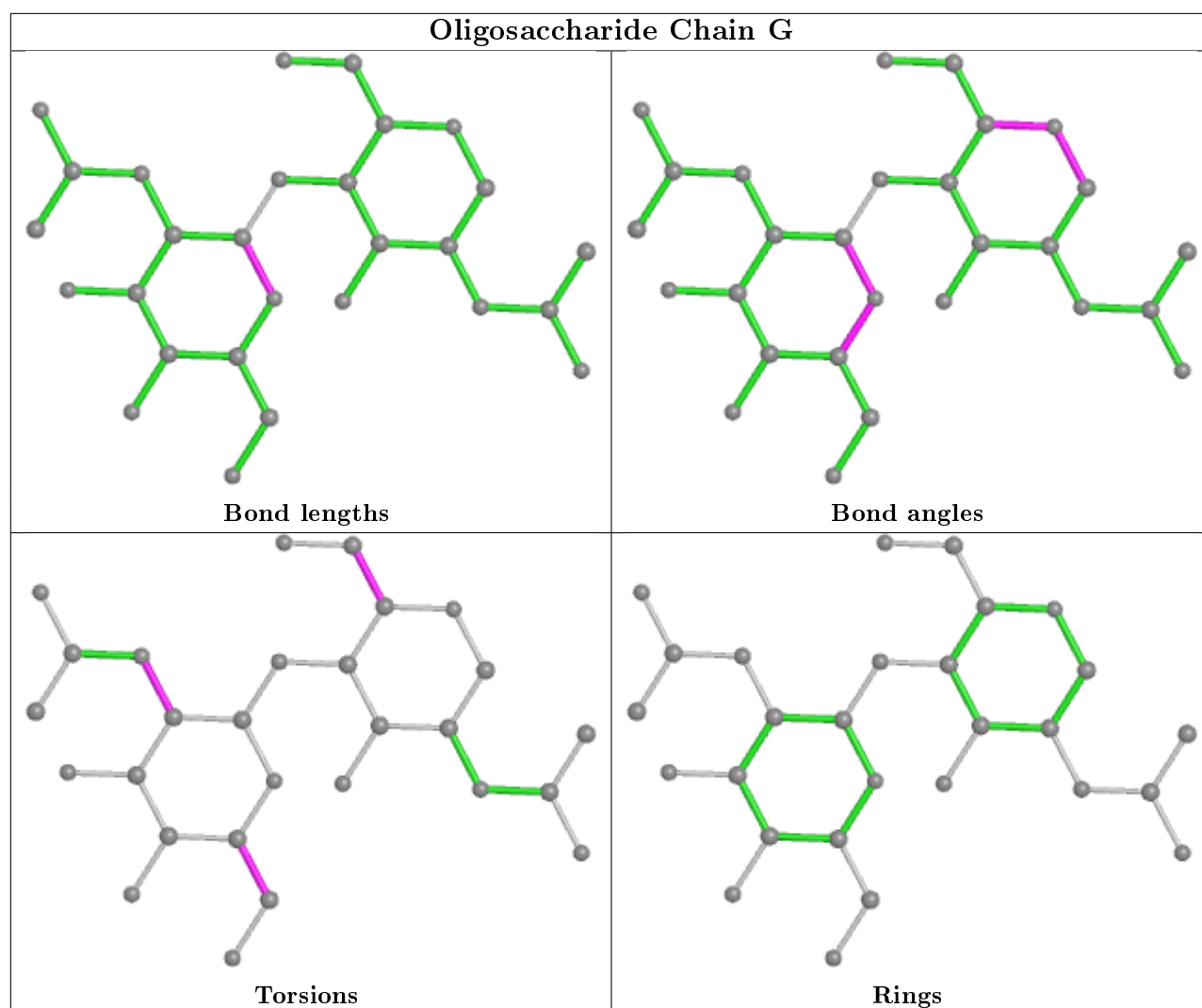
There are no ring outliers.

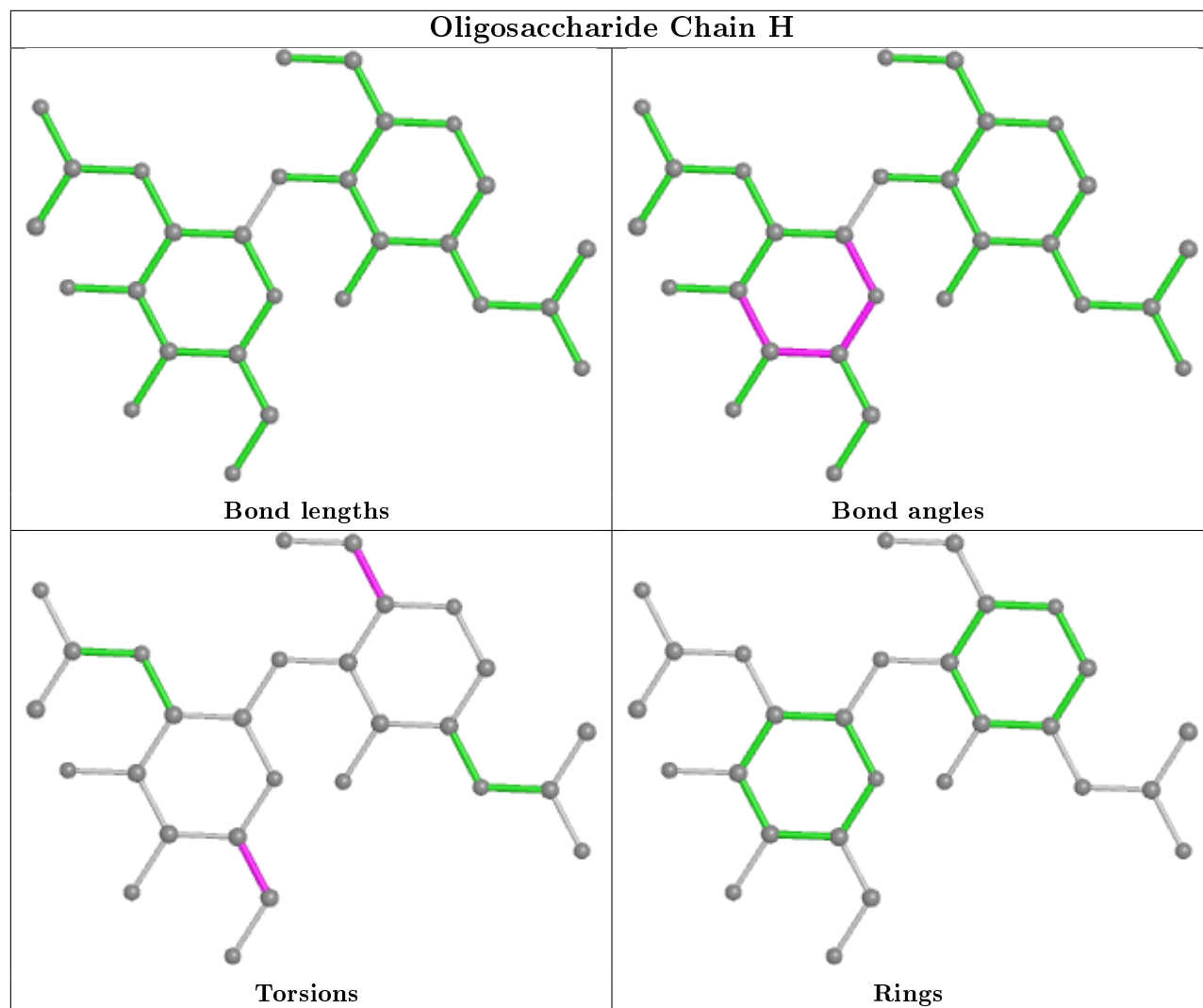
7 monomers are involved in 8 short contacts:

