



# Full wwPDB X-ray Structure Validation Report ⓘ

Aug 21, 2020 – 03:58 PM BST

PDB ID : 5LLS  
Title : Porcine dipeptidyl peptidase IV in complex with 8-(3-aminopiperidin-1-yl)-7-[(2-bromophenyl)methyl]-1,3-dimethyl-2,3,6,7-tetrahydro-1H-purine-2,6-dione  
Authors : Nar, H.; Blaesle, M.  
Deposited on : 2016-07-28  
Resolution : 2.41 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](mailto: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.

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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

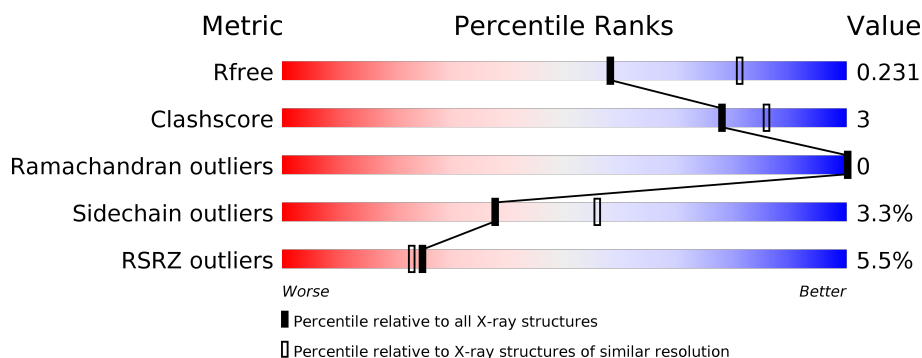
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.41 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.



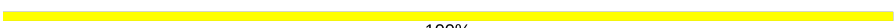


Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
$R_{free}$	130704	4647 (2.44-2.40)
Clashscore	141614	5161 (2.44-2.40)
Ramachandran outliers	138981	5073 (2.44-2.40)
Sidechain outliers	138945	5074 (2.44-2.40)
RSRZ outliers	127900	4543 (2.44-2.40)

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  $\geq 3$ , 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	766	<div> <div>4%</div> <div> <div></div> <div>85%</div> <div>10%</div> <div>5%</div> </div> </div>
1	B	766	<div> <div>4%</div> <div> <div></div> <div>85%</div> <div>10%</div> <div>5%</div> </div> </div>
1	C	766	<div> <div>6%</div> <div> <div></div> <div>85%</div> <div>10%</div> <div>5%</div> </div> </div>
1	D	766	<div> <div>6%</div> <div> <div></div> <div>86%</div> <div>9%</div> <div>5%</div> </div> </div>
2	E	2	<div> <div></div> <div> <div>50%</div> <div>50%</div> </div> </div>
2	F	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%
2	H	2	 100%
2	I	2	 50% 50%
2	J	2	 100%
2	K	2	 100%
2	L	2	 50% 50%
2	M	2	 50% 50%

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
3	NAG	D	801	-	-	-	X

## 2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 25567 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 Dipeptidyl peptidase 4.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	727	Total	C	N	O	S	0	0	0
			5960	3822	985	1130	23			
1	B	727	Total	C	N	O	S	0	0	0
			5960	3822	985	1130	23			
1	C	727	Total	C	N	O	S	0	0	0
			5960	3822	985	1130	23			
1	D	727	Total	C	N	O	S	0	0	0
			5960	3822	985	1130	23			

- Molecule 2 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
2	E	2	Total	C	N	O	0	0	0
			28	16	2	10			
2	F	2	Total	C	N	O	0	0	0
			28	16	2	10			
2	G	2	Total	C	N	O	0	0	0
			28	16	2	10			
2	H	2	Total	C	N	O	0	0	0
			28	16	2	10			
2	I	2	Total	C	N	O	0	0	0
			28	16	2	10			
2	J	2	Total	C	N	O	0	0	0
			28	16	2	10			
2	K	2	Total	C	N	O	0	0	0
			28	16	2	10			
2	L	2	Total	C	N	O	0	0	0
			28	16	2	10			

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

- Molecule 3 is 2-acetamido-2-deoxy-beta-D-glucopyranose (three-letter code: NAG) (formula:  $C_8H_{15}NO_6$ ).



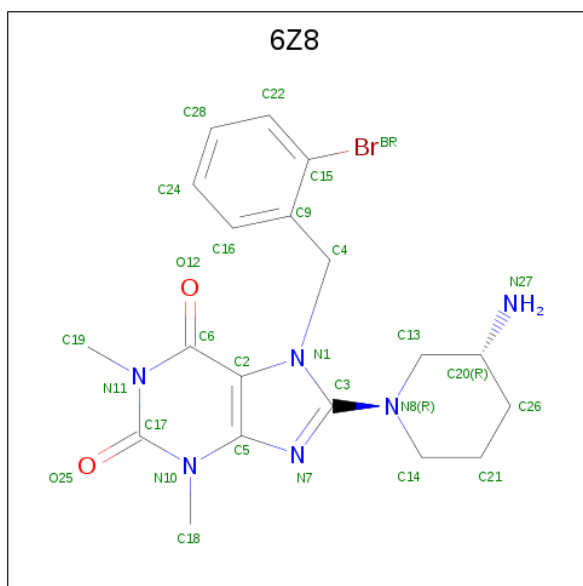
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			14	8	1	5		
3	A	1	Total	C	N	O	0	0
			14	8	1	5		
3	A	1	Total	C	N	O	0	0
			14	8	1	5		
3	A	1	Total	C	N	O	0	0
			14	8	1	5		
3	B	1	Total	C	N	O	0	0
			14	8	1	5		
3	B	1	Total	C	N	O	0	0
			14	8	1	5		
3	B	1	Total	C	N	O	0	0
			14	8	1	5		
3	C	1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

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

- Molecule 4 is 8-[(3 {R})-3-azanylpiperidin-1-yl]-7-[(2-bromophenyl)methyl]-1,3-dimethyl-purine-2,6-dione (three-letter code: 6Z8) (formula: C<sub>19</sub>H<sub>23</sub>BrN<sub>6</sub>O<sub>2</sub>).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total 28	Br 1	C 19	N 6	O 2	0	0
4	A	1	Total 28	Br 1	C 19	N 6	O 2	0	0
4	B	1	Total 28	Br 1	C 19	N 6	O 2	0	0
4	B	1	Total 28	Br 1	C 19	N 6	O 2	0	0
4	C	1	Total 28	Br 1	C 19	N 6	O 2	0	0
4	C	1	Total 28	Br 1	C 19	N 6	O 2	0	0

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	D	1	Total	Br	C	N	O	0	0
			28	1	19	6	2		
4	D	1	Total	Br	C	N	O	0	0
			28	1	19	6	2		

- Molecule 5 is SULFATE ION (three-letter code: SO4) (formula: O<sub>4</sub>S).



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

- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	254	Total	O	0	0
			254	254		
6	B	280	Total	O	0	0
			280	280		
6	C	273	Total	O	0	0
			273	273		

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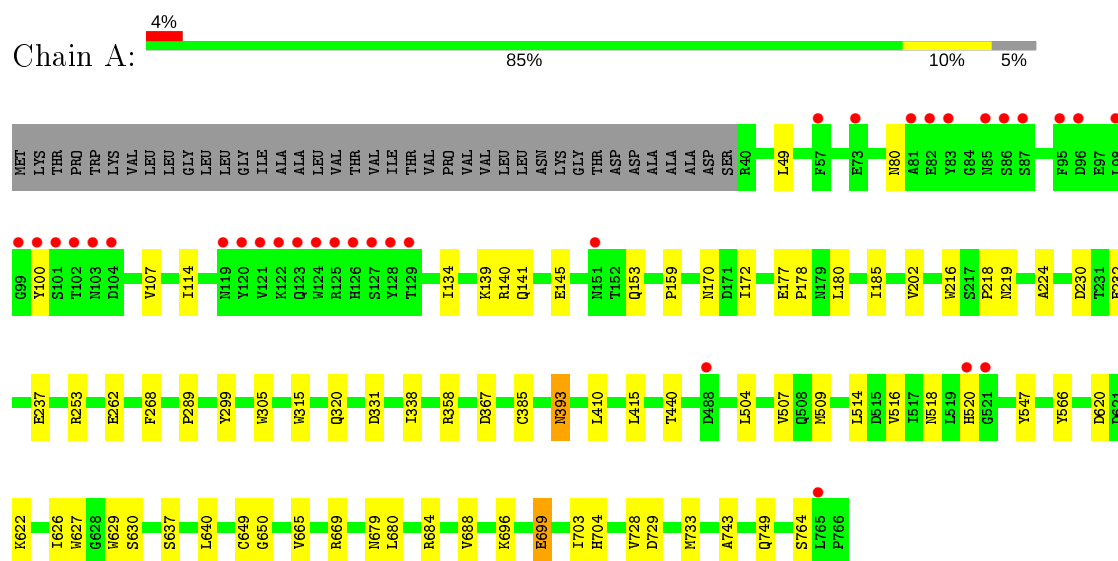
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	D	214	Total 214	O 214	0	0



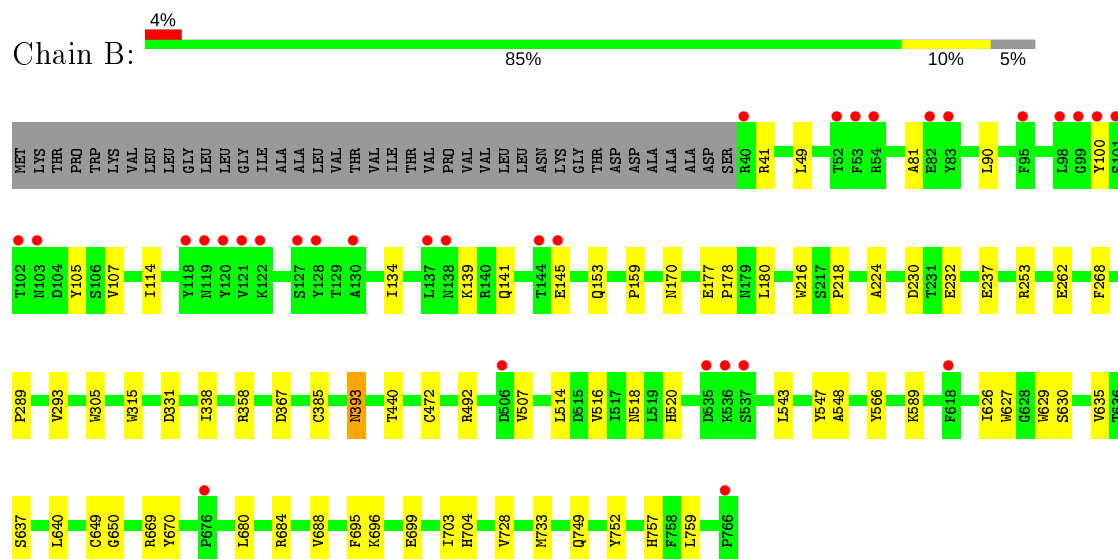
### 3 Residue-property plots [i](#)

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.

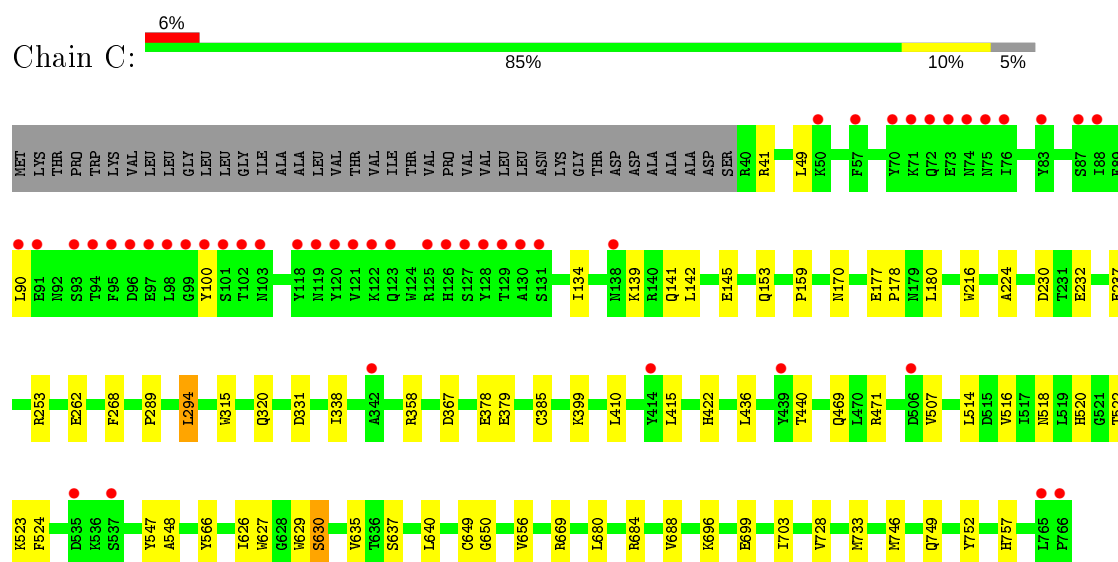
- Molecule 1: Dipeptidyl peptidase 4



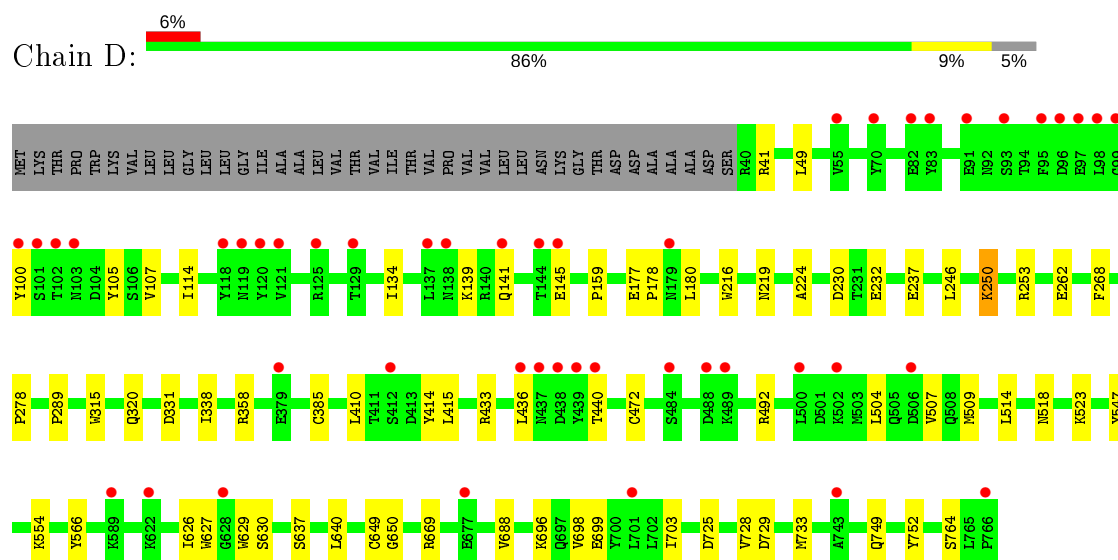
- Molecule 1: Dipeptidyl peptidase 4



- Molecule 1: Dipeptidyl peptidase 4



• Molecule 1: Dipeptidyl peptidase 4



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



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



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

Chain G:  100%

NAG1  
NAG2

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

Chain H:  100%

NAG1  
NAG2

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

Chain I:  50%  50%

NAG1  
NAG2

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

Chain J:  100%

NAG1  
NAG2

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

Chain K:  100%

NAG1  
NAG2

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

Chain L:  50%  50%

NAG1  
NAG2

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

Chain M:  50%  50%



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	61.90Å 117.68Å 133.12Å 112.70° 94.60° 91.20°	Depositor
Resolution (Å)	35.28 – 2.41 35.10 – 2.41	Depositor EDS
% Data completeness (in resolution range)	80.3 (35.28-2.41) 80.5 (35.10-2.41)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.70 (at 2.42Å)	Xtriage
Refinement program	BUSTER 2.11.6	Depositor
R, $R_{free}$	0.190 , 0.228 0.195 , 0.231	Depositor DCC
$R_{free}$ test set	5218 reflections (4.88%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	34.7	Xtriage
Anisotropy	0.537	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.35 , 55.7	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.93	EDS
Total number of atoms	25567	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	41.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.02% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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: 6Z8, NAG, SO4

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	0.42	0/6135	0.65	0/8345
1	B	0.43	0/6135	0.66	0/8345
1	C	0.41	0/6135	0.65	0/8345
1	D	0.40	0/6135	0.65	0/8345
All	All	0.42	0/24540	0.65	0/33380

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	5960	0	5658	34	0
1	B	5960	0	5658	36	0
1	C	5960	0	5658	37	0
1	D	5960	0	5658	32	0
2	E	28	0	25	0	0
2	F	28	0	25	0	0
2	G	28	0	25	0	0
2	H	28	0	25	0	0
2	I	28	0	25	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	J	28	0	25	0	0
2	K	28	0	25	0	0
2	L	28	0	25	0	0
2	M	28	0	25	0	0
3	A	56	0	52	1	0
3	B	56	0	52	0	0
3	C	70	0	65	0	0
3	D	28	0	26	1	0
4	A	56	0	0	2	0
4	B	56	0	0	0	0
4	C	56	0	0	2	0
4	D	56	0	0	1	0
5	A	5	0	0	0	0
5	B	5	0	0	0	0
5	C	5	0	0	0	0
5	D	5	0	0	0	0
6	A	254	0	0	1	0
6	B	280	0	0	0	0
6	C	273	0	0	2	0
6	D	214	0	0	1	0
All	All	25567	0	23052	137	0