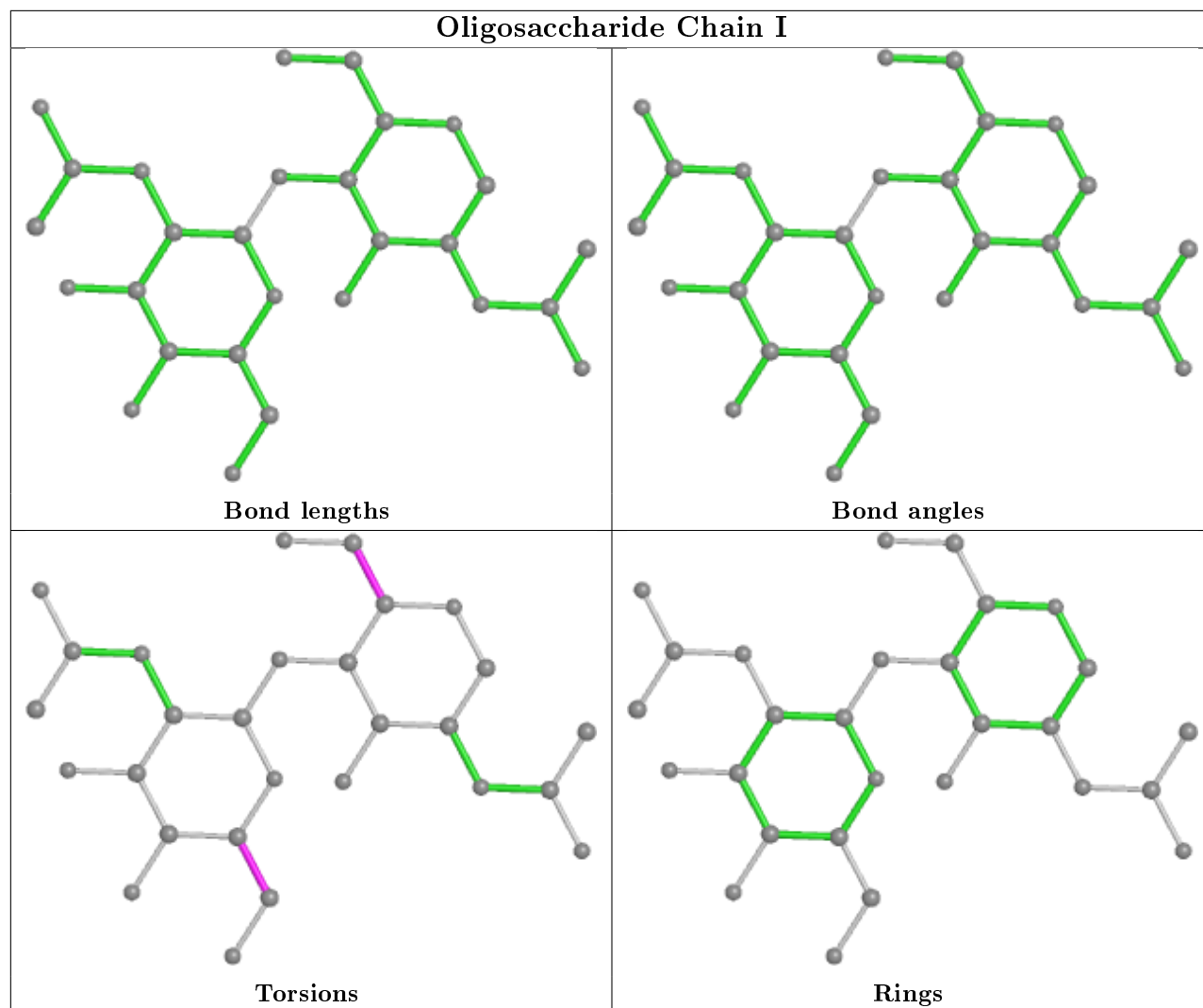
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	L	3	BMA	1	0
4	H	2	NAG	4	0
3	D	2	NAG	1	0
3	L	1	NAG	1	0
3	L	2	NAG	1	0
4	H	1	NAG	4	0
4	M	2	NAG	1	0

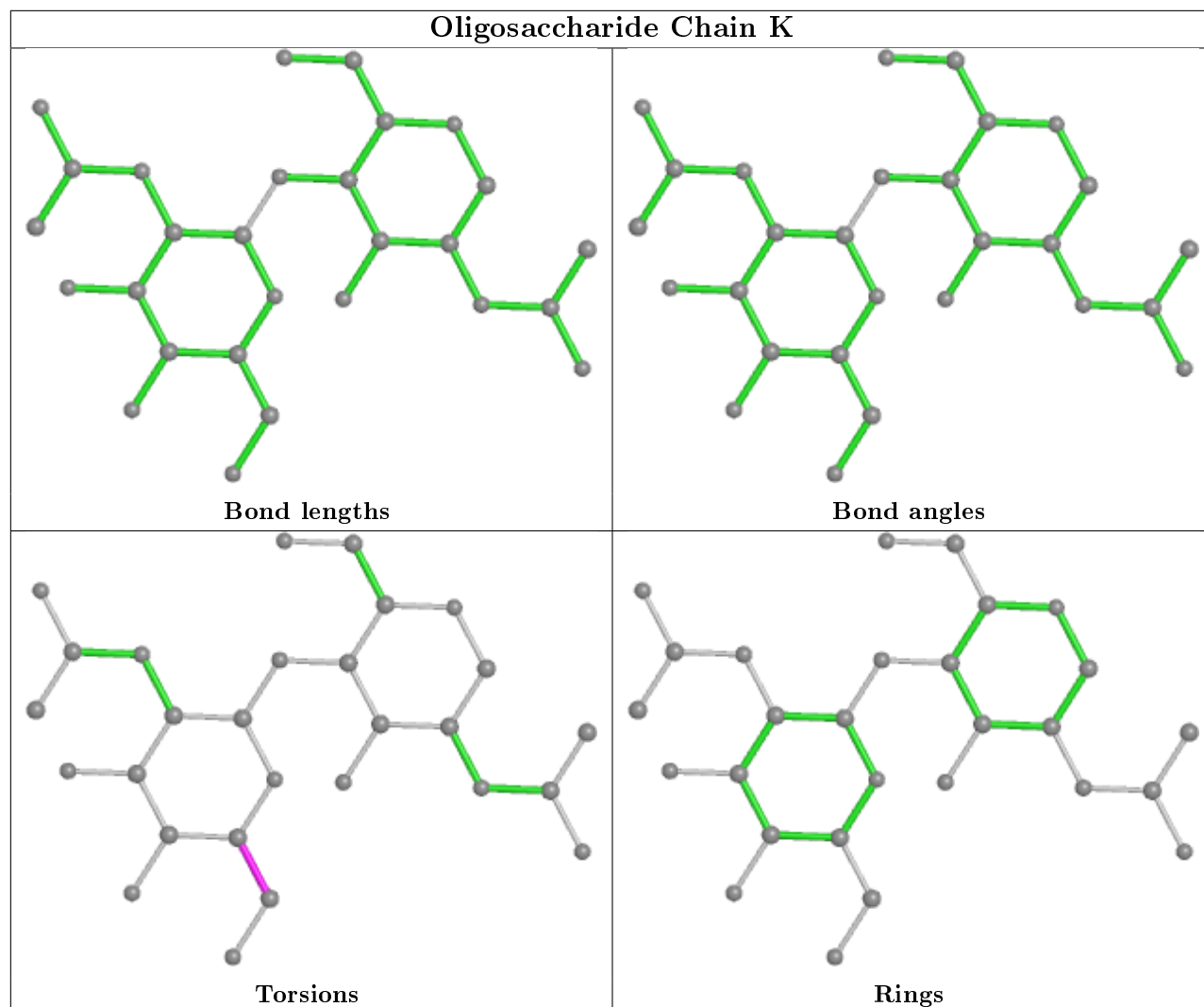
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.

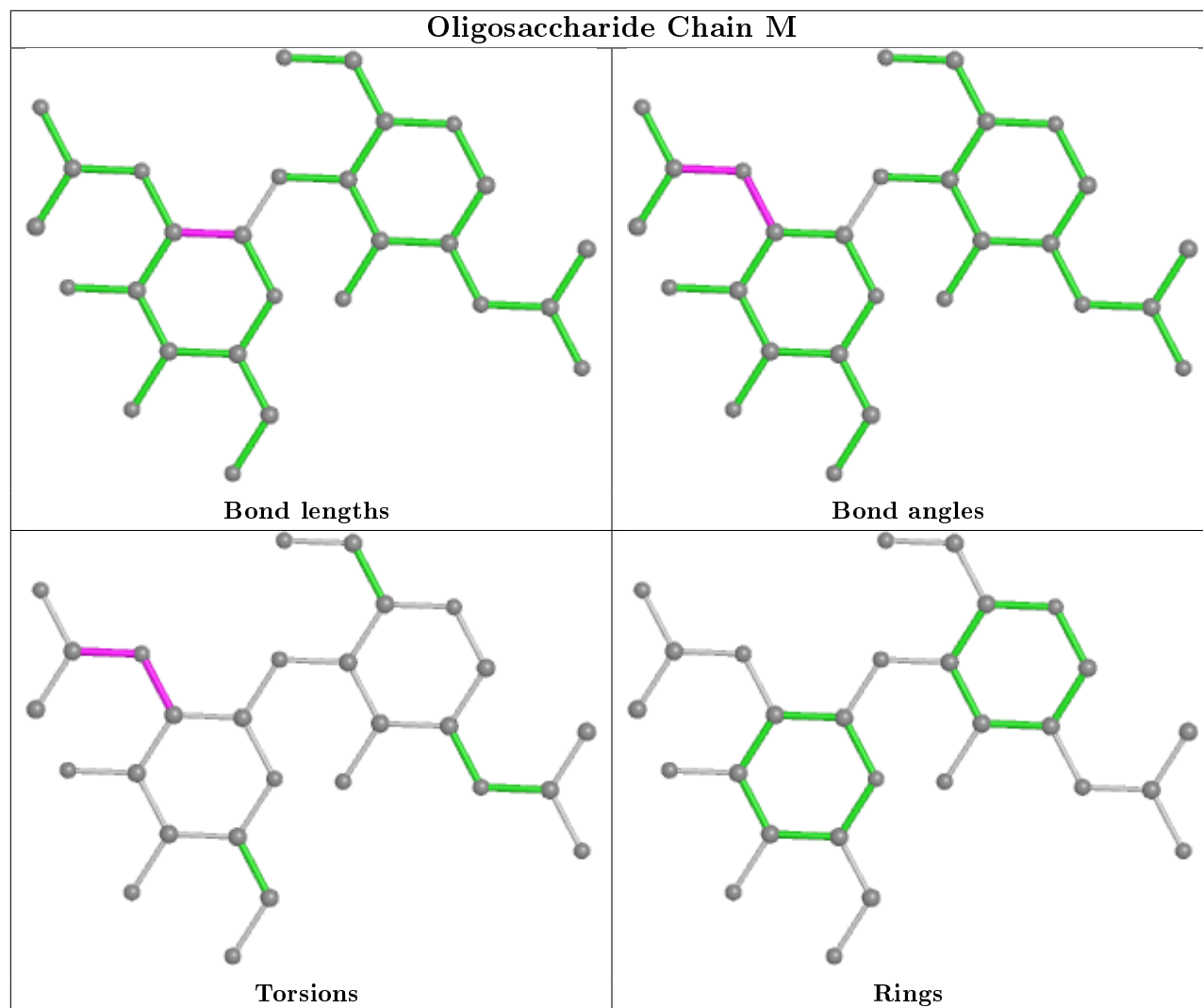


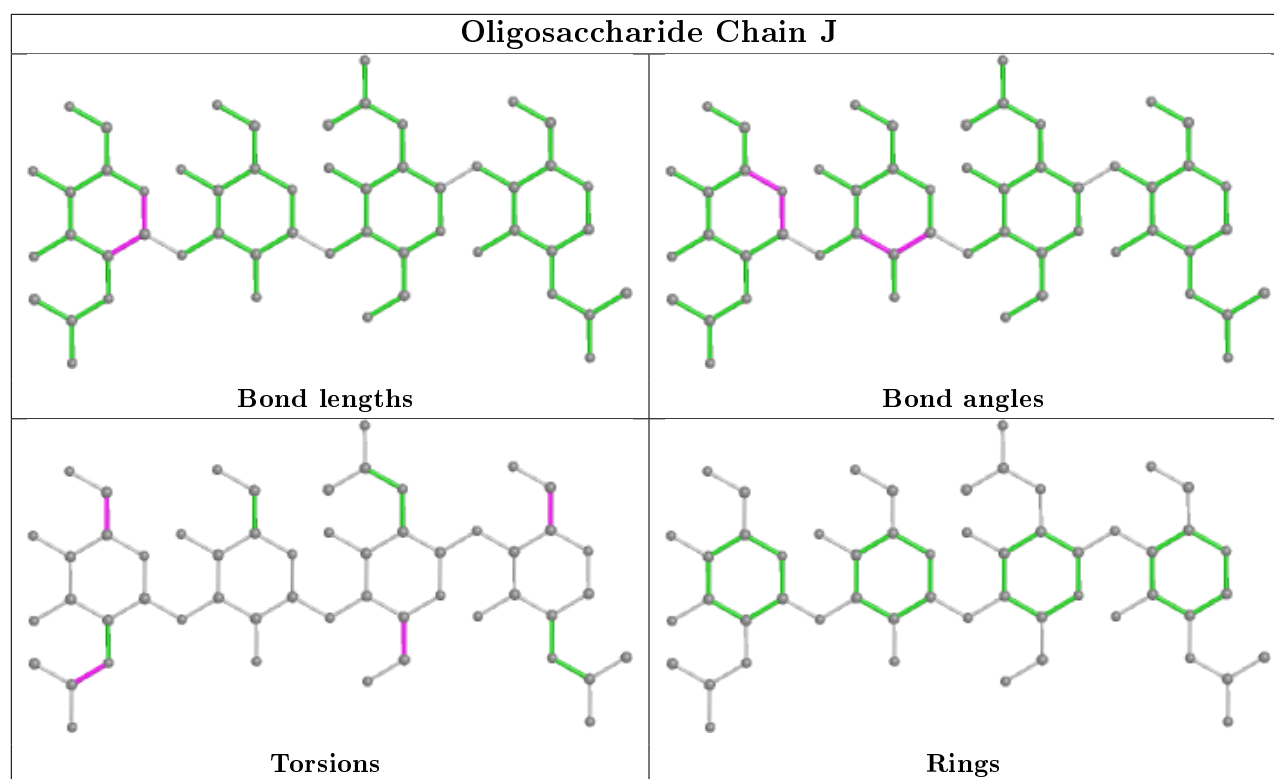












5.6 Ligand geometry [i](#)

Of 11 ligands modelled in this entry, 4 are monoatomic - leaving 7 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	$\# Z > 2$
9	NAG	A	714	1	14,14,15	0.32	0	17,19,21	0.51	0
9	NAG	B	714	1	14,14,15	0.24	0	17,19,21	0.59	0
8	EDO	B	705	-	3,3,3	0.47	0	2,2,2	0.31	0
8	EDO	A	706	-	3,3,3	0.46	0	2,2,2	0.31	0
8	EDO	B	706	-	3,3,3	0.46	0	2,2,2	0.33	0
9	NAG	B	715	1	14,14,15	0.35	0	17,19,21	0.54	0