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

All (137) 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:A:253:ARG:HH21	1:B:253:ARG:HH21	1.29	0.78
1:C:640:LEU:HD11	1:C:650:GLY:HA3	1.66	0.77
1:B:640:LEU:HD11	1:B:650:GLY:HA3	1.70	0.73
1:C:253:ARG:HH21	1:D:253:ARG:HH21	1.38	0.70
1:C:696:LYS:HG3	1:C:728:VAL:HG22	1.77	0.67
1:D:696:LYS:HG3	1:D:728:VAL:HG22	1.77	0.66
1:A:696:LYS:HG3	1:A:728:VAL:HG22	1.79	0.65
1:D:640:LEU:HD11	1:D:650:GLY:HA3	1.78	0.64
1:C:656:VAL:HG23	6:C:927:HOH:O	1.98	0.64
1:C:637:SER:HB3	1:C:688:VAL:HG11	1.80	0.62
1:A:507:VAL:HG13	1:A:509:MET:HG2	1.81	0.62
1:B:696:LYS:HG3	1:B:728:VAL:HG22	1.83	0.61
1:D:414:TYR:CD1	1:D:433:ARG:HD3	2.36	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:410:LEU:HD13	1:D:415:LEU:HD12	1.83	0.59
1:B:107:VAL:HG22	1:B:114:ILE:HD12	1.85	0.58
1:D:507:VAL:HG13	1:D:509:MET:HG2	1.85	0.57
1:C:41:ARG:HG3	1:C:507:VAL:HG22	1.87	0.57
1:A:637:SER:HB3	1:A:688:VAL:HG11	1.87	0.57
1:D:320:GLN:OE1	1:D:669:ARG:HD3	2.05	0.57
1:A:393:ASN:HD22	1:A:393:ASN:H	1.53	0.56
1:B:393:ASN:H	1:B:393:ASN:HD22	1.53	0.56
1:C:422:HIS:CE1	6:C:908:HOH:O	2.58	0.56
1:A:640:LEU:HD11	1:A:650:GLY:HA3	1.87	0.56
1:B:159:PRO:HD3	1:B:216:TRP:HB3	1.89	0.55
1:A:410:LEU:HD13	1:A:415:LEU:HD12	1.88	0.55
1:B:41:ARG:HG3	1:B:507:VAL:HG22	1.89	0.55
1:A:153:GLN:HE22	1:A:170:ASN:ND2	2.06	0.54
1:A:159:PRO:HD3	1:A:216:TRP:HB3	1.90	0.54
1:A:504:LEU:HA	1:A:507:VAL:HG12	1.88	0.54
1:B:81:ALA:HA	1:B:492:ARG:HH22	1.72	0.53
1:A:320:GLN:OE1	1:A:669:ARG:HD3	2.08	0.53
1:B:289:PRO:HB2	1:B:293:VAL:HG23	1.89	0.53
1:D:415:LEU:HB2	1:D:436:LEU:HD21	1.89	0.53
1:C:159:PRO:HD3	1:C:216:TRP:HB3	1.90	0.53
1:D:637:SER:HB3	1:D:688:VAL:HG11	1.90	0.53
1:D:649:CYS:HB3	1:D:699:GLU:HB2	1.91	0.53
1:C:649:CYS:HB3	1:C:699:GLU:HB2	1.91	0.53
1:C:320:GLN:OE1	1:C:669:ARG:HD3	2.09	0.52
1:C:49:LEU:HD22	1:C:749:GLN:HA	1.91	0.52
1:A:680:LEU:HD11	1:A:684:ARG:HH21	1.74	0.52
1:B:649:CYS:HB3	1:B:699:GLU:HB2	1.92	0.52
1:D:159:PRO:HD3	1:D:216:TRP:HB3	1.92	0.52
1:B:177:GLU:HB2	1:B:180:LEU:HB2	1.92	0.51
1:A:134:ILE:HD13	1:A:178:PRO:HB3	1.92	0.51
1:C:177:GLU:HB2	1:C:180:LEU:HB2	1.92	0.51
1:D:504:LEU:HA	1:D:507:VAL:HG12	1.92	0.51
1:D:331:ASP:HB2	1:D:338:ILE:HD12	1.93	0.51
1:C:410:LEU:HD13	1:C:415:LEU:HD12	1.93	0.51
1:A:331:ASP:HB2	1:A:338:ILE:HD12	1.93	0.51
1:D:49:LEU:HD22	1:D:749:GLN:HA	1.93	0.50
1:C:134:ILE:HD13	1:C:178:PRO:HB3	1.93	0.50
1:A:630:SER:HB2	4:A:809:6Z8:O12	2.12	0.50
1:C:153:GLN:HE22	1:C:170:ASN:HD22	1.58	0.50
1:A:649:CYS:HB3	1:A:699:GLU:HB2	1.94	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:153:GLN:HE22	1:C:170:ASN:ND2	2.10	0.50
1:D:134:ILE:HD13	1:D:178:PRO:HB3	1.94	0.50
1:D:107:VAL:HG22	1:D:114:ILE:HD12	1.93	0.49
1:B:696:LYS:CG	1:B:728:VAL:HG22	2.42	0.49
1:D:177:GLU:HB2	1:D:180:LEU:HB2	1.93	0.49
1:B:153:GLN:HE22	1:B:170:ASN:ND2	2.10	0.49
1:A:139:LYS:HG3	1:A:141:GLN:HB2	1.94	0.49
1:A:177:GLU:HB2	1:A:180:LEU:HB2	1.93	0.49
1:A:729:ASP:OD2	1:B:757:HIS:HD2	1.96	0.49
1:B:134:ILE:HD13	1:B:178:PRO:HB3	1.94	0.49
1:B:637:SER:HB3	1:B:688:VAL:HG11	1.93	0.49
1:B:49:LEU:HD22	1:B:749:GLN:HA	1.94	0.48
1:D:139:LYS:HG3	1:D:141:GLN:HB2	1.95	0.48
1:D:278:PRO:HB2	3:D:806:NAG:O7	2.14	0.48
1:A:49:LEU:HD22	1:A:749:GLN:HA	1.94	0.47
1:D:49:LEU:HD21	1:D:752:TYR:HB2	1.97	0.47
1:B:331:ASP:HB2	1:B:338:ILE:HD12	1.95	0.47
1:B:680:LEU:HD11	1:B:684:ARG:HH21	1.80	0.47
1:C:331:ASP:HB2	1:C:338:ILE:HD12	1.95	0.47
1:A:224:ALA:HB1	1:A:268:PHE:CZ	2.48	0.47
1:B:703:ILE:HG12	1:B:733:MET:HB3	1.97	0.47
1:C:139:LYS:HG3	1:C:141:GLN:HB2	1.97	0.47
1:B:224:ALA:HB1	1:B:268:PHE:CZ	2.50	0.47
1:C:289:PRO:HG2	1:C:294:LEU:HG	1.97	0.46
1:C:378:GLU:CD	1:C:378:GLU:H	2.19	0.46
1:C:469:GLN:OE1	1:C:471:ARG:HD2	2.16	0.46
1:C:522:THR:HB	1:C:524:PHE:CE2	2.50	0.46
1:B:695:PHE:HB3	1:B:728:VAL:HG11	1.98	0.45
1:B:139:LYS:HG3	1:B:141:GLN:HB2	1.98	0.45
1:C:680:LEU:HD11	1:C:684:ARG:HH21	1.81	0.44
1:C:746:MET:HG3	1:D:725:ASP:HA	1.98	0.44
1:B:105:TYR:HB2	1:B:114:ILE:HD11	2.00	0.44
1:C:703:ILE:HG12	1:C:733:MET:HB3	1.98	0.44
1:A:743:ALA:HB3	6:A:916:HOH:O	2.17	0.44
1:D:105:TYR:HB2	1:D:114:ILE:HD11	1.99	0.44
1:B:393:ASN:ND2	1:B:393:ASN:H	2.15	0.44
1:A:80:ASN:HB2	3:A:801:NAG:H82	2.00	0.44
1:B:543:LEU:HD22	1:B:759:LEU:HD11	2.00	0.44
1:C:49:LEU:HD21	1:C:752:TYR:HB2	1.98	0.44
1:B:49:LEU:HD21	1:B:752:TYR:HB2	1.99	0.44
1:D:703:ILE:HG12	1:D:733:MET:HB3	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:703:ILE:HG12	1:A:733:MET:HB3	2.00	0.43
4:C:808:6Z8:C4	4:C:808:6Z8:C13	2.96	0.43
1:B:669:ARG:HD2	1:B:670:TYR:CZ	2.53	0.43
1:B:289:PRO:HB3	1:B:315:TRP:CD2	2.54	0.43
1:C:224:ALA:HB1	1:C:268:PHE:CZ	2.53	0.43
1:D:224:ALA:HB1	1:D:268:PHE:CZ	2.53	0.43
1:A:393:ASN:H	1:A:393:ASN:ND2	2.15	0.43
1:C:237:GLU:HG2	1:C:253:ARG:HG2	2.00	0.43
1:A:237:GLU:HG2	1:A:253:ARG:HG2	2.00	0.42
1:C:630:SER:HB2	4:C:808:6Z8:O12	2.19	0.42
1:C:232:GLU:HB3	1:C:262:GLU:HG2	2.01	0.42
1:B:626:ILE:O	1:B:650:GLY:HA2	2.20	0.42
1:D:246:LEU:HD21	1:D:250:LYS:HE3	2.01	0.42
1:C:289:PRO:HB3	1:C:315:TRP:CD2	2.55	0.42
1:C:640:LEU:HD11	1:C:650:GLY:CA	2.44	0.42
4:A:809:6Z8:C13	4:A:809:6Z8:C4	2.97	0.42
1:D:232:GLU:HB3	1:D:262:GLU:HG2	2.02	0.42
1:C:415:LEU:HB2	1:C:436:LEU:HD21	2.01	0.42
1:C:548:ALA:HB3	1:C:635:VAL:HG21	2.02	0.42
1:A:218:PRO:HD3	1:A:305:TRP:HB3	2.01	0.41
1:D:289:PRO:HB3	1:D:315:TRP:CD2	2.55	0.41
1:D:626:ILE:O	1:D:650:GLY:HA2	2.21	0.41
1:A:289:PRO:HB3	1:A:315:TRP:CD2	2.55	0.41
1:A:626:ILE:O	1:A:650:GLY:HA2	2.21	0.41
1:A:107:VAL:HG22	1:A:114:ILE:HD12	2.03	0.41
1:A:299:TYR:CZ	1:A:665:VAL:HG22	2.56	0.41
1:B:232:GLU:HB3	1:B:262:GLU:HG2	2.03	0.41
1:B:759:LEU:HA	1:B:759:LEU:HD23	1.95	0.41
1:A:172:ILE:HB	1:A:185:ILE:HB	2.01	0.41
1:B:49:LEU:HB3	1:B:749:GLN:HG2	2.01	0.41
1:B:548:ALA:HB3	1:B:635:VAL:HG21	2.03	0.41
1:A:620:ASP:OD1	1:A:622:LYS:HE3	2.21	0.41
1:B:218:PRO:HD3	1:B:305:TRP:HB3	2.03	0.41
1:C:626:ILE:O	1:C:650:GLY:HA2	2.21	0.41
1:C:757:HIS:HD2	1:D:729:ASP:OD2	2.04	0.41
1:D:237:GLU:HG2	1:D:253:ARG:HG2	2.02	0.41
1:D:554:LYS:HE2	6:D:910:HOH:O	2.21	0.41
1:B:237:GLU:HG2	1:B:253:ARG:HG2	2.02	0.41
1:C:49:LEU:HB3	1:C:749:GLN:HG2	2.03	0.41
1:D:640:LEU:HB3	1:D:698:VAL:HG21	2.03	0.40
1:A:232:GLU:HB3	1:A:262:GLU:HG2	2.04	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:D:811:6Z8:C13	4:D:811:6Z8:C4	2.99	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	725/766 (95%)	699 (96%)	26 (4%)	0	100	100
1	B	725/766 (95%)	699 (96%)	26 (4%)	0	100	100
1	C	725/766 (95%)	699 (96%)	26 (4%)	0	100	100
1	D	725/766 (95%)	696 (96%)	29 (4%)	0	100	100
All	All	2900/3064 (95%)	2793 (96%)	107 (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	651/682 (96%)	628 (96%)	23 (4%)	36	53
1	B	651/682 (96%)	630 (97%)	21 (3%)	39	57
1	C	651/682 (96%)	629 (97%)	22 (3%)	37	54
1	D	651/682 (96%)	631 (97%)	20 (3%)	40	58

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
All	All	2604/2728 (96%)	2518 (97%)	86 (3%)	38 56

All (86) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	100	TYR
1	A	140	ARG
1	A	145	GLU
1	A	202	VAL
1	A	219	ASN
1	A	230	ASP
1	A	358	ARG
1	A	367	ASP
1	A	385	CYS
1	A	393	ASN
1	A	440	THR
1	A	514	LEU
1	A	516	VAL
1	A	518	ASN
1	A	520	HIS
1	A	547	TYR
1	A	566	TYR
1	A	627	TRP
1	A	629	TRP
1	A	679	ASN
1	A	699	GLU
1	A	704	HIS
1	A	764	SER
1	B	90	LEU
1	B	100	TYR
1	B	145	GLU
1	B	230	ASP
1	B	358	ARG
1	B	367	ASP
1	B	385	CYS
1	B	393	ASN
1	B	440	THR
1	B	472	CYS
1	B	514	LEU
1	B	516	VAL
1	B	518	ASN
1	B	520	HIS

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Mol	Chain	Res	Type
1	B	547	TYR
1	B	566	TYR
1	B	589	LYS
1	B	627	TRP
1	B	629	TRP
1	B	630	SER
1	B	704	HIS
1	C	90	LEU
1	C	100	TYR
1	C	142	LEU
1	C	145	GLU
1	C	230	ASP
1	C	294	LEU
1	C	358	ARG
1	C	367	ASP
1	C	379	GLU
1	C	385	CYS
1	C	399	LYS
1	C	440	THR
1	C	514	LEU
1	C	516	VAL
1	C	518	ASN
1	C	520	HIS
1	C	523	LYS
1	C	547	TYR
1	C	566	TYR
1	C	627	TRP
1	C	629	TRP
1	C	630	SER
1	D	41	ARG
1	D	100	TYR
1	D	145	GLU
1	D	219	ASN
1	D	230	ASP
1	D	250	LYS
1	D	358	ARG
1	D	385	CYS
1	D	440	THR
1	D	472	CYS
1	D	492	ARG
1	D	514	LEU
1	D	518	ASN

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Mol	Chain	Res	Type
1	D	523	LYS
1	D	547	TYR
1	D	566	TYR
1	D	627	TRP
1	D	629	TRP
1	D	630	SER
1	D	764	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (40) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	75	ASN
1	A	169	ASN
1	A	170	ASN
1	A	192	ASN
1	A	247	GLN
1	A	369	ASN
1	A	393	ASN
1	A	463	ASN
1	A	483	HIS
1	A	679	ASN
1	A	757	HIS
1	B	75	ASN
1	B	169	ASN
1	B	170	ASN
1	B	192	ASN
1	B	247	GLN
1	B	345	HIS
1	B	393	ASN
1	B	463	ASN
1	B	483	HIS
1	C	75	ASN
1	C	169	ASN
1	C	170	ASN
1	C	192	ASN
1	C	247	GLN
1	C	422	HIS
1	C	463	ASN
1	C	483	HIS
1	C	704	HIS
1	C	757	HIS
1	D	75	ASN

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Mol	Chain	Res	Type
1	D	169	ASN
1	D	170	ASN
1	D	192	ASN
1	D	247	GLN
1	D	435	GLN
1	D	463	ASN
1	D	483	HIS
1	D	586	GLN
1	D	704	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 ⓘ

18 monosaccharides are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z  > 2$	Counts	RMSZ	$\# Z  > 2$
2	NAG	E	1	1,2	14,14,15	0.29	0	17,19,21	0.70	0
2	NAG	E	2	2	14,14,15	0.30	0	17,19,21	0.64	1 (5%)
2	NAG	F	1	1,2	14,14,15	0.29	0	17,19,21	0.85	1 (5%)
2	NAG	F	2	2	14,14,15	0.28	0	17,19,21	0.74	0
2	NAG	G	1	1,2	14,14,15	0.33	0	17,19,21	0.87	1 (5%)
2	NAG	G	2	2	14,14,15	0.32	0	17,19,21	0.77	1 (5%)
2	NAG	H	1	1,2	14,14,15	0.30	0	17,19,21	0.76	0
2	NAG	H	2	2	14,14,15	0.29	0	17,19,21	0.61	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
2	NAG	I	1	1,2	14,14,15	0.26	0	17,19,21	0.89	1 (5%)
2	NAG	I	2	2	14,14,15	0.30	0	17,19,21	0.53	0
2	NAG	J	1	1,2	14,14,15	0.28	0	17,19,21	0.75	1 (5%)
2	NAG	J	2	2	14,14,15	0.36	0	17,19,21	1.00	2 (11%)
2	NAG	K	1	1,2	14,14,15	0.27	0	17,19,21	0.60	0
2	NAG	K	2	2	14,14,15	0.30	0	17,19,21	0.59	0
2	NAG	L	1	1,2	14,14,15	0.27	0	17,19,21	1.28	1 (5%)
2	NAG	L	2	2	14,14,15	0.31	0	17,19,21	0.56	0
2	NAG	M	1	1,2	14,14,15	0.30	0	17,19,21	0.91	1 (5%)
2	NAG	M	2	2	14,14,15	0.28	0	17,19,21	0.56	0

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	NAG	E	1	1,2	-	0/6/23/26	0/1/1/1
2	NAG	E	2	2	-	2/6/23/26	0/1/1/1
2	NAG	F	1	1,2	-	1/6/23/26	0/1/1/1
2	NAG	F	2	2	-	2/6/23/26	0/1/1/1
2	NAG	G	1	1,2	-	0/6/23/26	0/1/1/1
2	NAG	G	2	2	-	0/6/23/26	0/1/1/1
2	NAG	H	1	1,2	-	0/6/23/26	0/1/1/1
2	NAG	H	2	2	-	0/6/23/26	0/1/1/1
2	NAG	I	1	1,2	-	0/6/23/26	0/1/1/1
2	NAG	I	2	2	-	0/6/23/26	0/1/1/1
2	NAG	J	1	1,2	-	0/6/23/26	0/1/1/1
2	NAG	J	2	2	-	2/6/23/26	0/1/1/1
2	NAG	K	1	1,2	-	0/6/23/26	0/1/1/1
2	NAG	K	2	2	-	2/6/23/26	0/1/1/1
2	NAG	L	1	1,2	-	2/6/23/26	0/1/1/1
2	NAG	L	2	2	-	0/6/23/26	0/1/1/1
2	NAG	M	1	1,2	-	0/6/23/26	0/1/1/1
2	NAG	M	2	2	-	1/6/23/26	0/1/1/1

There are no bond length outliers.

All (10) bond angle outliers are listed below:



Mol	Chain	Res	Type	Atoms	Z	Observed( $^{\circ}$ )	Ideal( $^{\circ}$ )
2	L	1	NAG	C1-O5-C5	4.83	118.73	112.19
2	G	1	NAG	C1-O5-C5	2.97	116.21	112.19
2	I	1	NAG	C1-O5-C5	2.85	116.05	112.19
2	F	1	NAG	C1-O5-C5	2.84	116.04	112.19
2	M	1	NAG	C1-O5-C5	2.72	115.88	112.19
2	J	1	NAG	C1-O5-C5	2.70	115.85	112.19
2	G	2	NAG	C1-O5-C5	2.53	115.62	112.19
2	J	2	NAG	C2-N2-C7	2.44	126.38	122.90
2	E	2	NAG	C1-O5-C5	2.14	115.10	112.19
2	J	2	NAG	C1-O5-C5	2.06	114.99	112.19

There are no chirality outliers.

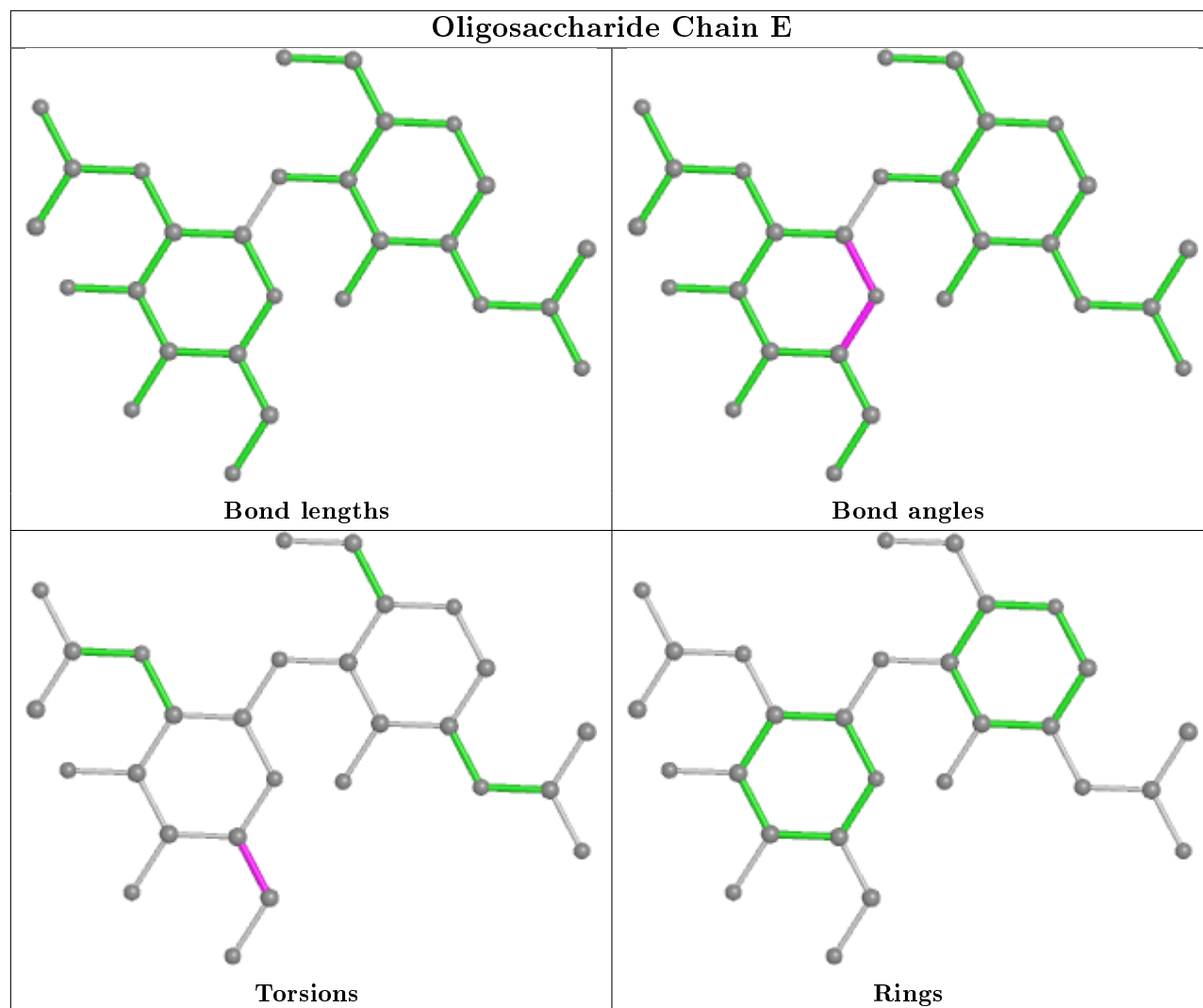
All (12) torsion outliers are listed below:

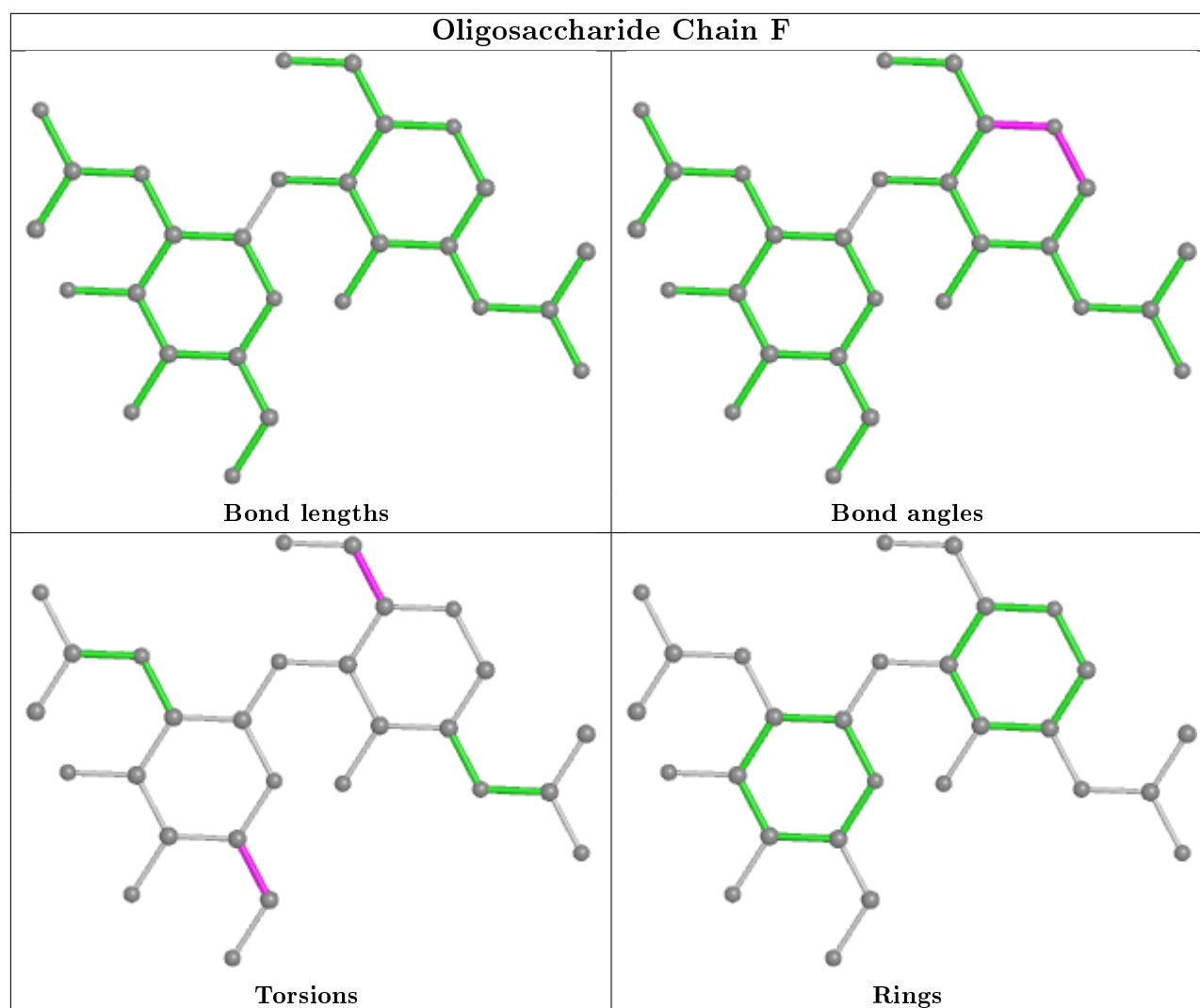
Mol	Chain	Res	Type	Atoms
2	L	1	NAG	O5-C5-C6-O6
2	F	2	NAG	O5-C5-C6-O6
2	F	1	NAG	O5-C5-C6-O6
2	K	2	NAG	C4-C5-C6-O6
2	K	2	NAG	O5-C5-C6-O6
2	L	1	NAG	C4-C5-C6-O6
2	E	2	NAG	C4-C5-C6-O6
2	E	2	NAG	O5-C5-C6-O6
2	M	2	NAG	C4-C5-C6-O6
2	F	2	NAG	C4-C5-C6-O6
2	J	2	NAG	C1-C2-N2-C7
2	J	2	NAG	C3-C2-N2-C7

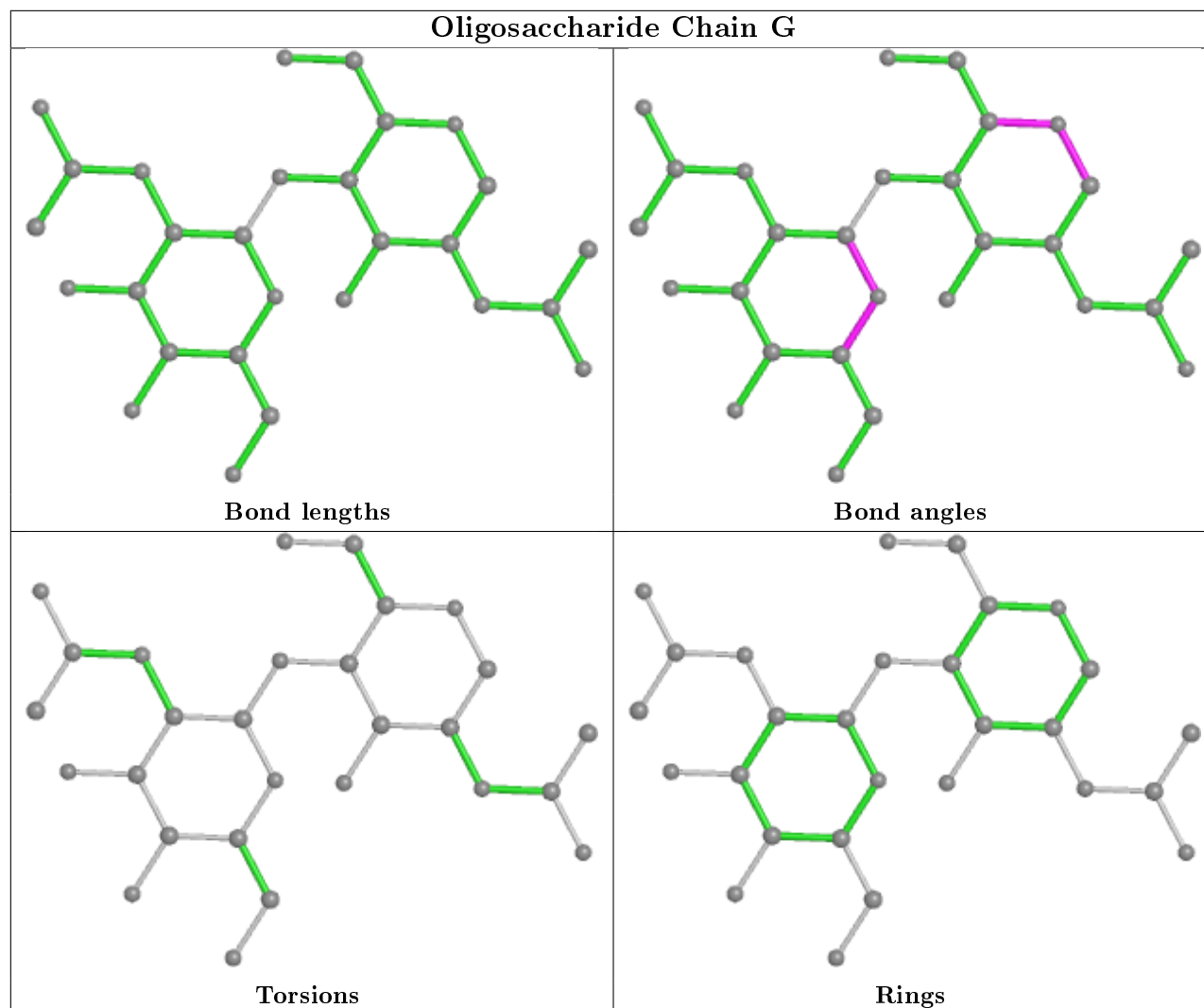
There are no ring outliers.

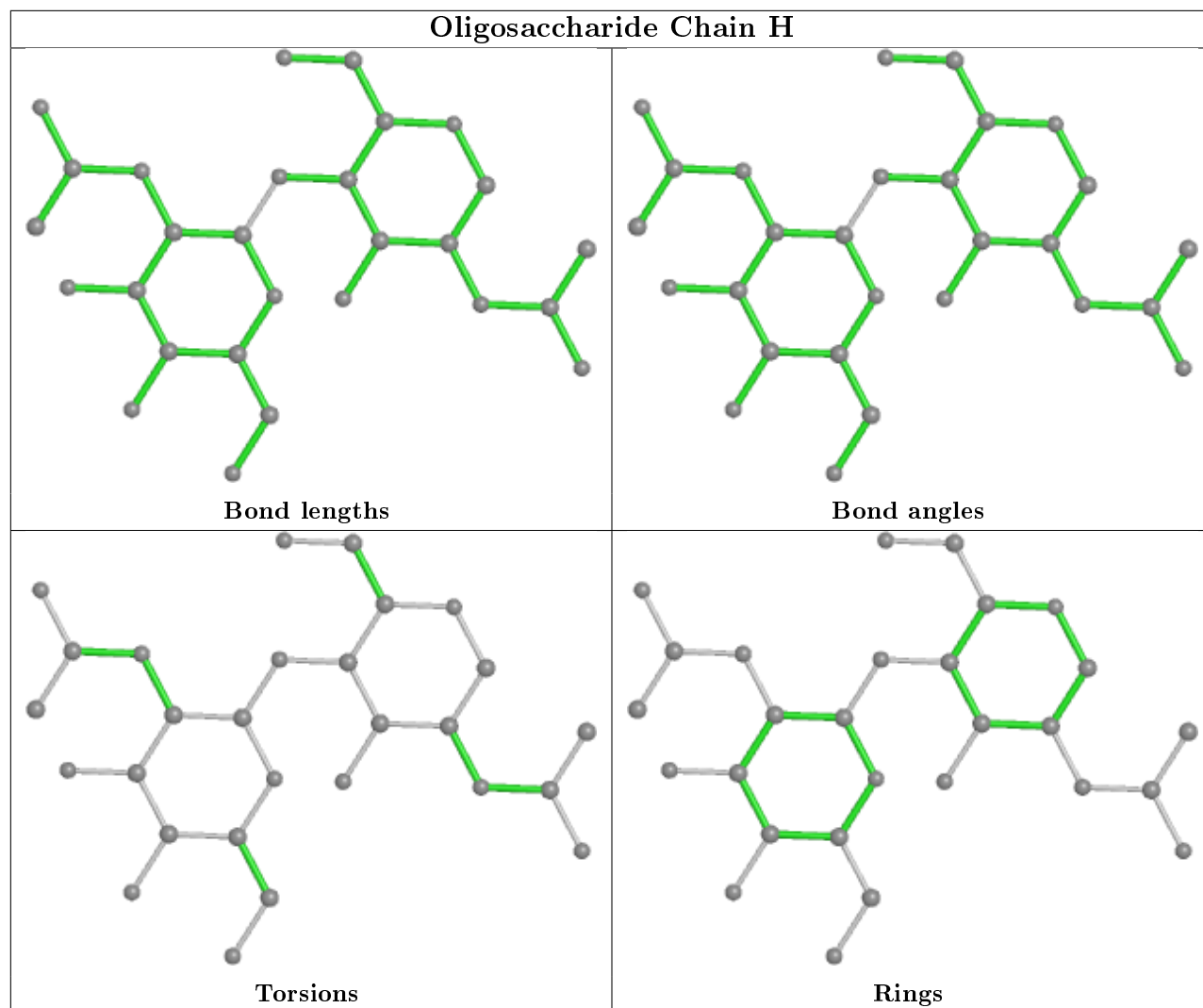
No monomer is involved in short contacts.

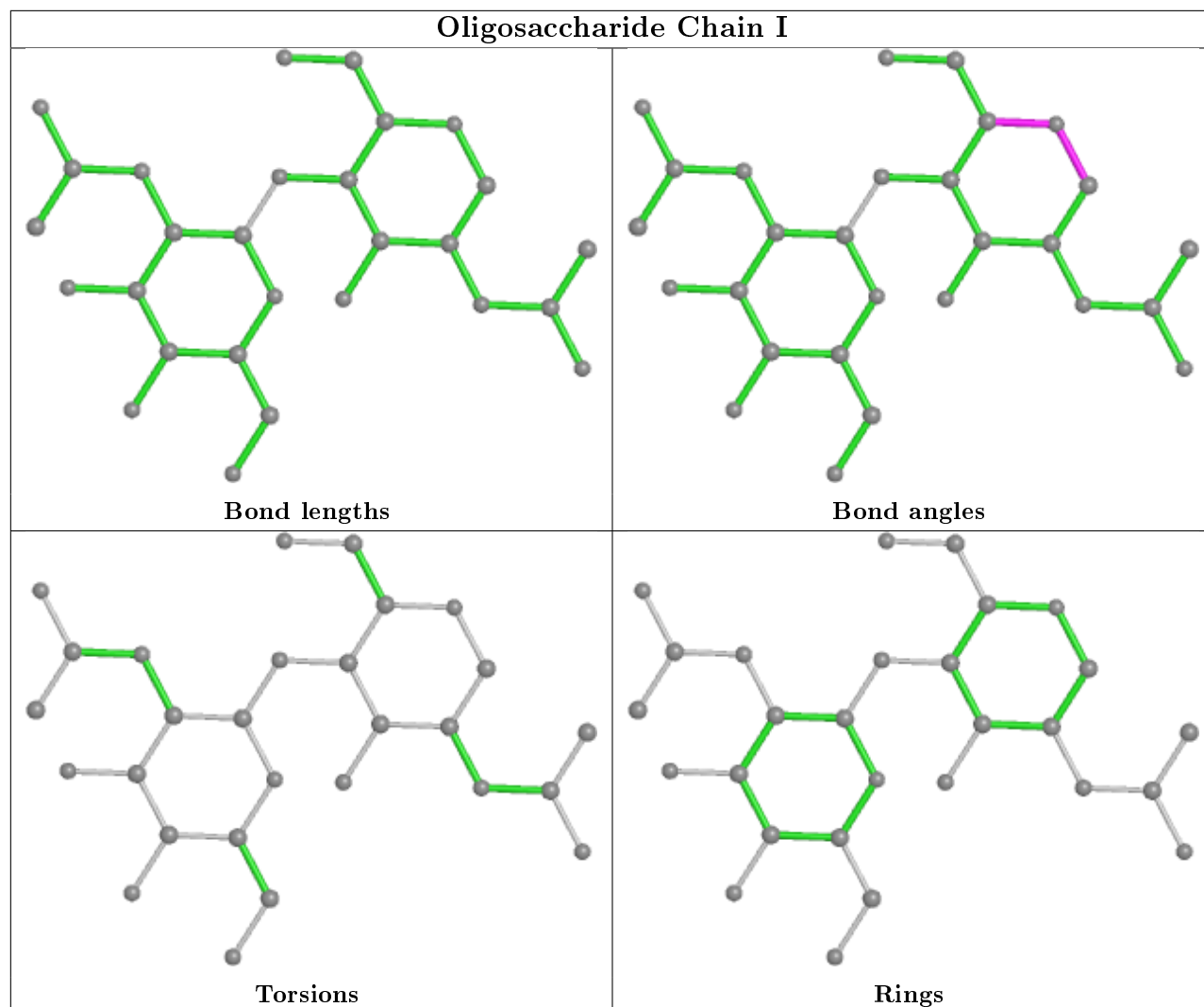
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.