8	EDO	B	707	-	3,3,3	0.46	0	2,2,2	0.32	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
9	NAG	A	714	1	-	0/6/23/26	0/1/1/1
9	NAG	B	714	1	-	2/6/23/26	0/1/1/1
8	EDO	B	705	-	-	0/1/1/1	-
8	EDO	A	706	-	-	0/1/1/1	-
8	EDO	B	706	-	-	1/1/1/1	-
9	NAG	B	715	1	-	1/6/23/26	0/1/1/1
8	EDO	B	707	-	-	0/1/1/1	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (4) torsion outliers are listed below:

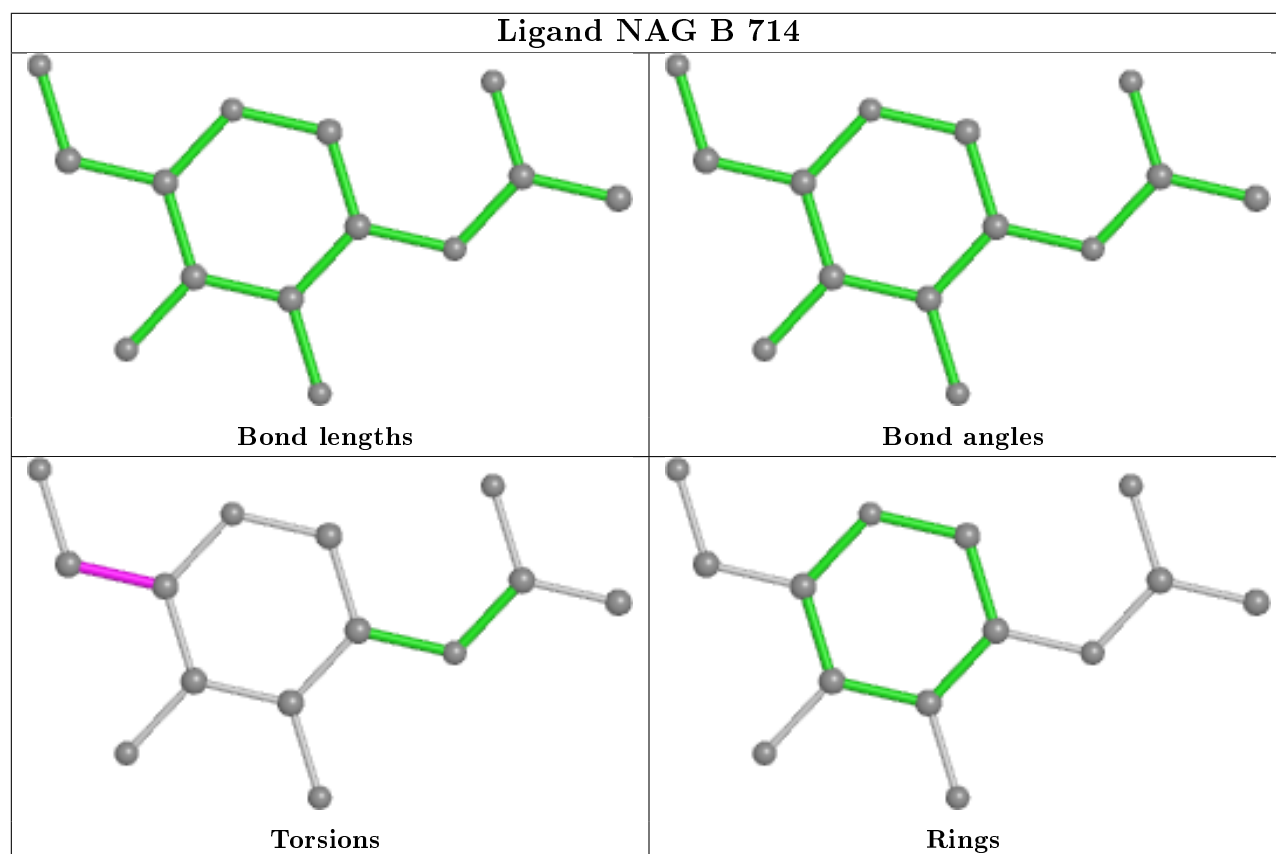
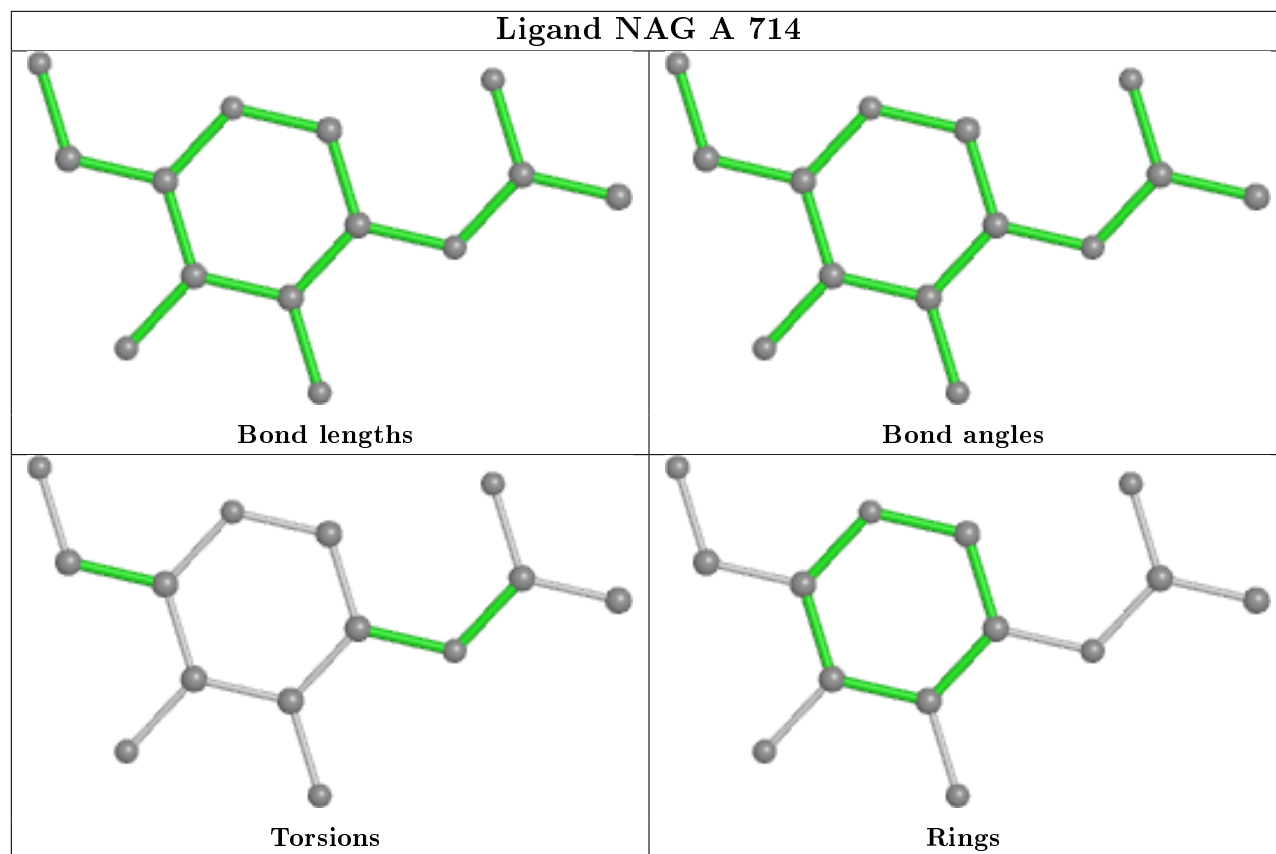
Mol	Chain	Res	Type	Atoms
9	B	714	NAG	O5-C5-C6-O6
9	B	714	NAG	C4-C5-C6-O6
9	B	715	NAG	C4-C5-C6-O6
8	B	706	EDO	O1-C1-C2-O2