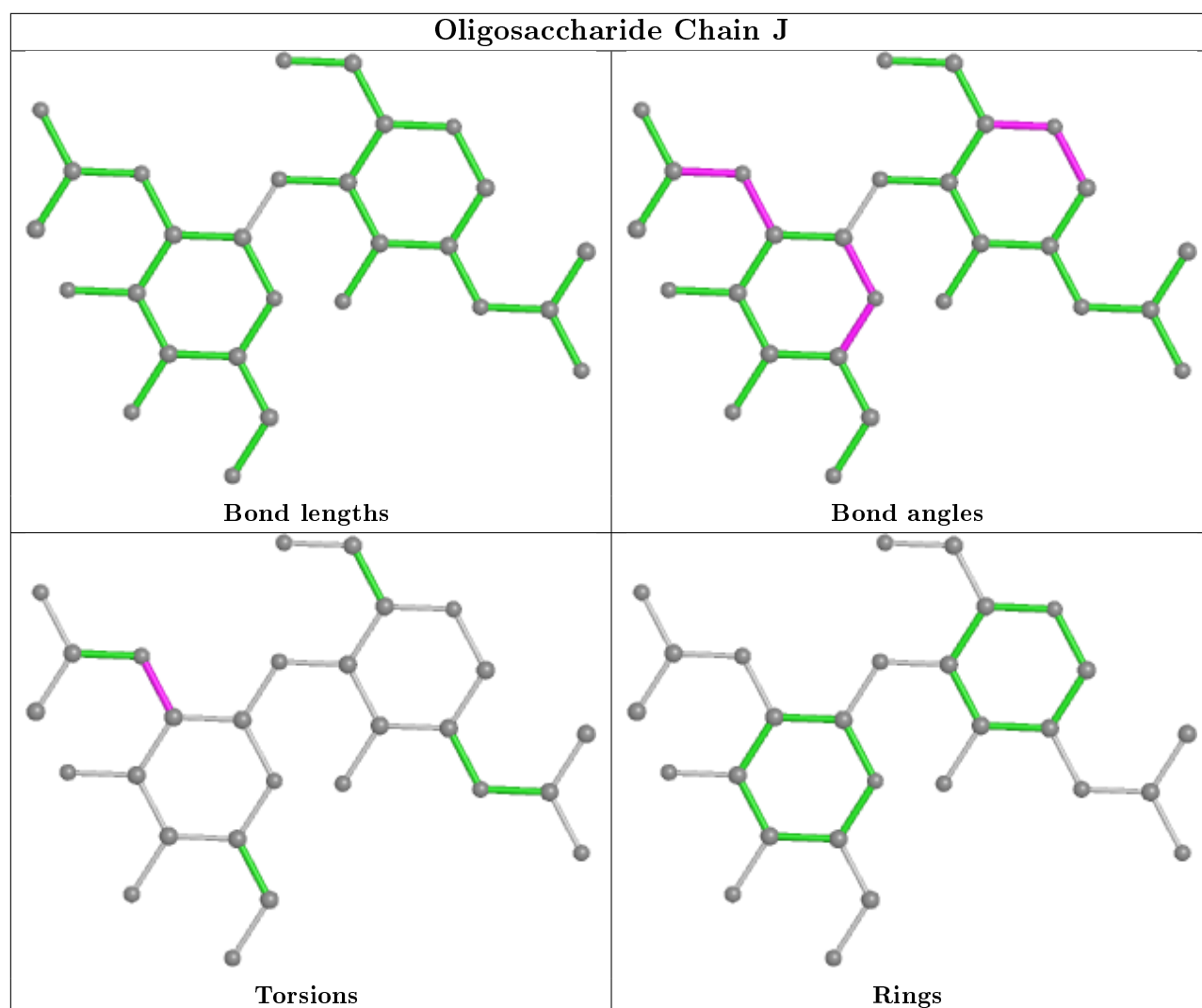


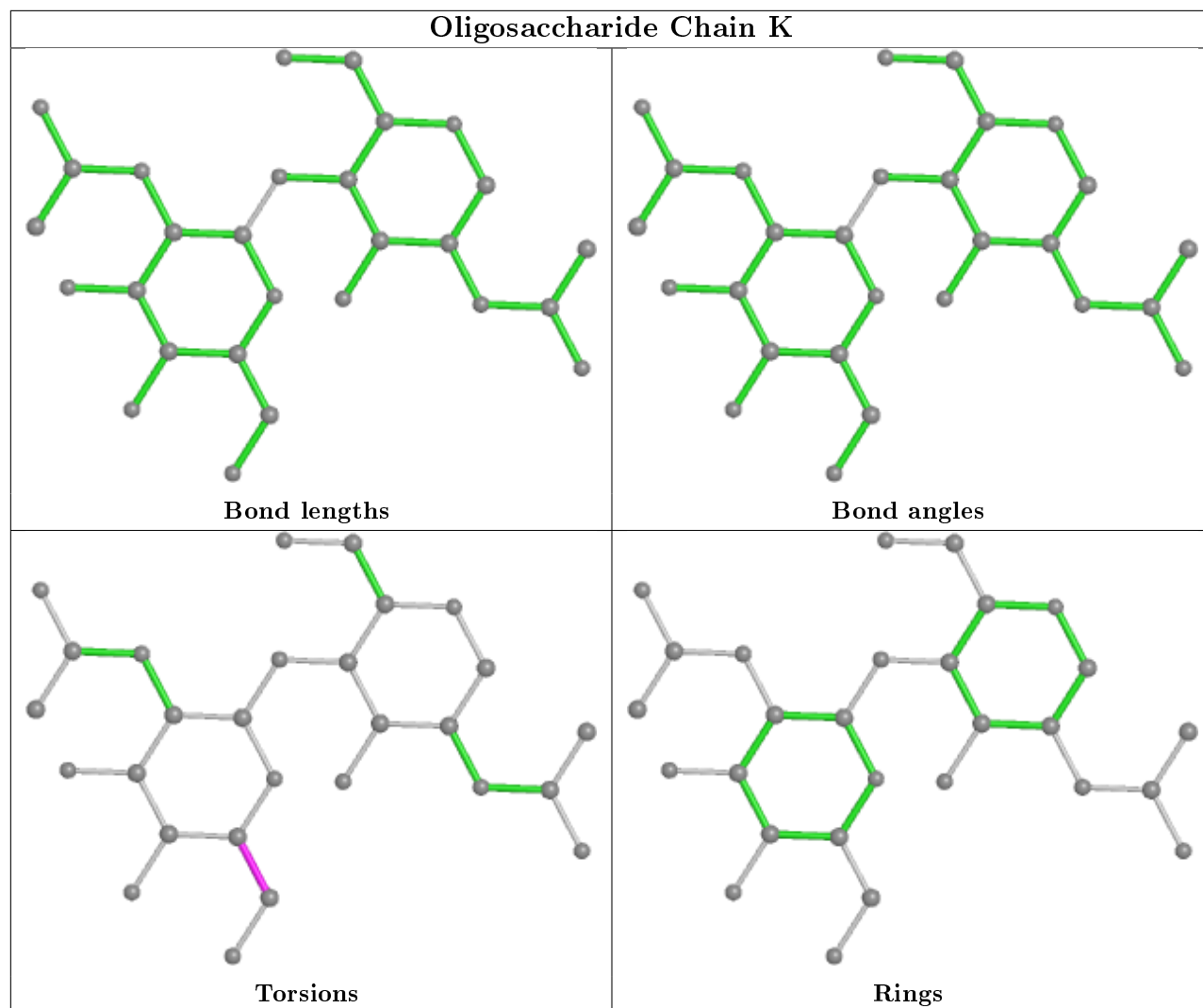




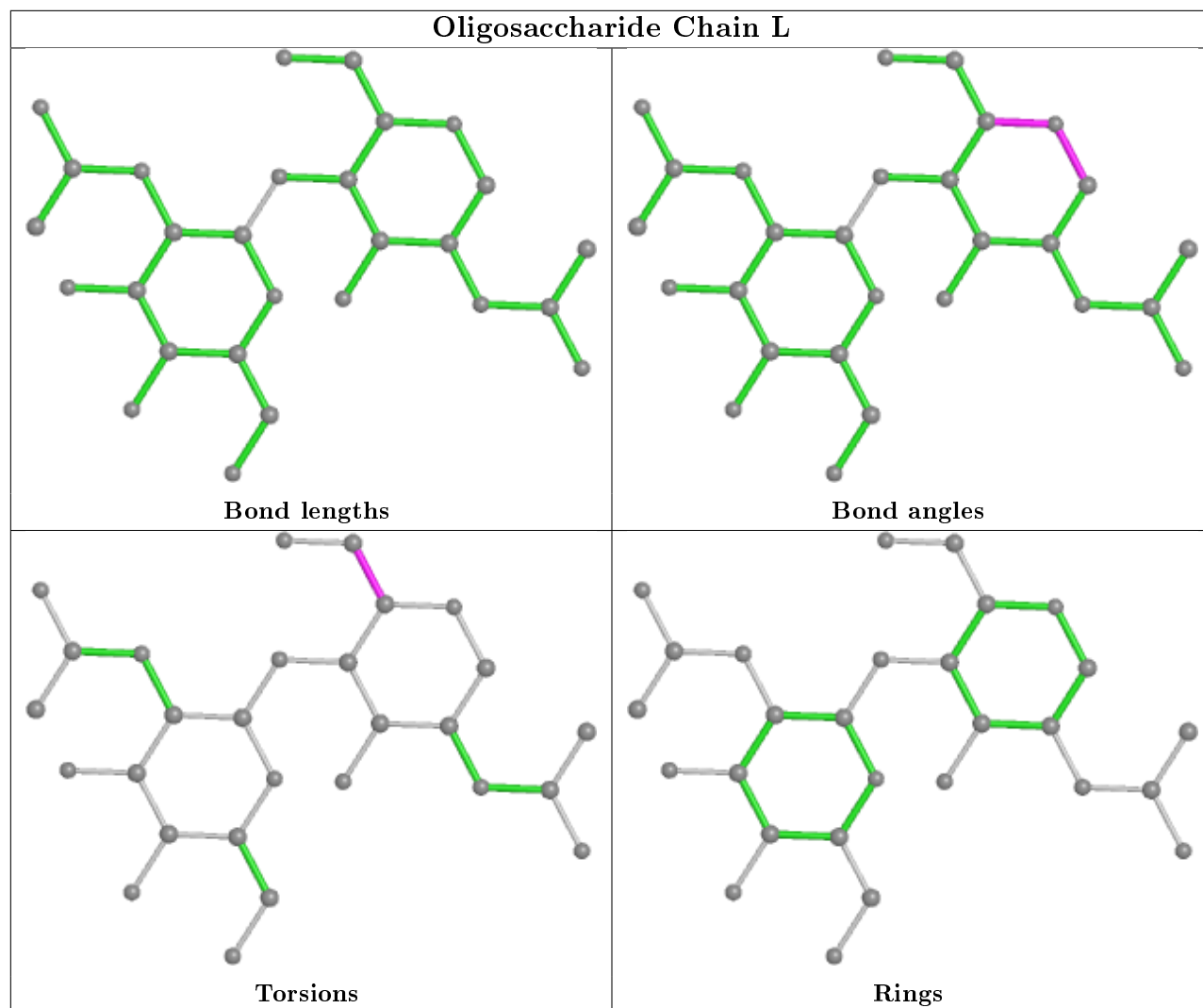


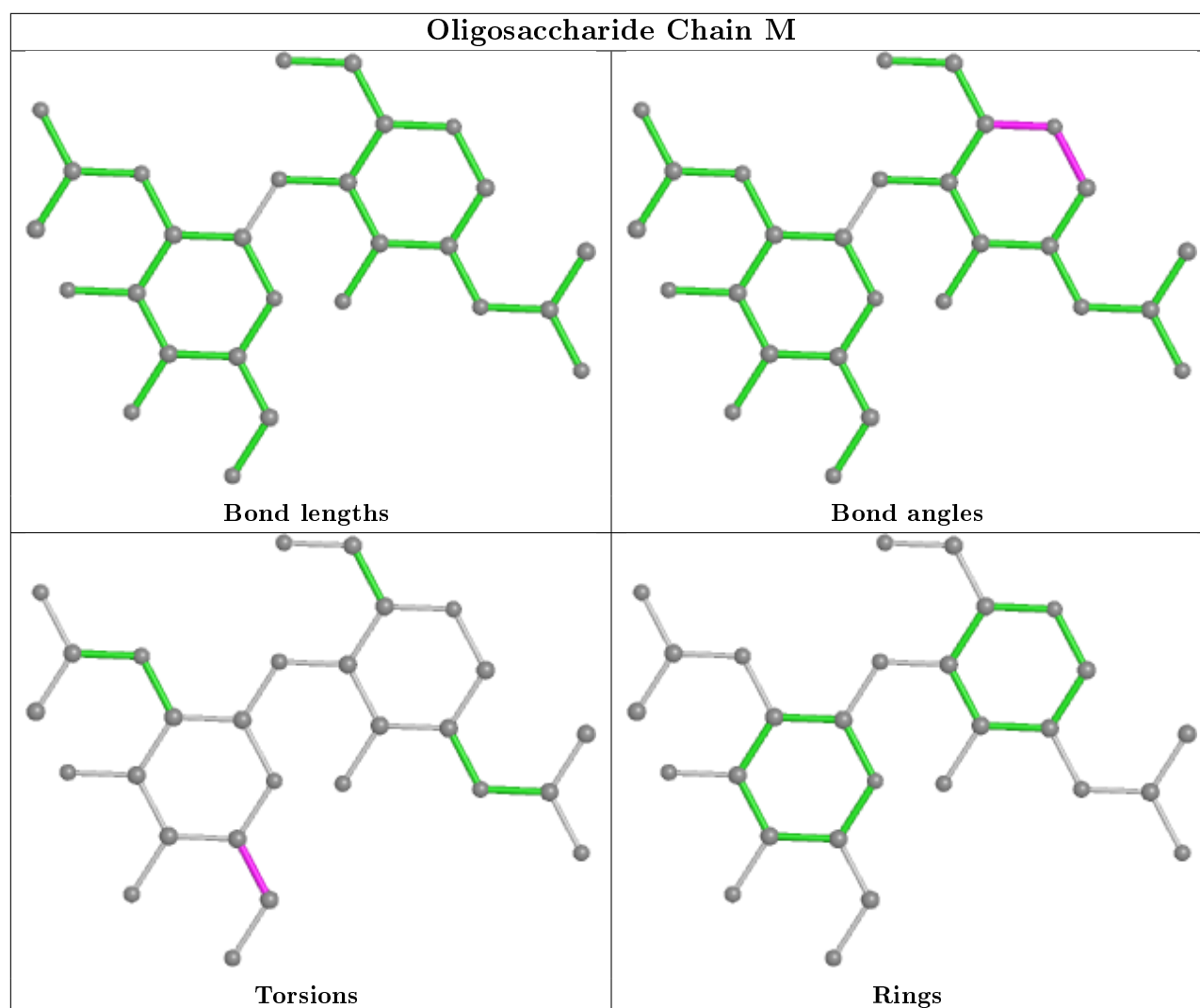












## 5.6 Ligand geometry [i](#)

27 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	$\# Z  > 2$
5	SO4	D	812	-	4,4,4	0.13	0	6,6,6	0.11	0
4	6Z8	D	813	-	24,31,31	1.12	3 (12%)	27,46,46	1.59	3 (11%)
5	SO4	A	810	-	4,4,4	0.24	0	6,6,6	0.28	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
3	NAG	C	802	1	14,14,15	0.28	0	17,19,21	0.63	0
3	NAG	D	806	1	14,14,15	0.34	0	17,19,21	0.78	1 (5%)
3	NAG	B	805	1	14,14,15	0.29	0	17,19,21	0.63	0
3	NAG	C	803	1	14,14,15	0.34	0	17,19,21	0.68	0
4	6Z8	A	809	-	24,31,31	1.23	3 (12%)	27,46,46	1.57	4 (14%)
3	NAG	A	802	1	14,14,15	0.28	0	17,19,21	0.62	0
3	NAG	A	805	1	14,14,15	0.32	0	17,19,21	0.80	1 (5%)
3	NAG	D	801	1	14,14,15	0.39	0	17,19,21	1.81	2 (11%)
3	NAG	B	801	1	14,14,15	0.34	0	17,19,21	0.65	1 (5%)
3	NAG	A	801	1	14,14,15	0.32	0	17,19,21	0.79	1 (5%)
3	NAG	C	805	1	14,14,15	0.46	0	17,19,21	1.77	4 (23%)
4	6Z8	A	811	-	24,31,31	1.12	3 (12%)	27,46,46	1.64	4 (14%)
4	6Z8	D	811	-	24,31,31	1.21	3 (12%)	27,46,46	1.73	5 (18%)
4	6Z8	B	811	-	24,31,31	1.15	3 (12%)	27,46,46	1.62	4 (14%)
5	SO4	C	809	-	4,4,4	0.31	0	6,6,6	0.31	0
4	6Z8	C	808	-	24,31,31	1.11	3 (12%)	27,46,46	1.59	3 (11%)
3	NAG	B	806	1	14,14,15	0.33	0	17,19,21	0.88	2 (11%)
5	SO4	B	810	-	4,4,4	0.12	0	6,6,6	0.13	0
3	NAG	C	801	1	14,14,15	0.35	0	17,19,21	0.82	1 (5%)
3	NAG	A	806	1	14,14,15	0.26	0	17,19,21	0.62	1 (5%)
4	6Z8	C	810	-	24,31,31	1.10	3 (12%)	27,46,46	1.54	3 (11%)
3	NAG	C	804	1	14,14,15	0.30	0	17,19,21	0.66	0
4	6Z8	B	809	-	24,31,31	1.19	3 (12%)	27,46,46	1.64	4 (14%)
3	NAG	B	804	1	14,14,15	0.34	0	17,19,21	0.92	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
4	6Z8	D	813	-	-	1/4/18/18	0/4/4/4
3	NAG	C	802	1	-	0/6/23/26	0/1/1/1
3	NAG	D	806	1	-	0/6/23/26	0/1/1/1
3	NAG	B	805	1	-	0/6/23/26	0/1/1/1
3	NAG	C	803	1	-	0/6/23/26	0/1/1/1
4	6Z8	A	809	-	-	0/4/18/18	0/4/4/4
3	NAG	A	802	1	-	0/6/23/26	0/1/1/1
3	NAG	A	805	1	-	0/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	NAG	D	801	1	-	0/6/23/26	0/1/1/1
3	NAG	B	801	1	-	1/6/23/26	0/1/1/1
3	NAG	A	801	1	-	1/6/23/26	0/1/1/1
3	NAG	C	805	1	-	2/6/23/26	0/1/1/1
4	6Z8	A	811	-	-	0/4/18/18	0/4/4/4
4	6Z8	D	811	-	-	0/4/18/18	0/4/4/4
4	6Z8	B	811	-	-	0/4/18/18	0/4/4/4
4	6Z8	C	808	-	-	0/4/18/18	0/4/4/4
3	NAG	B	806	1	-	0/6/23/26	0/1/1/1
3	NAG	C	801	1	-	0/6/23/26	0/1/1/1
3	NAG	A	806	1	-	1/6/23/26	0/1/1/1
4	6Z8	C	810	-	-	0/4/18/18	0/4/4/4
3	NAG	C	804	1	-	0/6/23/26	0/1/1/1
4	6Z8	B	809	-	-	0/4/18/18	0/4/4/4
3	NAG	B	804	1	-	0/6/23/26	0/1/1/1

All (24) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	811	6Z8	C5-N7	3.47	1.36	1.33
4	A	809	6Z8	C5-N7	3.46	1.36	1.33
4	D	811	6Z8	C5-N7	3.45	1.36	1.33
4	A	809	6Z8	C2-C5	-3.43	1.34	1.39
4	C	810	6Z8	C5-N7	3.39	1.36	1.33
4	A	811	6Z8	C5-N7	3.35	1.36	1.33
4	D	813	6Z8	C5-N7	3.30	1.36	1.33
4	B	809	6Z8	C5-N7	3.27	1.36	1.33
4	C	808	6Z8	C5-N7	3.10	1.36	1.33
4	C	808	6Z8	C2-C5	-2.96	1.35	1.39
4	D	811	6Z8	C2-C5	-2.95	1.35	1.39
4	B	809	6Z8	C2-C5	-2.79	1.35	1.39
4	A	811	6Z8	C6-N11	2.70	1.42	1.38
4	B	809	6Z8	C6-N11	2.62	1.42	1.38
4	D	813	6Z8	C6-N11	2.57	1.42	1.38
4	D	811	6Z8	C6-N11	2.56	1.42	1.38
4	B	811	6Z8	C6-N11	2.47	1.41	1.38
4	A	809	6Z8	C6-N11	2.45	1.41	1.38
4	C	810	6Z8	C2-C5	-2.42	1.36	1.39
4	C	810	6Z8	C6-N11	2.35	1.41	1.38
4	D	813	6Z8	C2-C5	-2.16	1.36	1.39
4	B	811	6Z8	C2-C5	-2.15	1.36	1.39
4	C	808	6Z8	C6-N11	2.08	1.41	1.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	811	6Z8	C2-C5	-2.01	1.36	1.39

All (45) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	801	NAG	C1-O5-C5	5.76	119.99	112.19
4	B	811	6Z8	C2-C6-N11	-5.36	112.49	118.20
4	A	811	6Z8	C2-C6-N11	-5.30	112.55	118.20
4	D	813	6Z8	C2-C6-N11	-5.29	112.56	118.20
4	B	809	6Z8	C2-C6-N11	-5.19	112.67	118.20
4	D	811	6Z8	C2-C6-N11	-5.16	112.70	118.20
4	C	810	6Z8	C2-C6-N11	-5.05	112.82	118.20
4	C	808	6Z8	C2-C6-N11	-5.03	112.84	118.20
4	C	808	6Z8	C6-C2-C5	5.02	123.19	119.96
4	A	809	6Z8	C2-C6-N11	-4.99	112.88	118.20
4	B	809	6Z8	C6-C2-C5	4.86	123.08	119.96
4	D	813	6Z8	C6-C2-C5	4.83	123.06	119.96
4	D	811	6Z8	C6-C2-C5	4.81	123.05	119.96
4	A	809	6Z8	C6-C2-C5	4.79	123.03	119.96
4	A	811	6Z8	C6-C2-C5	4.69	122.97	119.96
4	C	810	6Z8	C6-C2-C5	4.63	122.93	119.96
4	B	811	6Z8	C6-C2-C5	4.61	122.92	119.96
3	C	805	NAG	C1-C2-N2	4.33	117.88	110.49
3	C	805	NAG	C2-N2-C7	3.74	128.22	122.90
3	C	805	NAG	O5-C1-C2	-3.58	105.63	111.29
3	D	801	NAG	O5-C1-C2	3.58	116.94	111.29
4	D	811	6Z8	C20-C13-N8	-3.57	105.14	109.64
3	B	804	NAG	C1-O5-C5	3.02	116.28	112.19
3	C	801	NAG	C1-O5-C5	2.90	116.12	112.19
4	B	809	6Z8	C20-C13-N8	-2.78	106.14	109.64
4	D	811	6Z8	C26-C20-C13	-2.66	107.24	110.65
3	D	806	NAG	C1-O5-C5	2.33	115.34	112.19
3	A	805	NAG	O5-C1-C2	-2.32	107.62	111.29
3	A	801	NAG	C1-O5-C5	2.32	115.33	112.19
3	B	806	NAG	C1-O5-C5	2.30	115.30	112.19
4	A	809	6Z8	C18-N10-C5	2.29	121.50	118.25
4	C	810	6Z8	C18-N10-C5	2.28	121.48	118.25
3	B	801	NAG	C1-O5-C5	2.20	115.17	112.19
4	A	811	6Z8	C13-N8-C3	-2.17	117.38	122.03
4	C	808	6Z8	C18-N10-C5	2.17	121.33	118.25
3	C	805	NAG	C1-O5-C5	2.15	115.11	112.19
4	B	809	6Z8	C18-N10-C5	2.15	121.30	118.25

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	D	813	6Z8	C18-N10-C5	2.15	121.30	118.25
4	B	811	6Z8	C18-N10-C5	2.15	121.29	118.25
4	D	811	6Z8	C18-N10-C5	2.14	121.28	118.25
4	A	811	6Z8	C18-N10-C5	2.09	121.22	118.25
4	B	811	6Z8	C14-N8-C3	-2.05	117.64	122.03
3	A	806	NAG	C1-O5-C5	2.04	114.96	112.19
4	A	809	6Z8	C20-C13-N8	-2.03	107.09	109.64
3	B	806	NAG	O5-C1-C2	-2.01	108.11	111.29

There are no chirality outliers.

All (6) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	C	805	NAG	C1-C2-N2-C7
3	A	801	NAG	O5-C5-C6-O6
3	C	805	NAG	O5-C5-C6-O6
3	B	801	NAG	O5-C5-C6-O6
3	A	806	NAG	C4-C5-C6-O6
4	D	813	6Z8	N1-C4-C9-C16

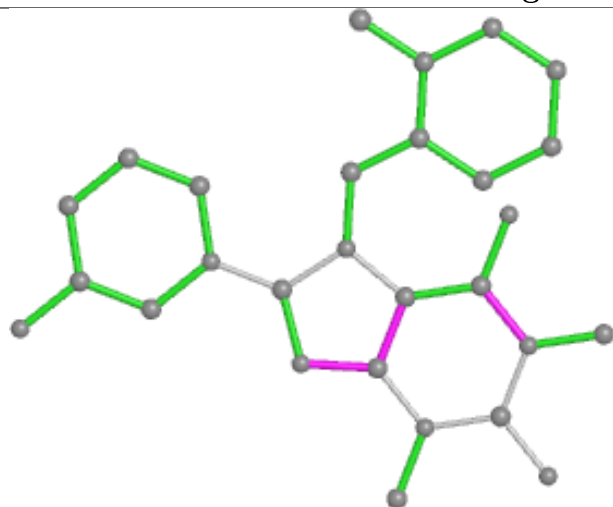
There are no ring outliers.

5 monomers are involved in 7 short contacts:

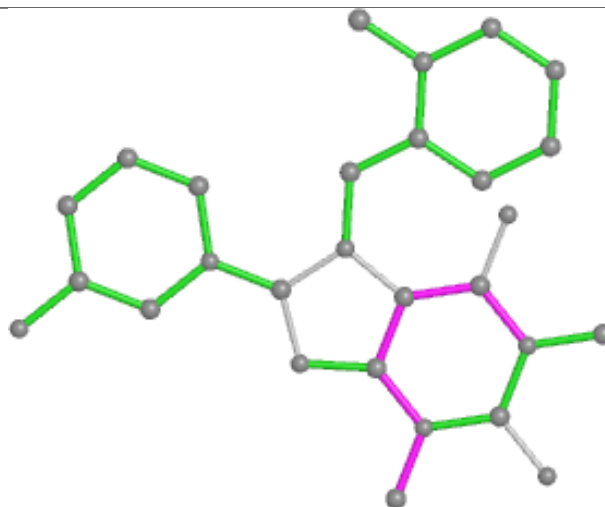
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	D	806	NAG	1	0
4	A	809	6Z8	2	0
3	A	801	NAG	1	0
4	D	811	6Z8	1	0
4	C	808	6Z8	2	0