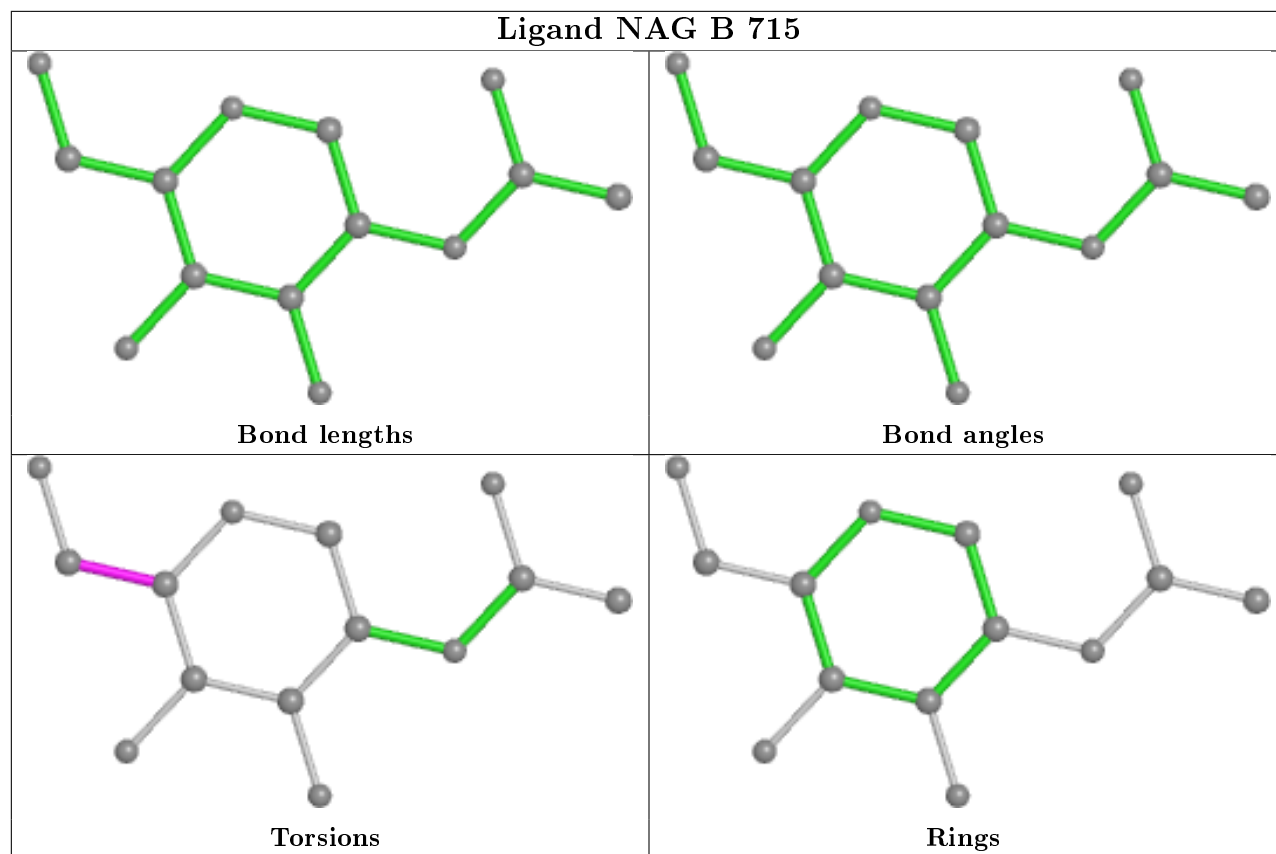
There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
9	B	714	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 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	596/597 (99%)	0.24	6 (1%) 82 82	64, 102, 148, 197	0
1	B	596/597 (99%)	0.26	7 (1%) 79 79	62, 100, 154, 244	0
2	E	193/217 (88%)	0.16	4 (2%) 63 63	78, 107, 168, 202	0
2	F	194/217 (89%)	0.28	6 (3%) 49 48	80, 112, 185, 233	0
All	All	1579/1628 (96%)	0.24	23 (1%) 73 74	62, 103, 160, 244	0

All (23) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	F	338	PHE	4.2
2	F	483	VAL	3.8
2	F	387	LEU	3.6
2	E	387	LEU	3.3
1	A	424	LEU	3.0
2	F	367	VAL	2.7
1	B	603	PHE	2.6
2	F	369	TYR	2.6
1	A	539	LEU	2.6
1	A	529	LEU	2.6
1	A	423	LEU	2.5
1	A	60	GLN	2.5
1	B	136	ASP	2.4
2	E	368	LEU	2.4
1	B	108	LEU	2.3
1	B	158	TYR	2.3
2	E	518	LEU	2.2
1	B	256	ILE	2.2
1	B	119	ILE	2.2
1	A	528	ALA	2.2
2	F	395	VAL	2.1

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Mol	Chain	Res	Type	RSRZ
1	B	84	PRO	2.0
2	E	370	ASN	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](#)