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

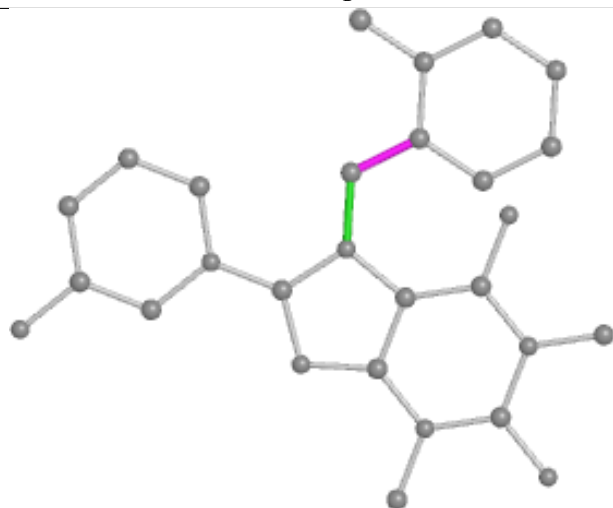
## Ligand 6Z8 D 813



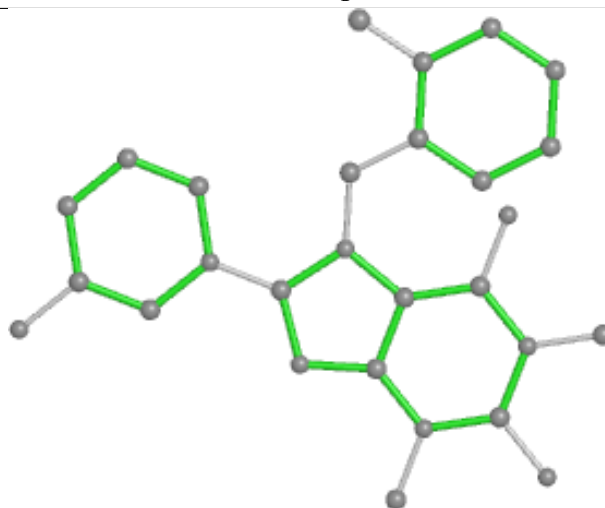
Bond lengths



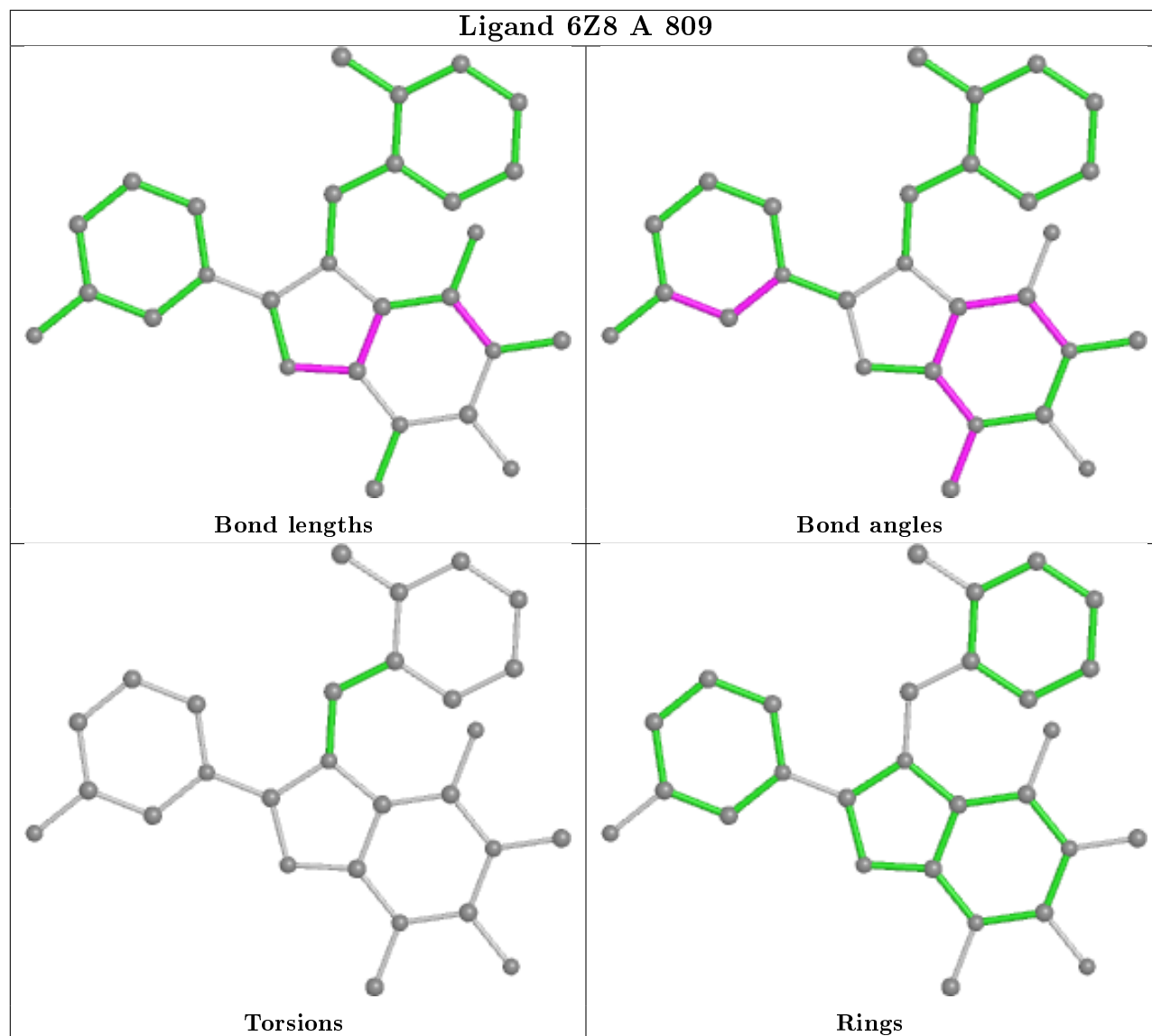
Bond angles



Torsions

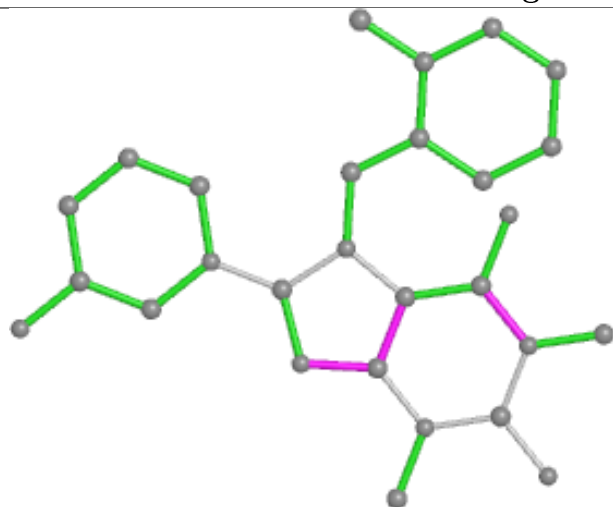


Rings

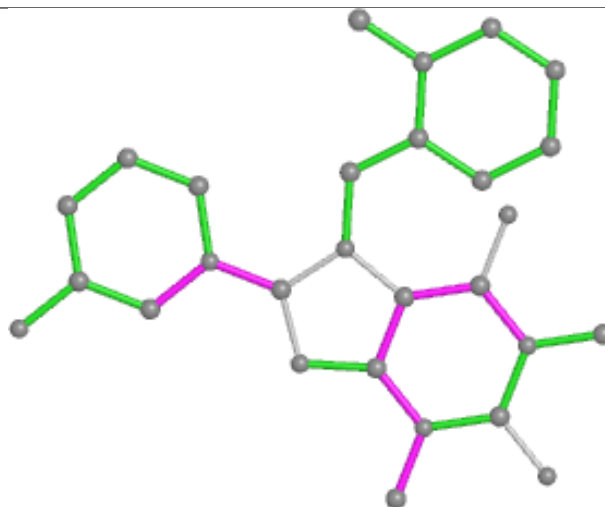




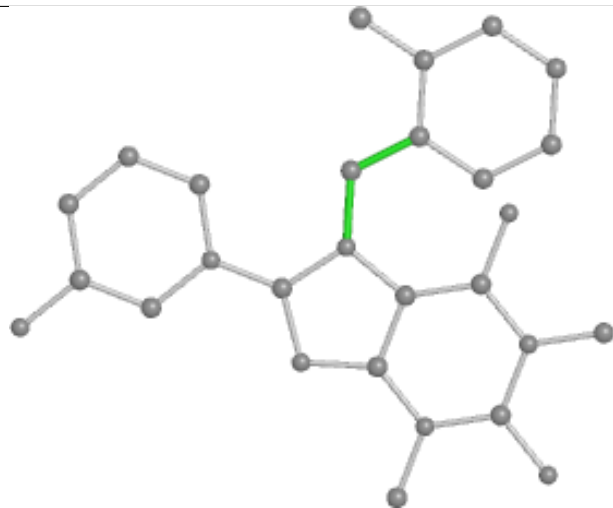
## Ligand 6Z8 A 811



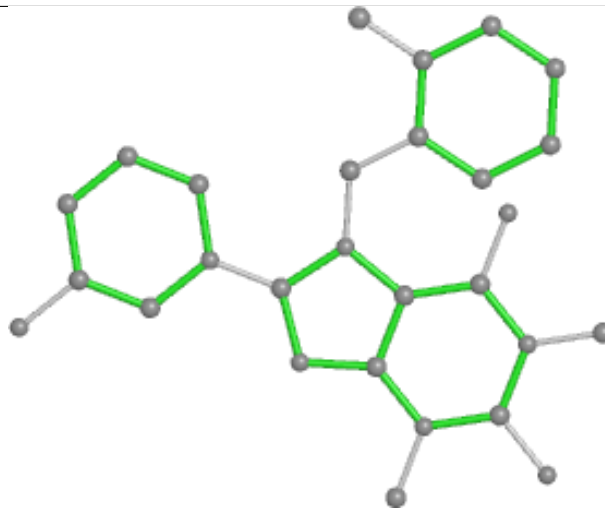
Bond lengths



Bond angles

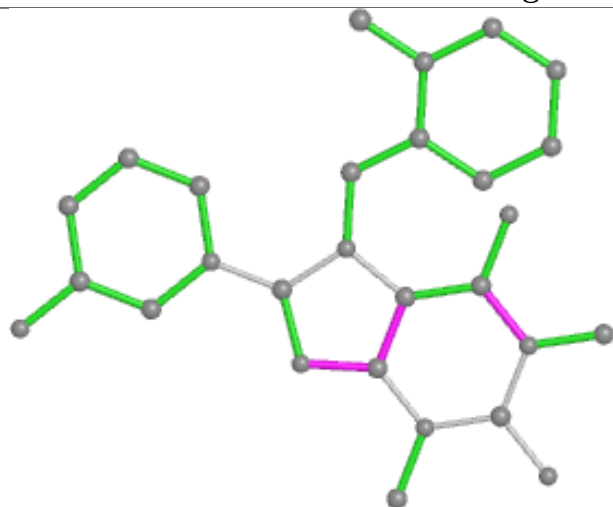


Torsions

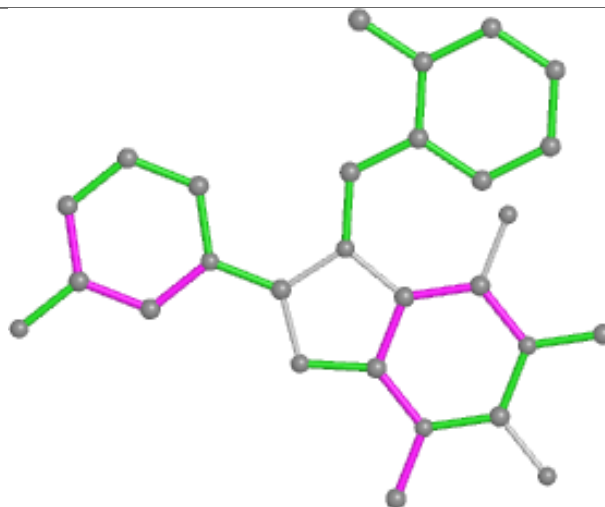


Rings

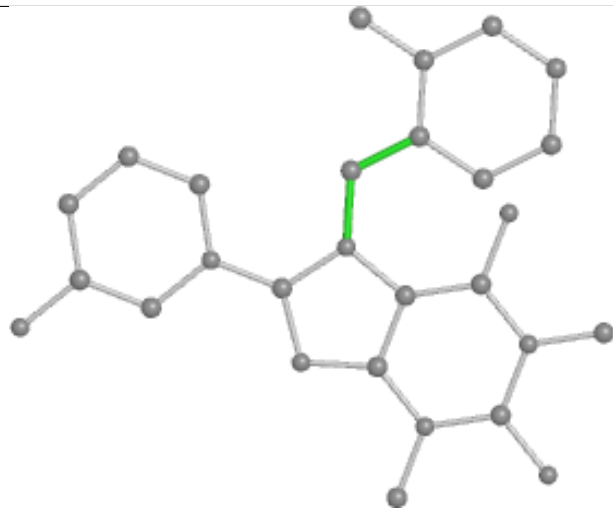
## Ligand 6Z8 D 811



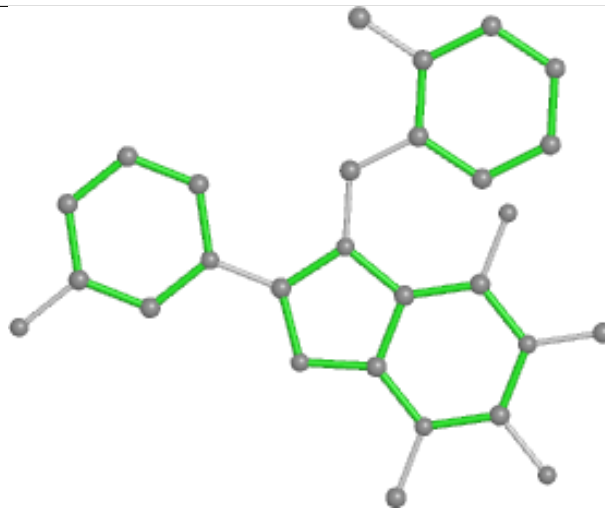
Bond lengths



Bond angles

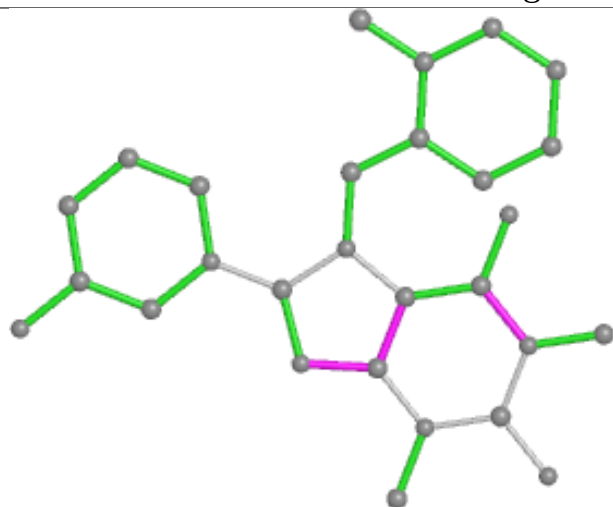


Torsions

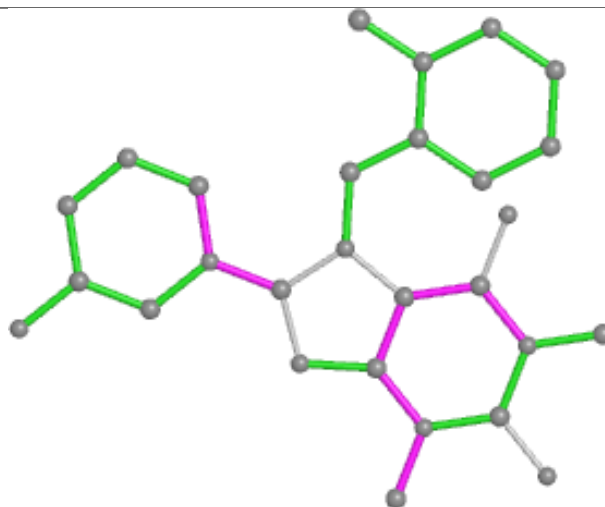


Rings

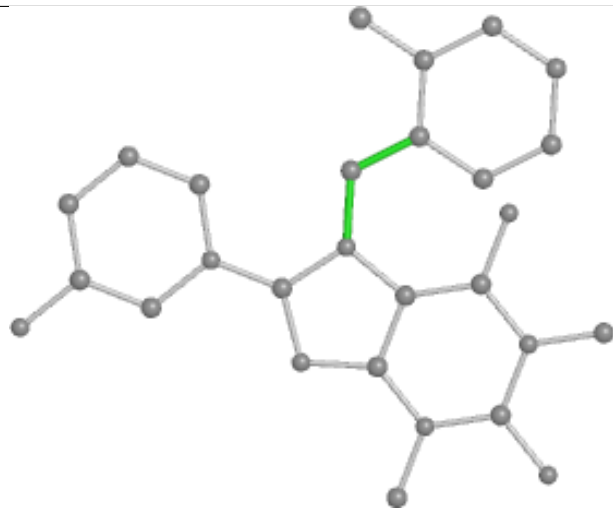
## Ligand 6Z8 B 811



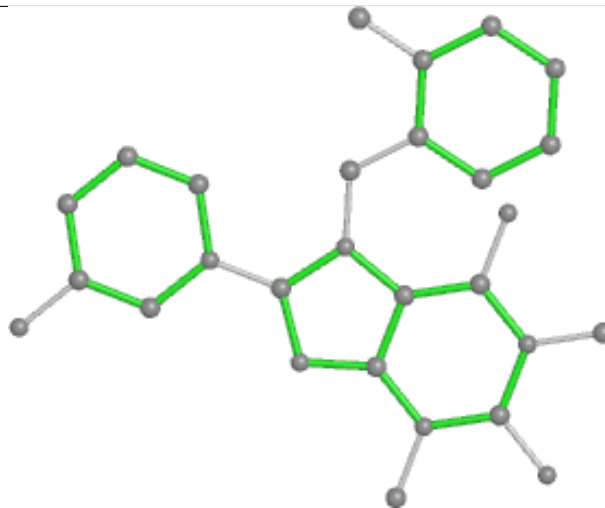
Bond lengths



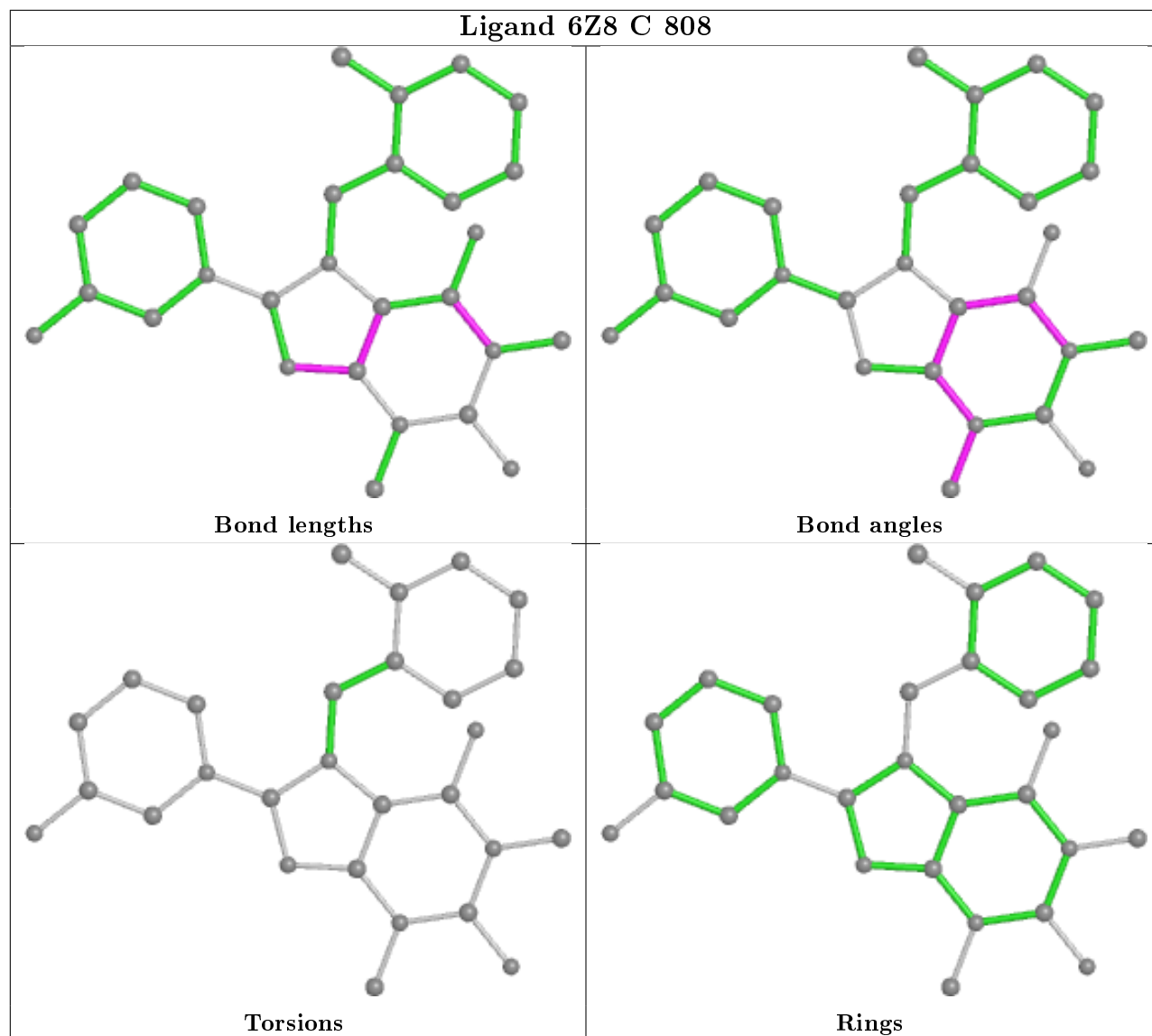
Bond angles

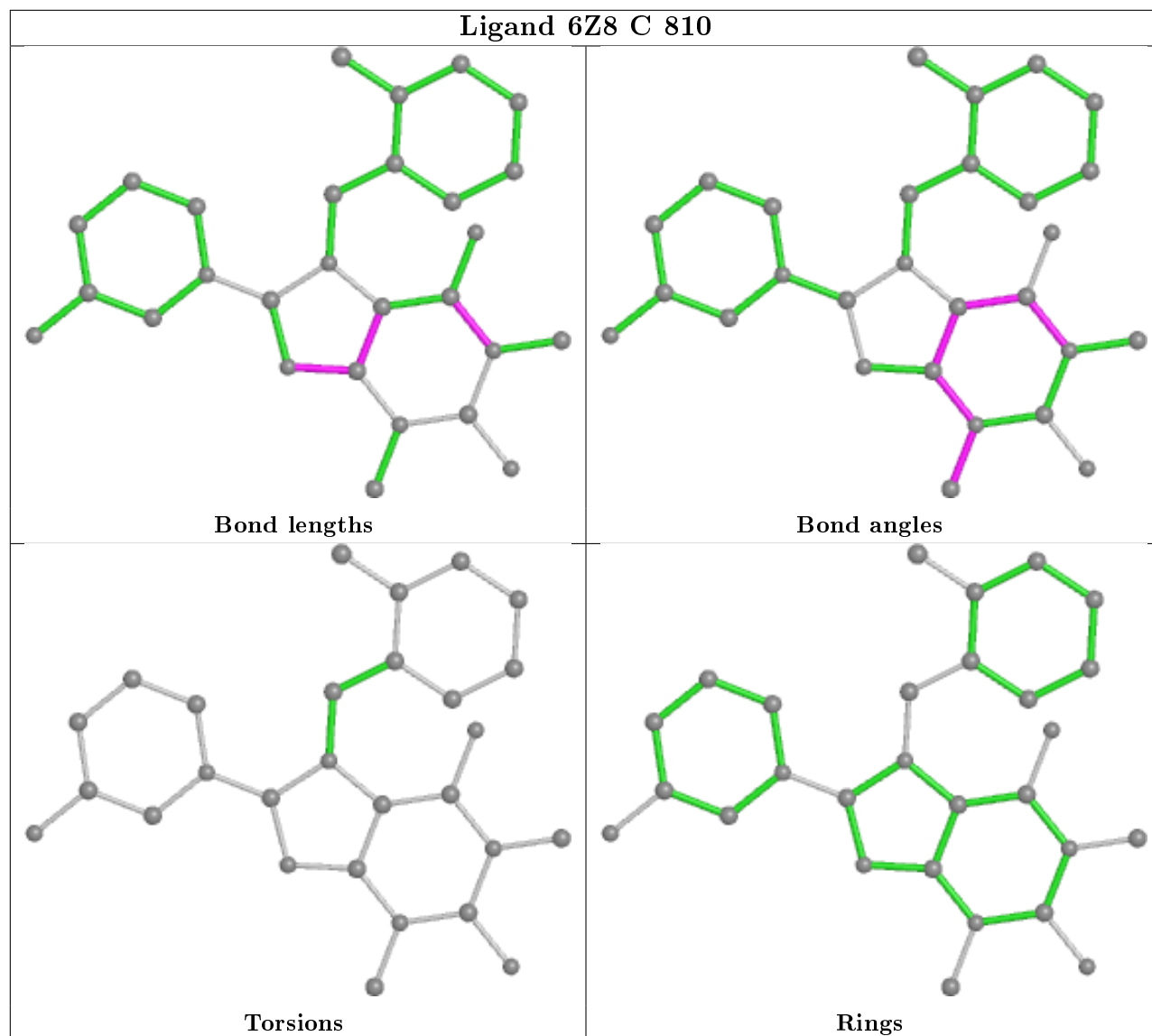


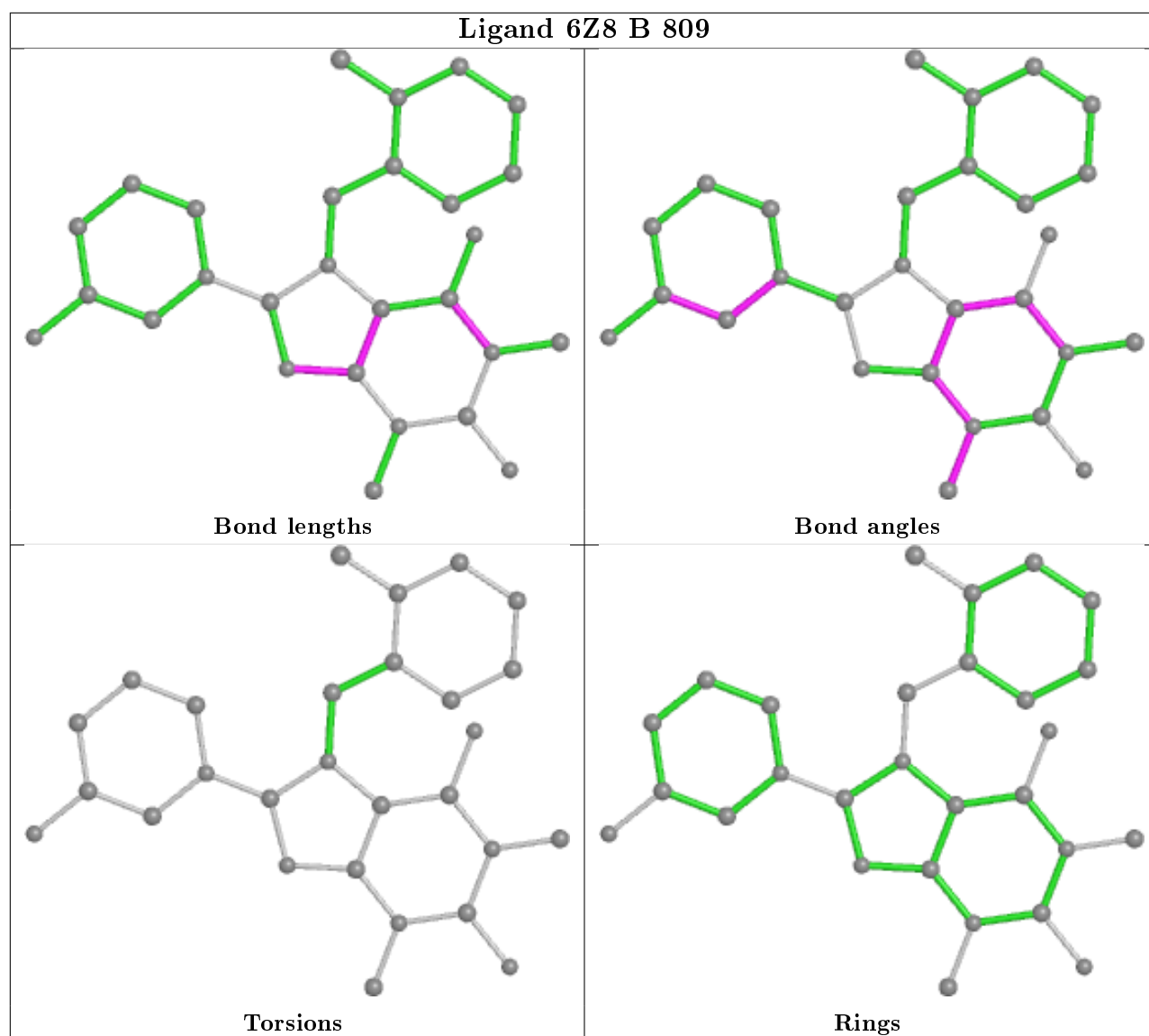
Torsions



Rings







## 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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	727/766 (94%)	0.07	33 (4%) 33 31	18, 34, 70, 121	0
1	B	727/766 (94%)	0.03	32 (4%) 34 32	17, 33, 68, 105	0
1	C	727/766 (94%)	0.16	47 (6%) 18 17	18, 36, 80, 140	0
1	D	727/766 (94%)	0.25	47 (6%) 18 17	17, 40, 79, 132	0
All	All	2908/3064 (94%)	0.13	159 (5%) 25 23	17, 36, 76, 140	0

All (159) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	100	TYR	13.5
1	A	99	GLY	12.8
1	D	100	TYR	12.6
1	A	101	SER	12.3
1	C	100	TYR	10.1
1	B	101	SER	9.1
1	C	98	LEU	8.9
1	C	101	SER	8.7
1	B	103	ASN	8.0
1	D	99	GLY	7.9
1	D	83	TYR	7.5
1	C	99	GLY	7.5
1	C	120	TYR	7.3
1	A	103	ASN	7.3
1	A	98	LEU	7.2
1	B	100	TYR	7.1
1	D	98	LEU	7.1
1	B	102	THR	6.9
1	B	506	ASP	6.9
1	C	96	ASP	6.9
1	C	95	PHE	6.7

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Mol	Chain	Res	Type	RSRZ
1	D	101	SER	6.6
1	D	102	THR	6.2
1	C	103	ASN	6.2
1	C	128	TYR	5.9
1	C	766	PRO	5.9
1	C	76	ILE	5.8
1	B	83	TYR	5.7
1	D	103	ASN	5.7
1	A	123	GLN	5.3
1	C	119	ASN	5.3
1	C	127	SER	5.2
1	A	102	THR	5.1
1	A	122	LYS	5.1
1	B	99	GLY	5.0
1	B	98	LEU	5.0
1	C	121	VAL	4.9
1	C	506	ASP	4.8
1	C	138	ASN	4.7
1	A	82	GLU	4.6
1	A	120	TYR	4.5
1	A	83	TYR	4.4
1	D	91	GLU	4.4
1	C	97	GLU	4.3
1	D	121	VAL	4.3
1	B	138	ASN	4.2
1	C	74	ASN	4.2
1	C	122	LYS	4.2
1	A	121	VAL	4.2
1	A	124	TRP	4.1
1	D	437	ASN	4.1
1	B	535	ASP	4.0
1	C	126	HIS	4.0
1	C	102	THR	4.0
1	A	128	TYR	3.9
1	B	54	ARG	3.9
1	B	121	VAL	3.9
1	D	82	GLU	3.9
1	B	52	THR	3.8
1	C	70	TYR	3.7
1	C	93	SER	3.7
1	B	82	GLU	3.7
1	D	120	TYR	3.7

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Mol	Chain	Res	Type	RSRZ
1	D	145	GLU	3.6
1	C	118	TYR	3.6
1	D	129	THR	3.5
1	B	144	THR	3.5
1	B	120	TYR	3.4
1	A	95	PHE	3.4
1	B	122	LYS	3.4
1	B	145	GLU	3.4
1	C	71	LYS	3.4
1	C	87	SER	3.3
1	B	766	PRO	3.2
1	D	438	ASP	3.2
1	D	138	ASN	3.1
1	D	141	GLN	3.1
1	D	97	GLU	3.1
1	C	130	ALA	3.1
1	D	412	SER	3.0
1	C	91	GLU	3.0
1	A	521	GLY	3.0
1	A	81	ALA	3.0
1	A	85	ASN	3.0
1	D	701	LEU	3.0
1	B	128	TYR	2.9
1	A	125	ARG	2.9
1	A	765	LEU	2.9
1	A	86	SER	2.9
1	B	95	PHE	2.9
1	C	57	PHE	2.9
1	D	125	ARG	2.9
1	D	70	TYR	2.8
1	B	127	SER	2.8
1	C	72	GLN	2.8
1	C	537	SER	2.8
1	A	119	ASN	2.8
1	B	119	ASN	2.8
1	D	95	PHE	2.7
1	D	436	LEU	2.7
1	D	439	TYR	2.7
1	D	144	THR	2.7
1	D	489	LYS	2.7
1	C	129	THR	2.6
1	A	73	GLU	2.6

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Mol	Chain	Res	Type	RSRZ
1	A	57	PHE	2.6
1	C	88	ILE	2.6
1	A	520	HIS	2.6
1	B	118	TYR	2.6
1	C	90	LEU	2.6
1	C	75	ASN	2.6
1	C	125	ARG	2.6
1	D	96	ASP	2.6
1	D	484	SER	2.5
1	B	536	LYS	2.5
1	D	506	ASP	2.5
1	A	127	SER	2.5
1	B	53	PHE	2.5
1	D	488	ASP	2.5
1	D	677	GLU	2.5
1	A	87	SER	2.4
1	C	94	THR	2.4
1	D	93	SER	2.4
1	B	40	ARG	2.4
1	D	379	GLU	2.4
1	D	766	PRO	2.4
1	D	589	LYS	2.4
1	D	55	VAL	2.3
1	A	96	ASP	2.3
1	C	342	ALA	2.3
1	C	439	TYR	2.3
1	A	129	THR	2.3
1	C	83	TYR	2.3
1	A	104	ASP	2.3
1	C	123	GLN	2.3
1	A	151	ASN	2.3
1	D	502	LYS	2.3
1	B	130	ALA	2.3
1	B	618	PHE	2.2
1	D	179	ASN	2.2
1	B	537	SER	2.2
1	A	488	ASP	2.2
1	C	765	LEU	2.2
1	C	131	SER	2.2
1	A	126	HIS	2.2
1	D	622	LYS	2.1
1	C	535	ASP	2.1

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Mol	Chain	Res	Type	RSRZ
1	D	500	LEU	2.1
1	C	414	TYR	2.1
1	D	118	TYR	2.1
1	D	743	ALA	2.1
1	D	119	ASN	2.1
1	D	137	LEU	2.1
1	B	676	PRO	2.0
1	C	50	LYS	2.0
1	D	440	THR	2.0
1	D	628	GLY	2.0
1	B	137	LEU	2.0
1	C	73	GLU	2.0

## 6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

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

## 6.3 Carbohydrates ⓘ

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, 95<sup>th</sup> 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.

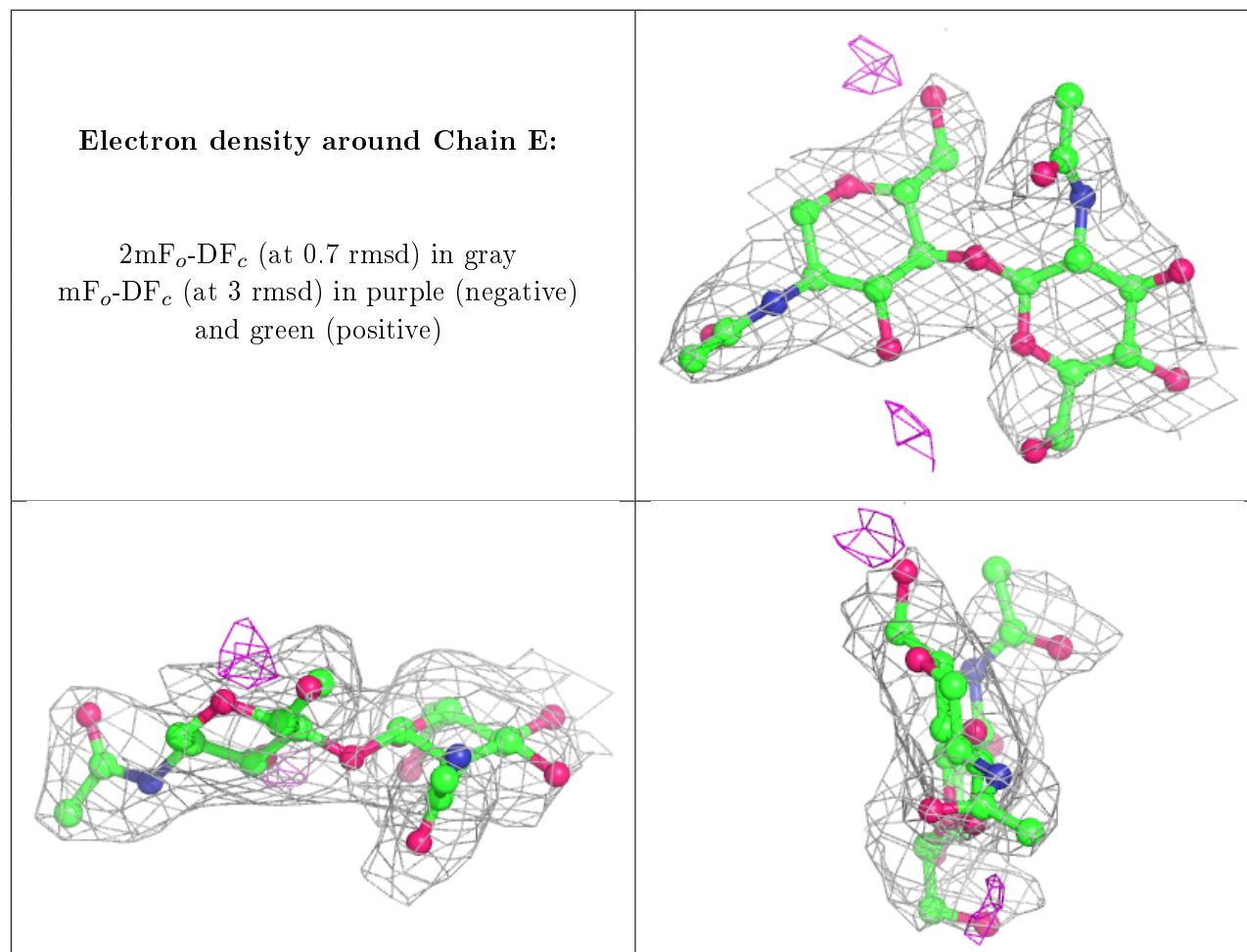
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å <sup>2</sup> )	Q<0.9
2	NAG	J	2	14/15	0.70	0.33	125,130,132,134	0
2	NAG	L	2	14/15	0.73	0.35	74,76,80,82	0
2	NAG	J	1	14/15	0.79	0.20	112,115,118,122	0
2	NAG	K	2	14/15	0.80	0.39	66,74,80,85	0
2	NAG	L	1	14/15	0.83	0.18	64,70,74,74	0
2	NAG	F	1	14/15	0.86	0.18	45,62,70,73	0
2	NAG	F	2	14/15	0.86	0.32	70,77,81,85	0
2	NAG	M	1	14/15	0.88	0.22	41,51,56,63	0
2	NAG	E	2	14/15	0.88	0.21	66,70,77,79	0
2	NAG	M	2	14/15	0.89	0.34	68,71,75,79	0
2	NAG	G	2	14/15	0.90	0.17	57,59,61,61	0
2	NAG	I	2	14/15	0.90	0.32	53,61,64,65	0
2	NAG	K	1	14/15	0.92	0.20	49,58,65,68	0
2	NAG	H	2	14/15	0.92	0.17	61,63,66,67	0
2	NAG	I	1	14/15	0.94	0.20	42,46,51,54	0
2	NAG	E	1	14/15	0.95	0.12	38,44,50,57	0

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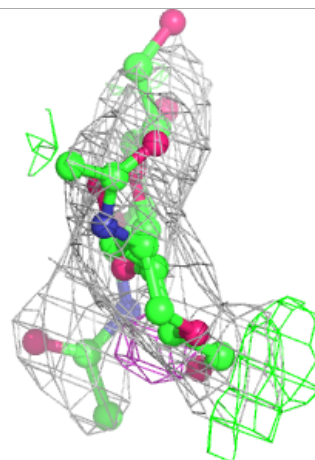
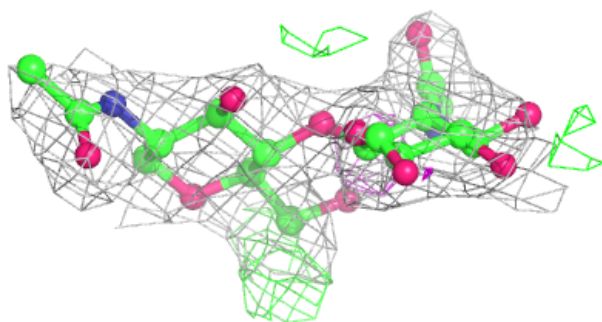
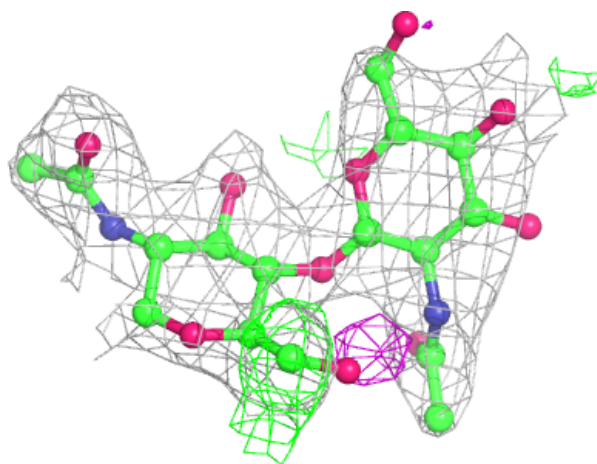
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
2	NAG	H	1	14/15	0.95	0.12	33,41,55,56	0
2	NAG	G	1	14/15	0.95	0.16	48,54,61,64	0

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.



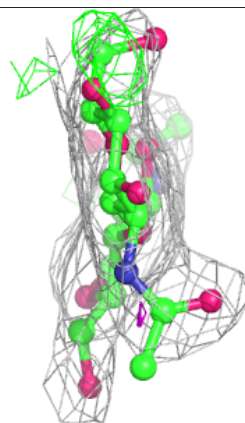
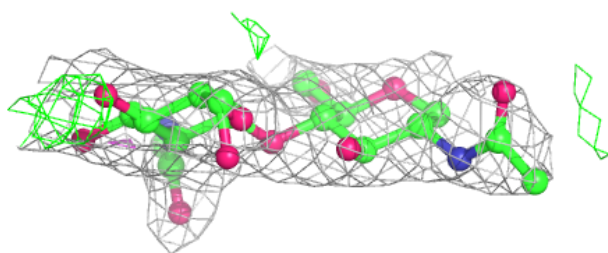
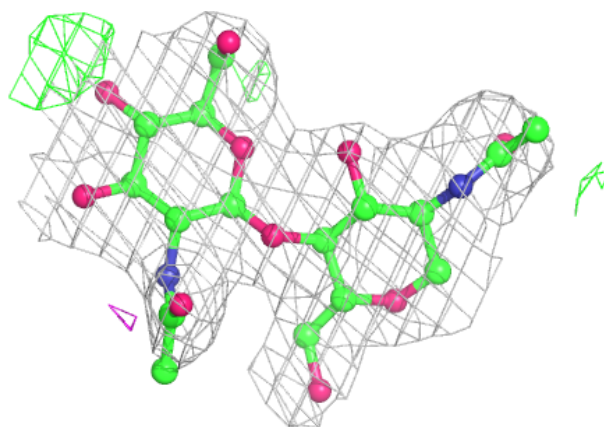
**Electron density around Chain F:**

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



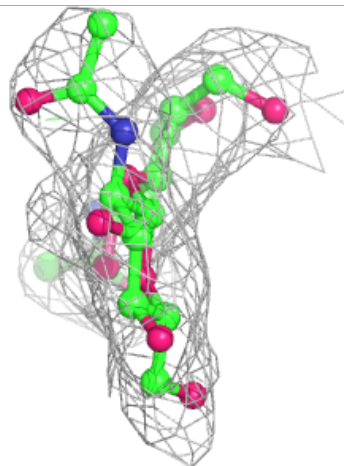
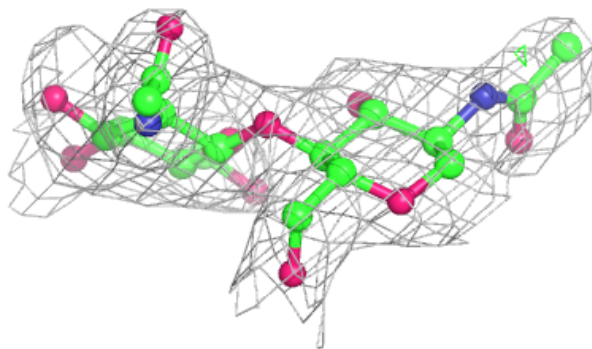
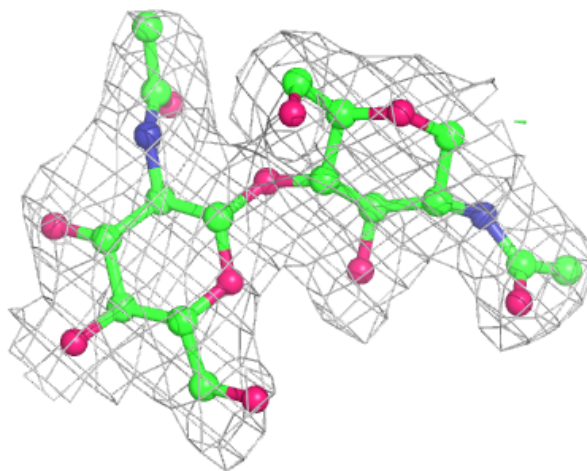
**Electron density around Chain G:**

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



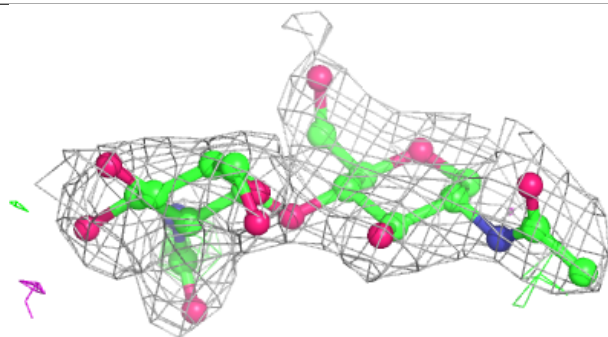
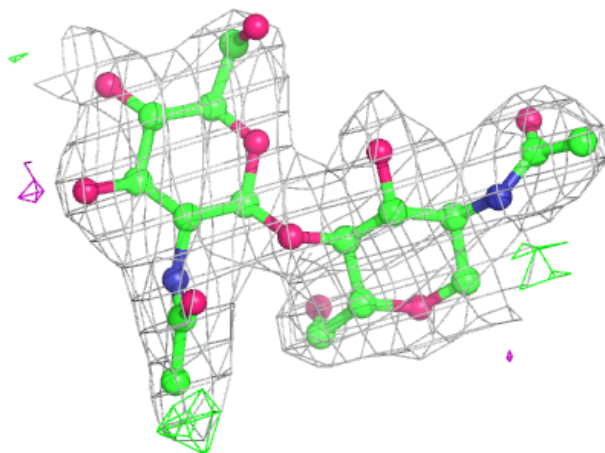
**Electron density around Chain H:**

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



**Electron density around Chain I:**

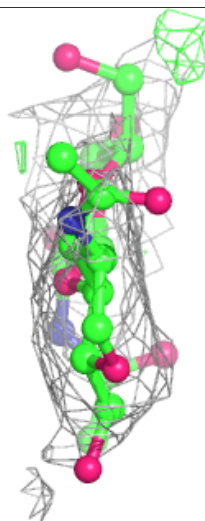
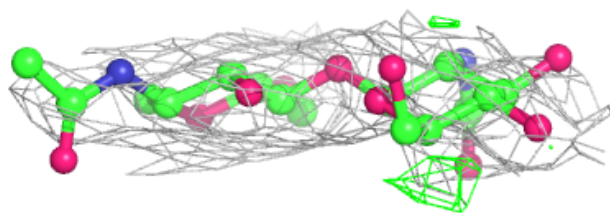
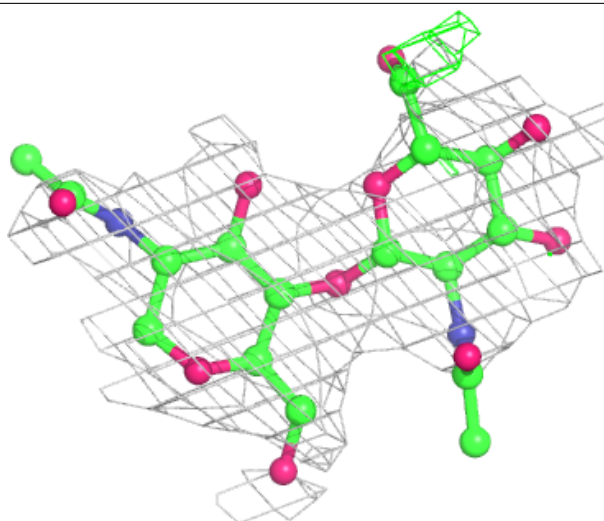
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)





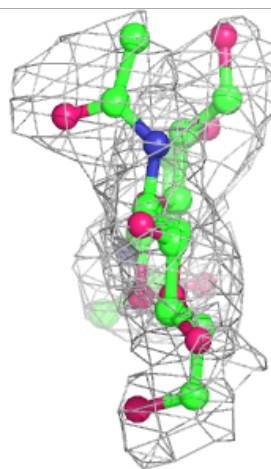
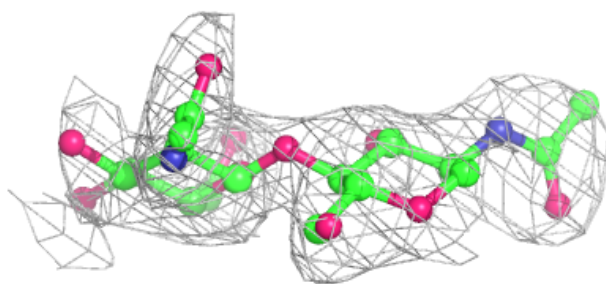
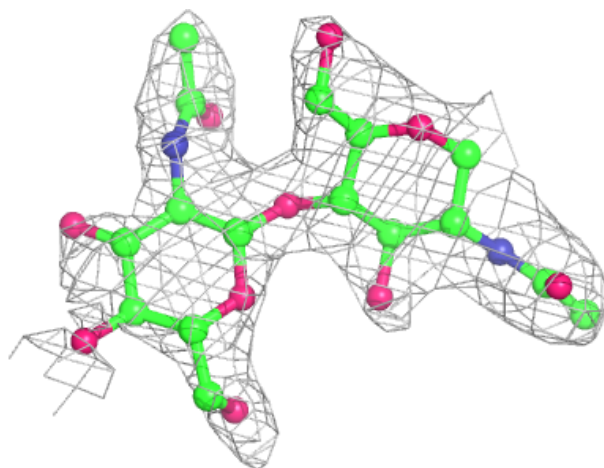
**Electron density around Chain J:**

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



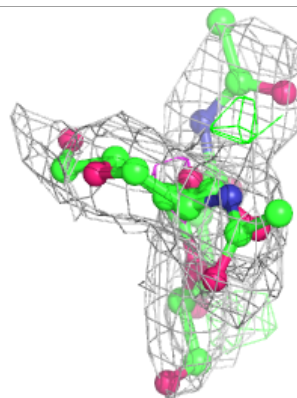
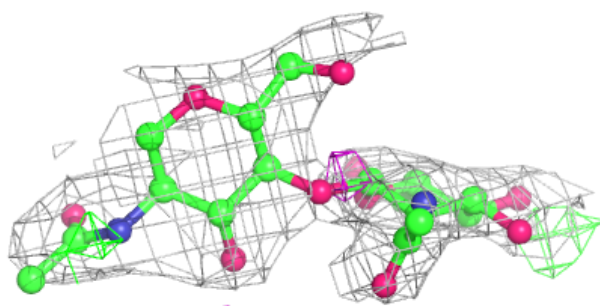
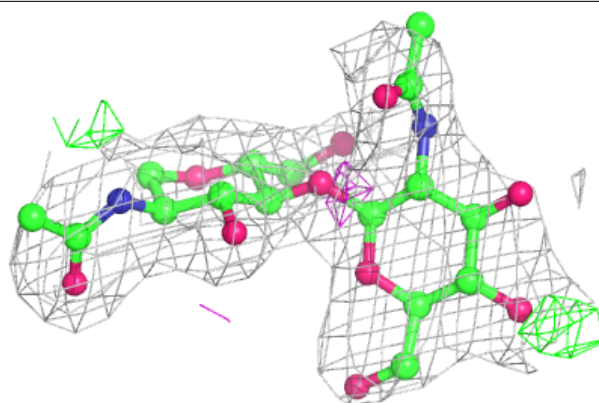
**Electron density around Chain K:**

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

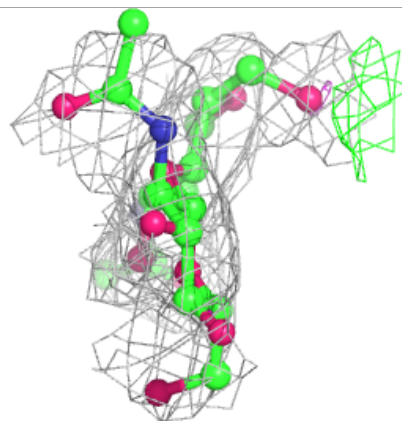
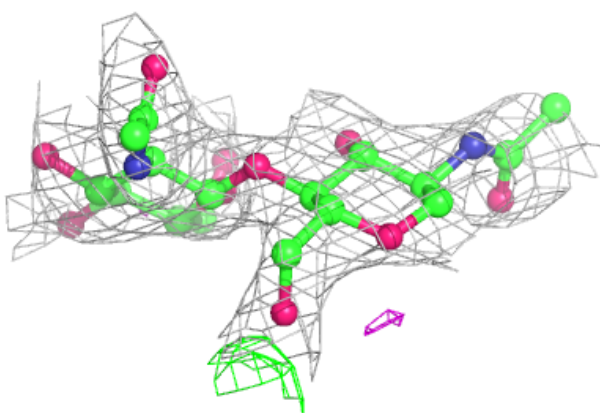
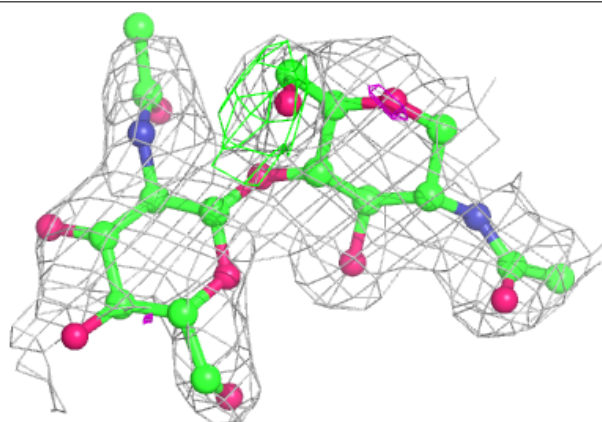