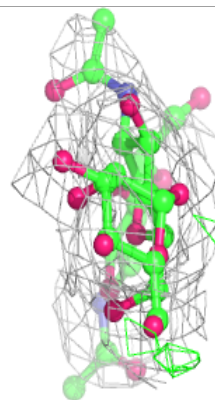
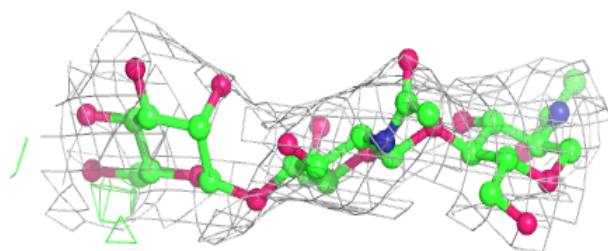
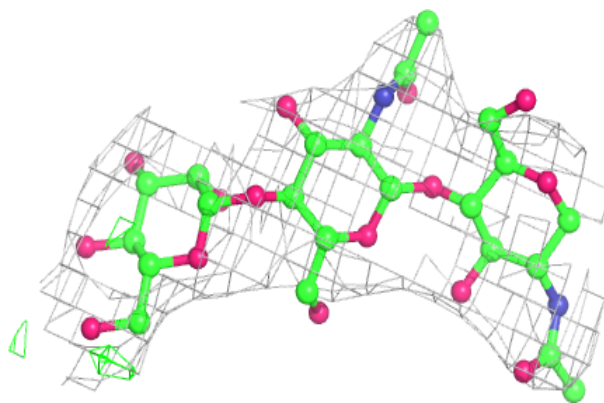
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q<0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	BMA	D	3	11/12	0.58	0.18	158,171,172,172	0
3	BMA	C	3	11/12	0.65	0.17	154,164,164,165	0
4	NAG	H	2	14/15	0.81	0.17	147,156,162,162	0
3	BMA	L	3	11/12	0.82	0.15	142,151,153,154	0
4	NAG	K	2	14/15	0.82	0.15	138,151,157,159	0
4	NAG	G	2	14/15	0.83	0.21	158,164,168,170	0
4	NAG	M	2	14/15	0.85	0.15	153,161,165,166	0
4	NAG	I	2	14/15	0.85	0.16	142,148,151,152	0
4	NAG	H	1	14/15	0.87	0.10	110,123,135,146	0
3	NAG	C	2	14/15	0.88	0.17	143,151,157,163	0
3	NAG	L	2	14/15	0.88	0.17	140,154,162,163	0
3	NAG	D	1	14/15	0.89	0.22	117,138,154,156	0
5	NAG	J	4	14/15	0.90	0.27	125,136,150,153	0
5	BMA	J	3	11/12	0.90	0.11	130,138,142,143	0
4	NAG	G	1	14/15	0.91	0.13	107,129,142,155	0
4	NAG	I	1	14/15	0.93	0.14	108,129,141,147	0
3	NAG	L	1	14/15	0.93	0.14	109,125,148,150	0
4	NAG	M	1	14/15	0.93	0.13	109,122,139,146	0
3	NAG	C	1	14/15	0.93	0.13	103,121,127,133	0
3	NAG	D	2	14/15	0.94	0.13	161,163,166,170	0
4	NAG	K	1	14/15	0.94	0.11	98,108,125,138	0
5	NAG	J	1	14/15	0.94	0.14	76,99,107,113	0
5	NAG	J	2	14/15	0.95	0.12	110,116,120,126	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.

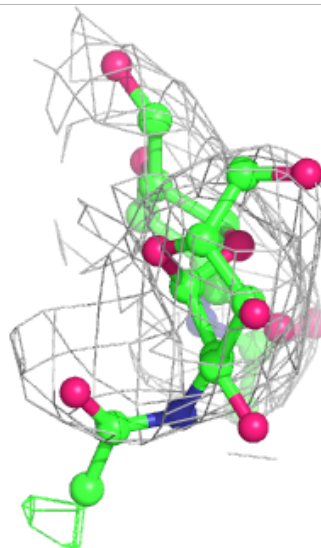
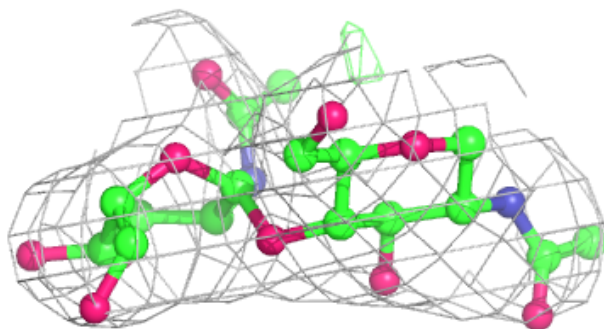
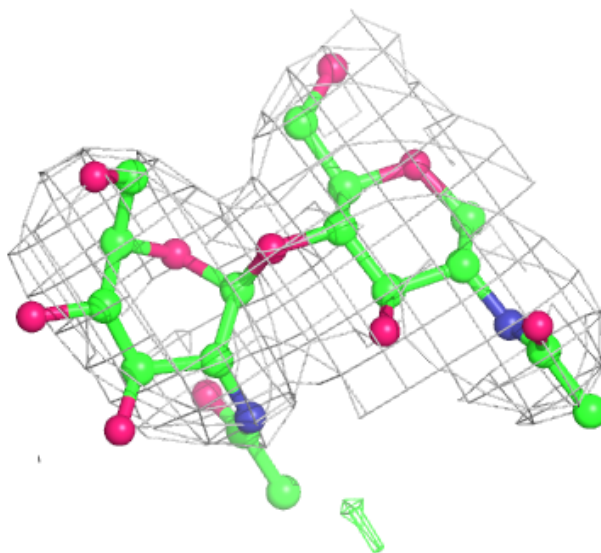
Electron density around Chain L:

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



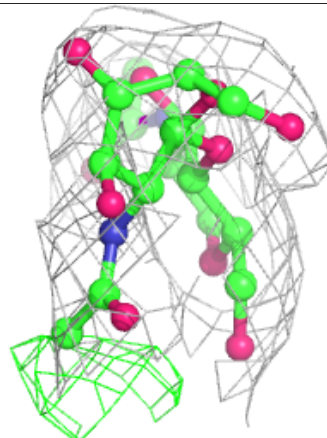
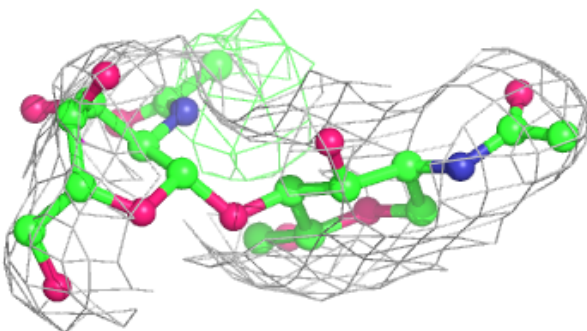
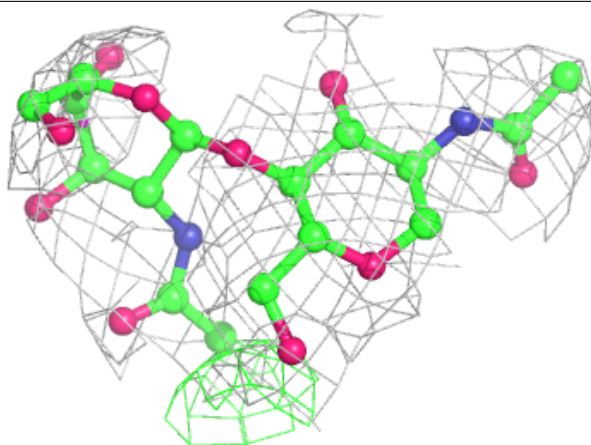
Electron density around Chain G:

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

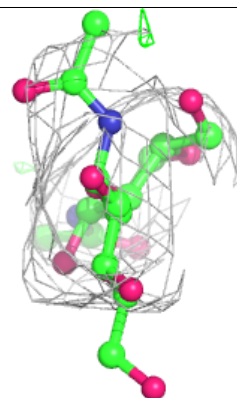
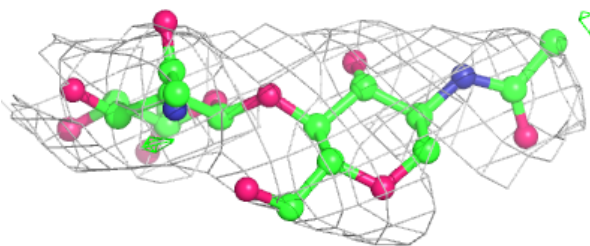
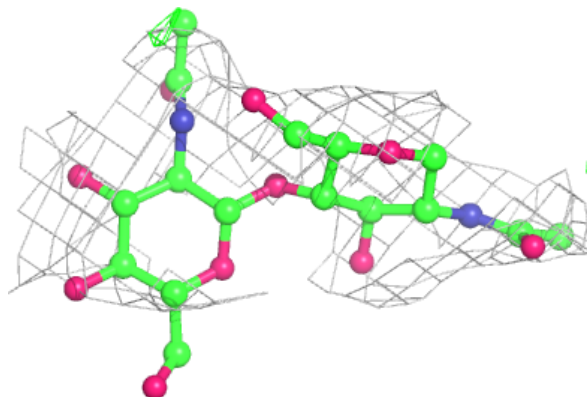


Electron density around Chain H:

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

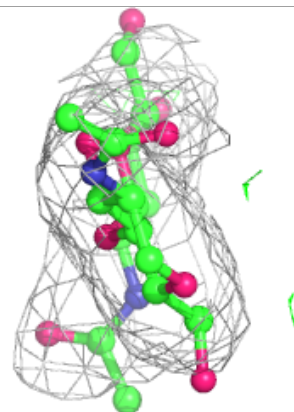
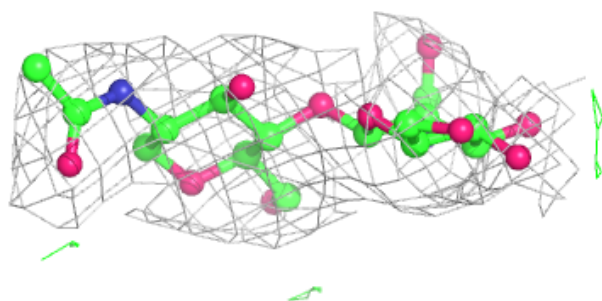
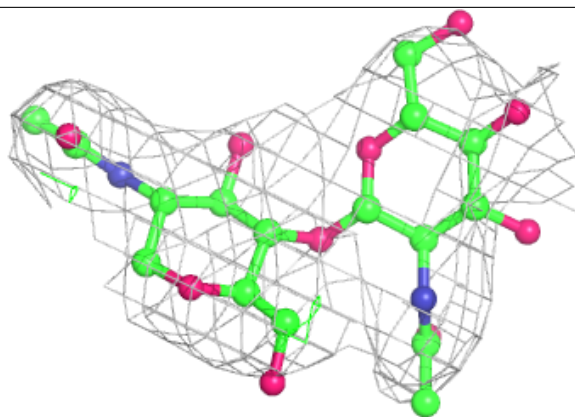
**Electron density around Chain I:**

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

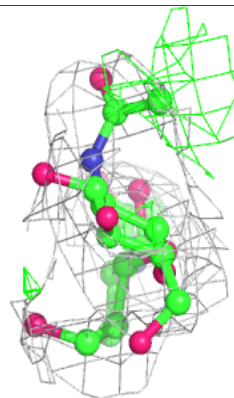
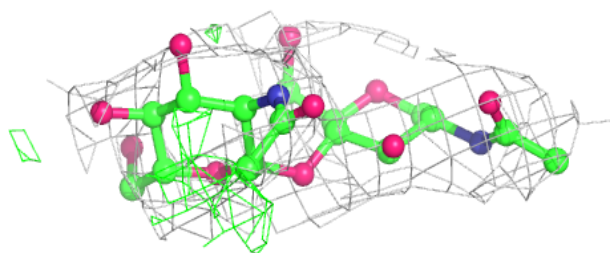
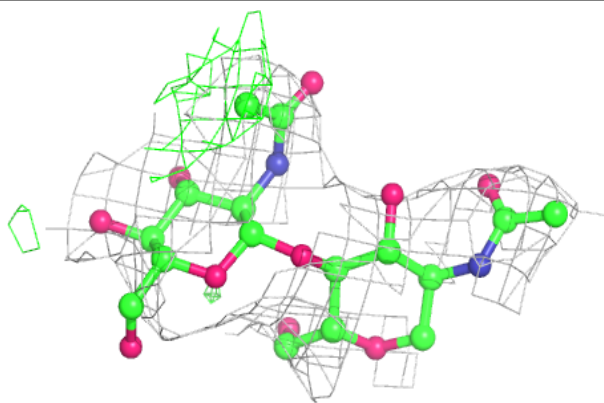


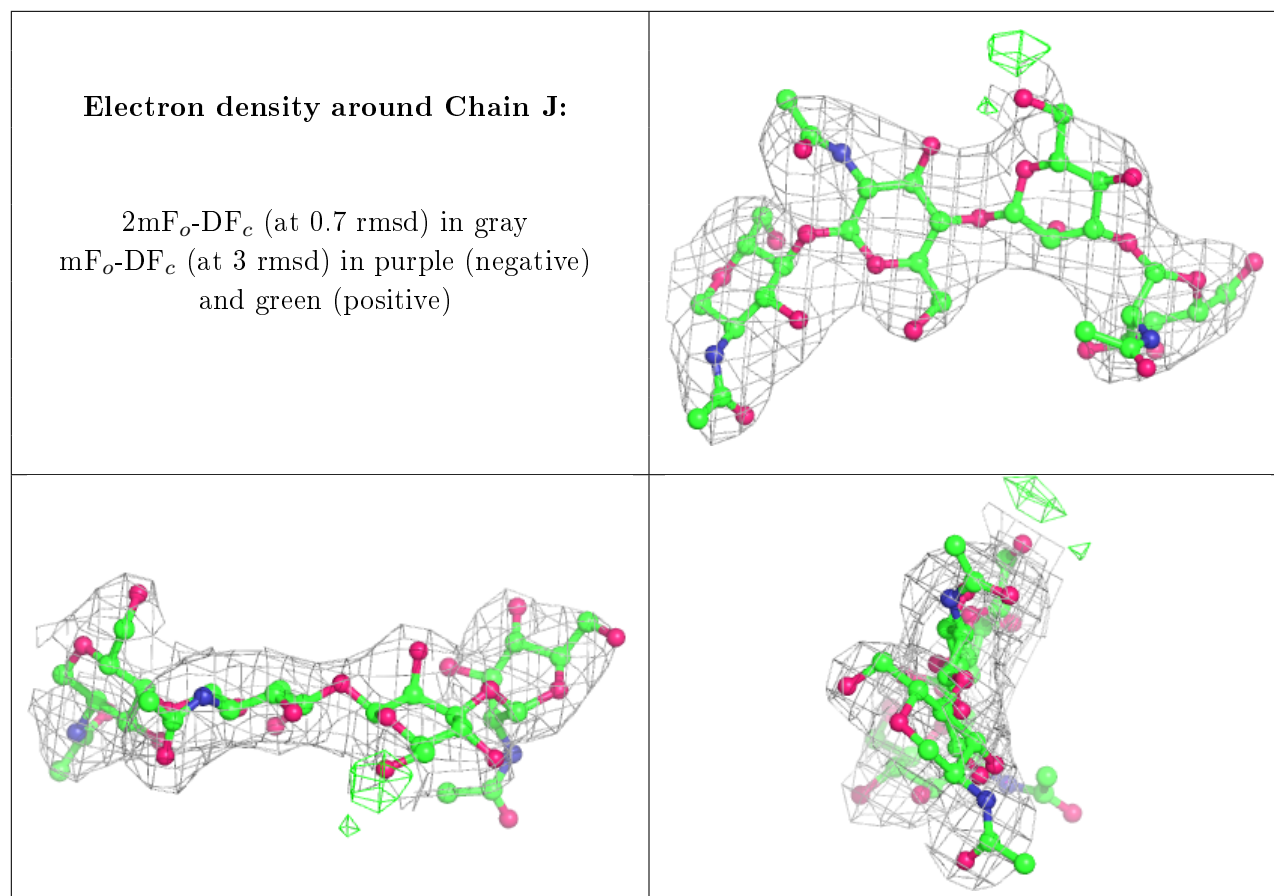
Electron density around Chain K:

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

**Electron density around Chain M:**

$2mF_o-DF_c$ (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.