**Electron density around Chain L:**

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

**Electron density around Chain 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 ⓘ

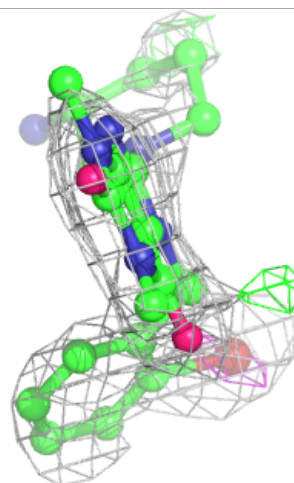
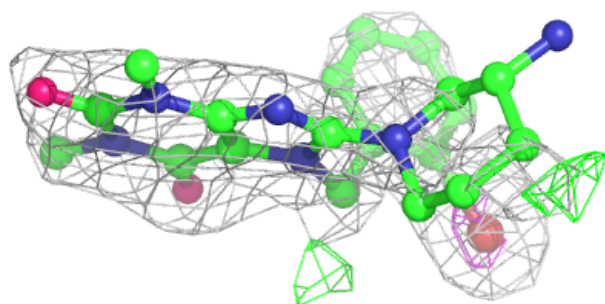
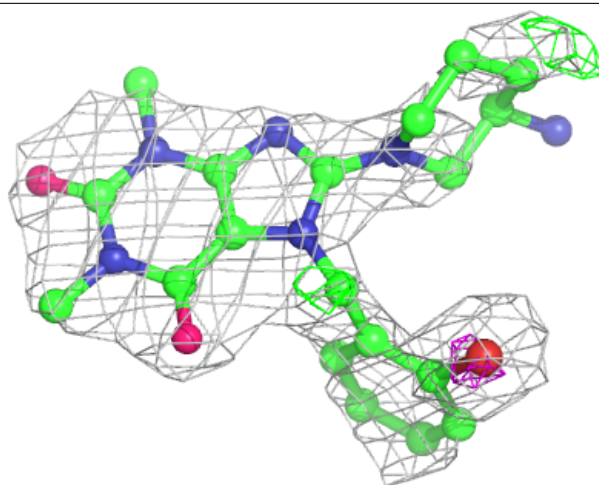
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å <sup>2</sup> )	Q<0.9
3	NAG	D	801	14/15	0.42	0.43	119,127,133,136	0
3	NAG	C	802	14/15	0.66	0.27	122,123,124,124	0
3	NAG	B	801	14/15	0.77	0.27	64,77,86,87	0
4	6Z8	D	813	28/28	0.78	0.29	28,45,58,68	28
3	NAG	A	805	14/15	0.79	0.25	73,75,78,79	0
3	NAG	A	801	14/15	0.79	0.31	83,90,95,96	0
3	NAG	D	806	14/15	0.80	0.28	94,102,107,110	0
3	NAG	B	805	14/15	0.80	0.20	72,74,76,77	0
4	6Z8	A	811	28/28	0.82	0.26	28,35,47,64	28
4	6Z8	B	811	28/28	0.82	0.27	24,42,51,68	28
3	NAG	C	804	14/15	0.82	0.21	49,53,57,57	0
4	6Z8	C	810	28/28	0.84	0.28	38,50,72,85	28
3	NAG	C	801	14/15	0.84	0.19	69,71,73,73	0
3	NAG	B	806	14/15	0.85	0.20	46,51,58,59	0
3	NAG	C	805	14/15	0.85	0.19	54,59,61,61	0
3	NAG	A	802	14/15	0.89	0.21	78,82,88,88	0
3	NAG	B	804	14/15	0.89	0.20	38,43,46,47	0
3	NAG	A	806	14/15	0.91	0.15	47,57,63,65	0
3	NAG	C	803	14/15	0.95	0.14	24,36,42,44	0
4	6Z8	D	811	28/28	0.97	0.15	22,33,38,49	0
4	6Z8	A	809	28/28	0.97	0.13	20,28,37,46	0
5	SO4	C	809	5/5	0.97	0.10	39,44,47,50	0
4	6Z8	C	808	28/28	0.97	0.14	21,29,38,44	0
4	6Z8	B	809	28/28	0.97	0.14	19,28,34,43	0
5	SO4	D	812	5/5	0.97	0.17	58,59,61,62	0
5	SO4	A	810	5/5	0.98	0.11	37,38,42,42	0
5	SO4	B	810	5/5	0.98	0.10	57,57,58,60	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.

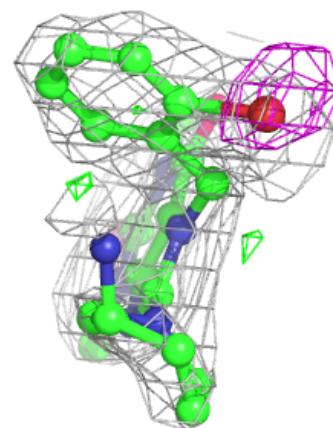
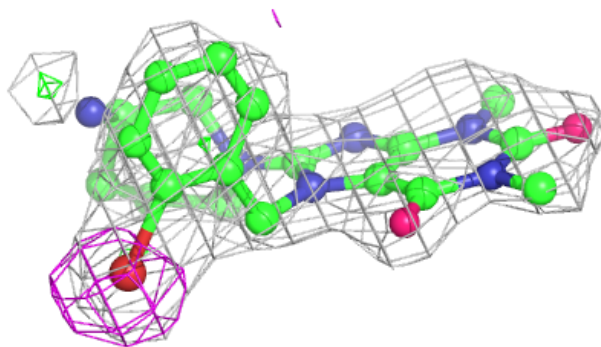
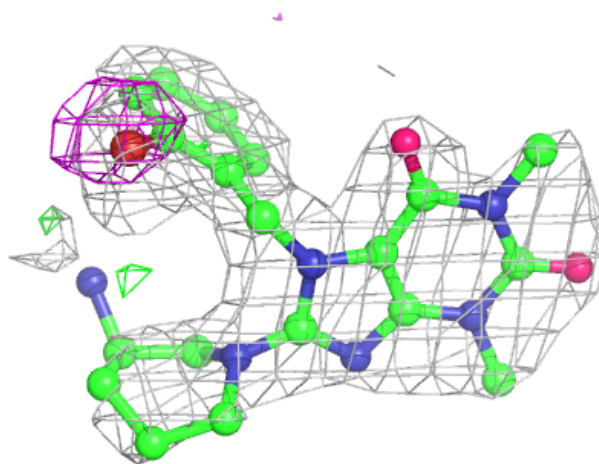
**Electron density around 6Z8 D 813:**

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



**Electron density around 6Z8 A 811:**

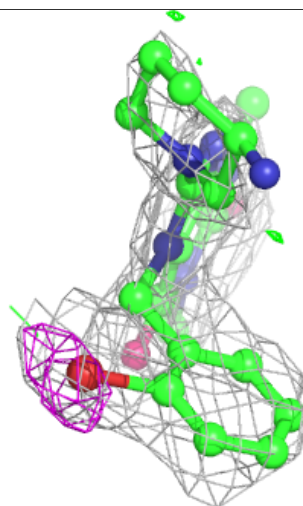
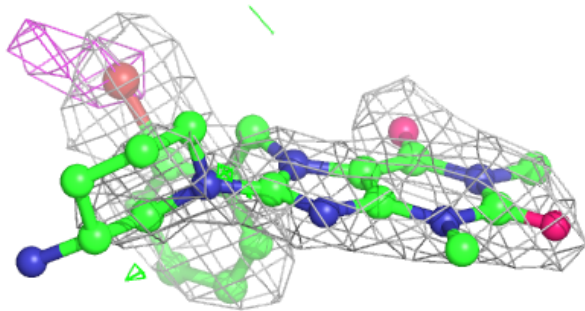
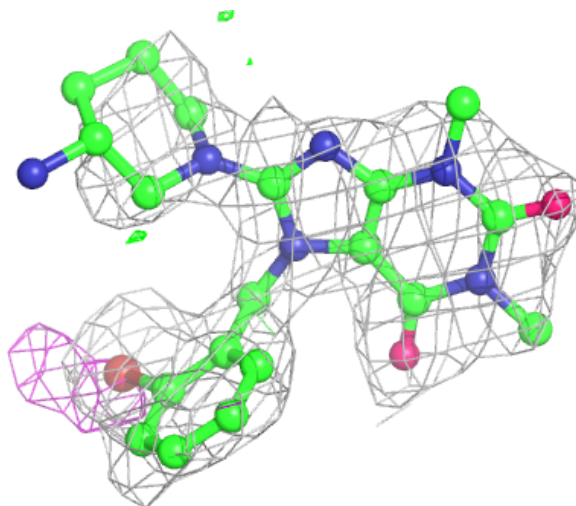
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)





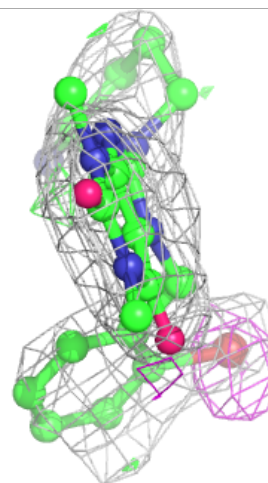
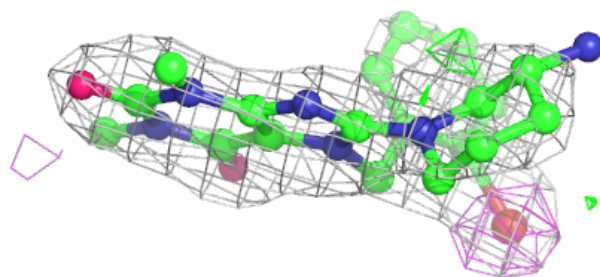
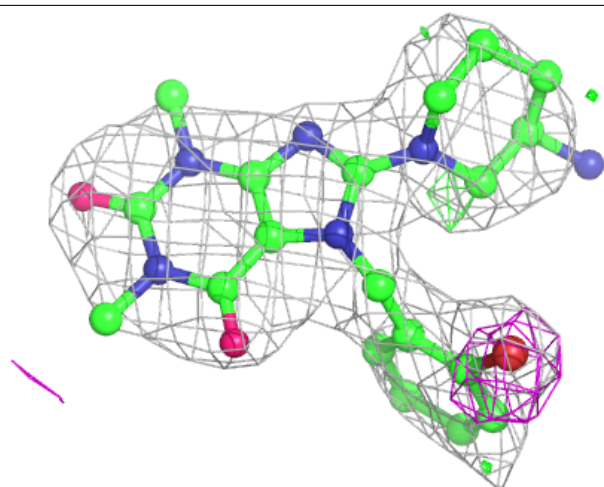
**Electron density around 6Z8 B 811:**

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



**Electron density around 6Z8 C 810:**

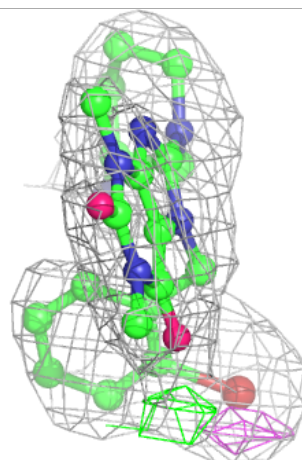
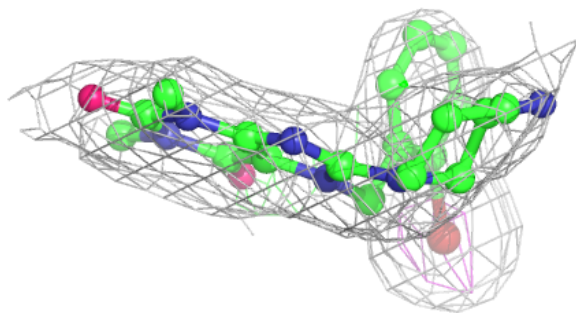
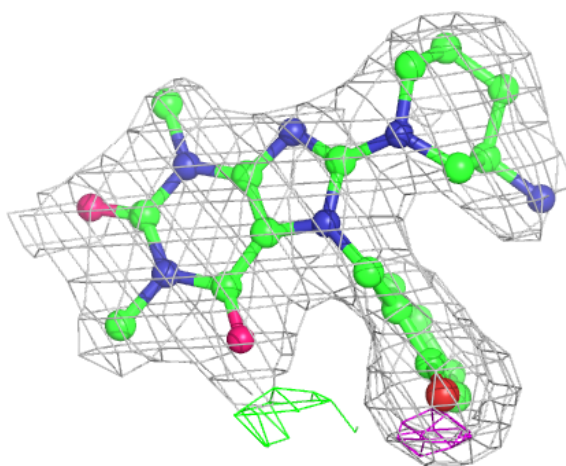
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)





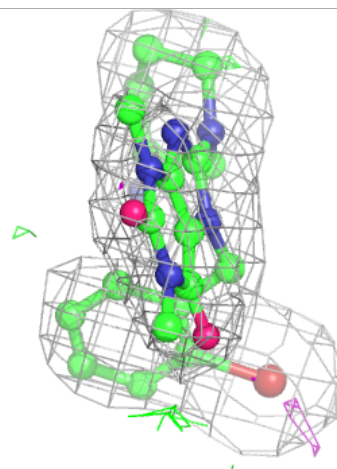
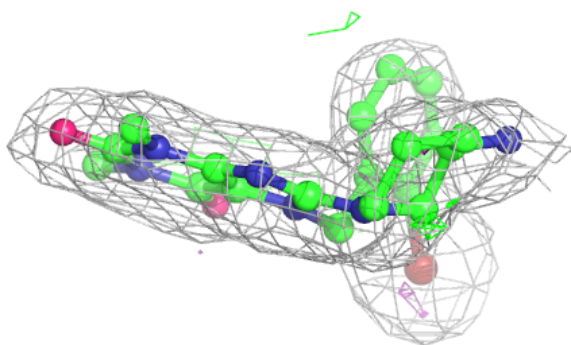
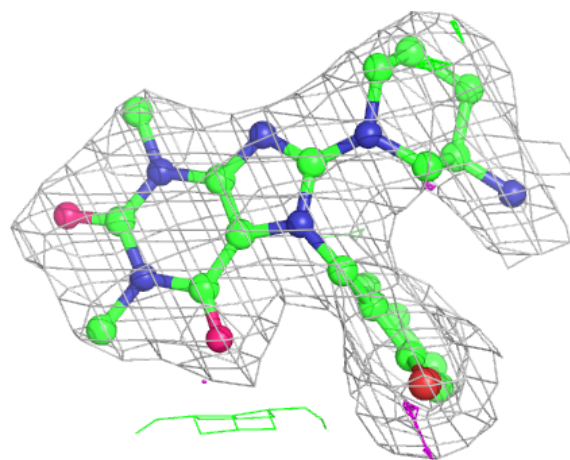
**Electron density around 6Z8 D 811:**

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



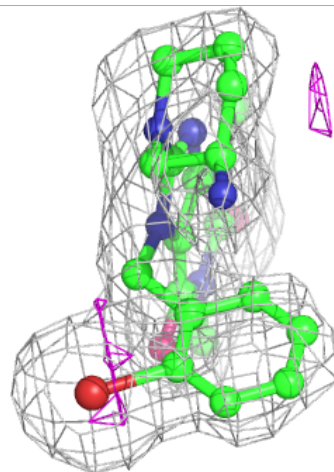
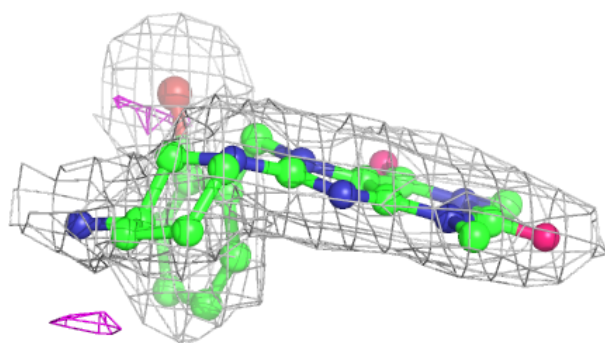
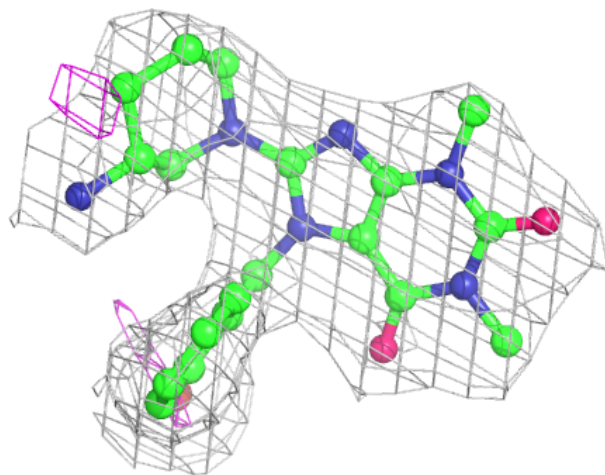
**Electron density around 6Z8 A 809:**

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



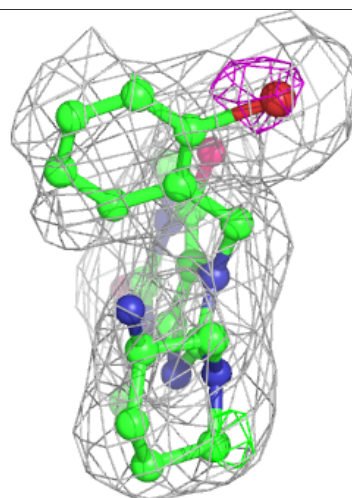
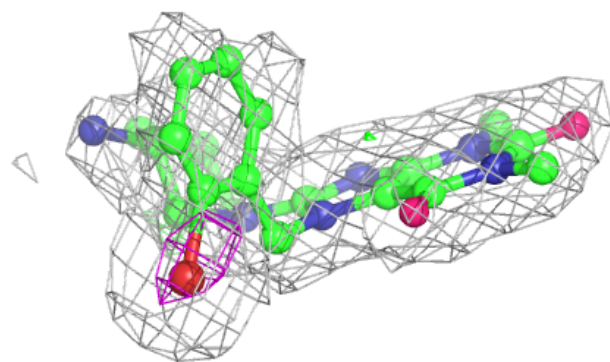
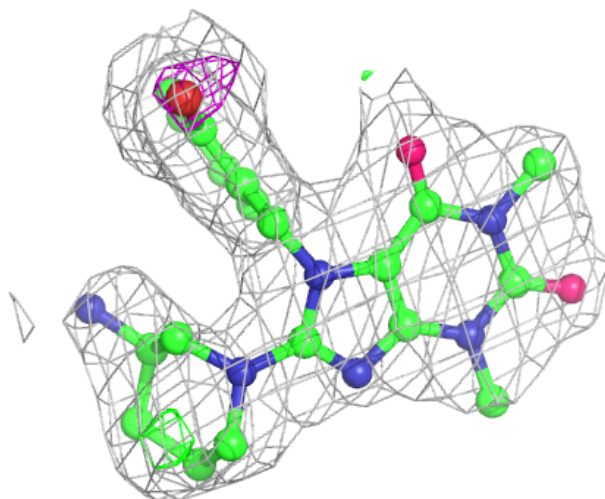
**Electron density around 6Z8 C 808:**

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



**Electron density around 6Z8 B 809:**

$2mF_o-DF_c$  (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.