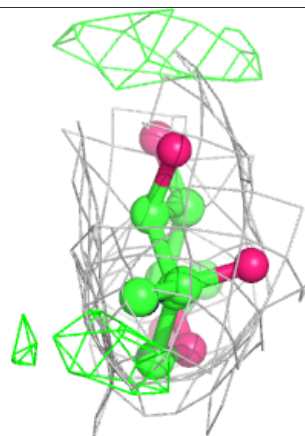
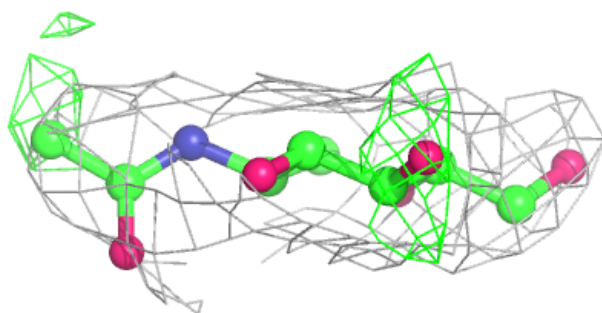
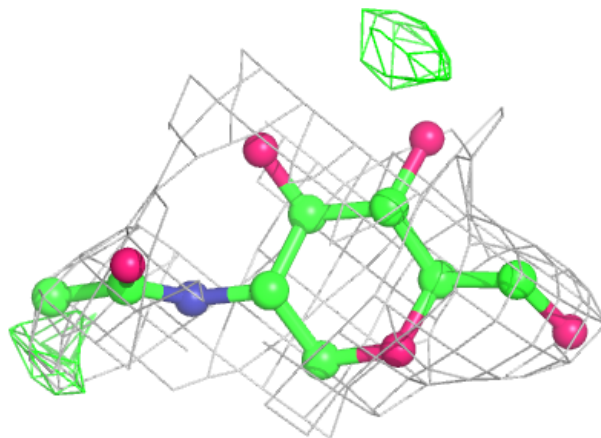
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
6	ZN	B	703	1/1	0.24	0.19	196,196,196,196	0
9	NAG	A	714	14/15	0.71	0.18	120,133,136,137	0
8	EDO	A	706	4/4	0.80	0.20	105,109,111,113	0
6	ZN	A	704	1/1	0.81	0.16	135,135,135,135	0
9	NAG	B	714	14/15	0.83	0.29	124,132,141,142	0
8	EDO	B	705	4/4	0.84	0.33	90,92,93,93	0
8	EDO	B	706	4/4	0.86	0.21	108,108,108,108	0
9	NAG	B	715	14/15	0.89	0.12	123,137,139,141	0
8	EDO	B	707	4/4	0.92	0.39	81,85,94,99	0
7	CL	B	704	1/1	0.96	0.15	102,102,102,102	0
7	CL	A	705	1/1	0.96	0.14	75,75,75,75	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.

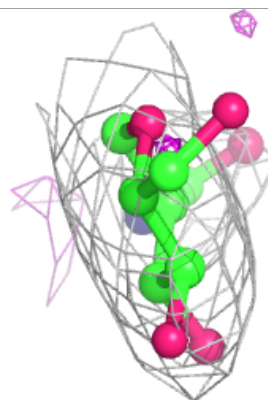
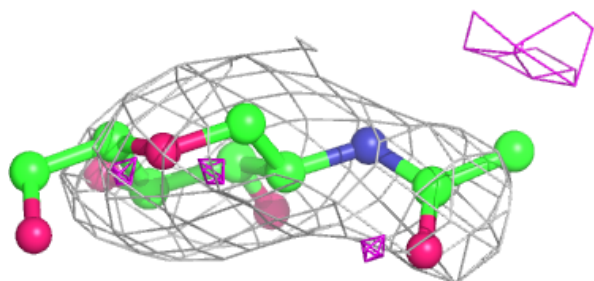
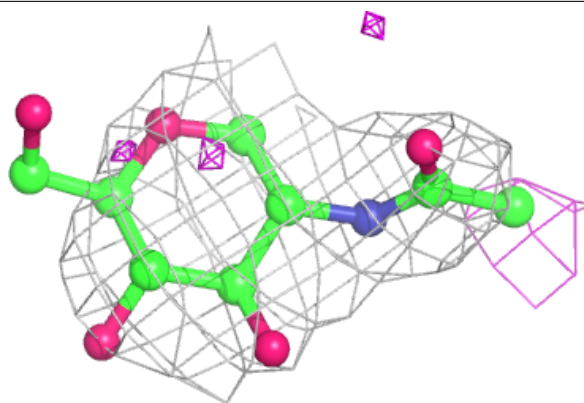
Electron density around NAG A 714:

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

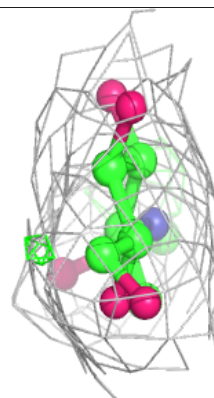
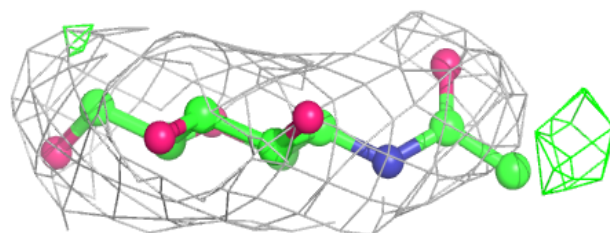
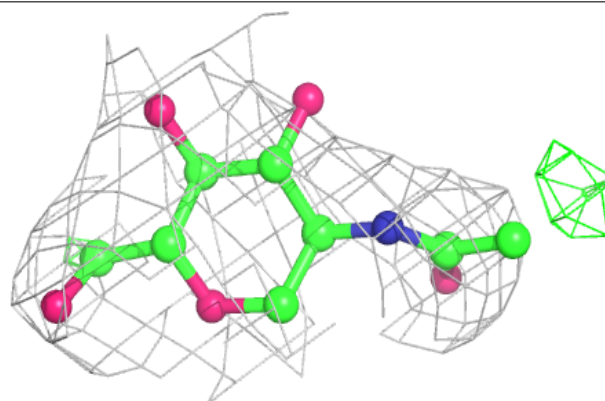


Electron density around NAG B 714:

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

**Electron density around NAG B 715:**

$2mF_o - DF_c$ (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.