



# Full wwPDB X-ray Structure Validation Report ⓘ

Aug 6, 2020 – 10:57 AM BST

PDB ID : 6ESY  
Title : Human butyrylcholinesterase in complex with thioflavine T  
Authors : Nachon, F.; Brazzolotto, X.; Wandhammer, M.; Trovaslet-Leroy, M.; Macdonald, I.R.; Darvesh, S.; Rosenberry, T.L.  
Deposited on : 2017-10-24  
Resolution : 2.80 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

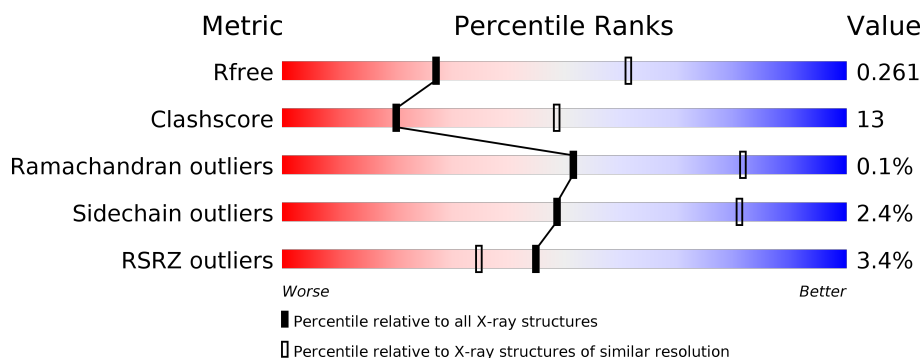
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.80 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.



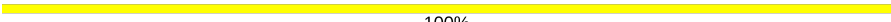
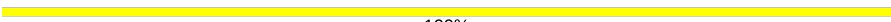
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
$R_{free}$	130704	3140 (2.80-2.80)
Clashscore	141614	3569 (2.80-2.80)
Ramachandran outliers	138981	3498 (2.80-2.80)
Sidechain outliers	138945	3500 (2.80-2.80)
RSRZ outliers	127900	3078 (2.80-2.80)

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	529	<div> <div>4%</div> <div> <div></div> <div>74%</div> <div>25%</div> </div> <div></div> </div>
1	B	529	<div> <div>3%</div> <div> <div></div> <div>72%</div> <div>26%</div> </div> <div></div> </div>
2	C	3	<div> <div>33%</div> <div>33%</div> <div>33%</div> </div>
2	H	3	<div> <div>33%</div> <div>33%</div> <div>33%</div> </div>
3	D	2	<div> <div>50%</div> <div>50%</div> </div>
3	E	2	<div> <div>50%</div> <div>50%</div> </div>

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Mol	Chain	Length	Quality of chain
3	G	2	 100%
4	F	2	 100%

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
2	NAG	C	2	-	-	-	X
2	FUC	C	3	X	-	-	X
2	FUC	H	3	X	-	-	-
4	FUC	F	2	X	-	-	-
6	TFX	B	611	-	-	X	-
6	TFX	B	612	-	-	X	-
7	CL	A	614	-	-	-	X

## 2 Entry composition [i](#)

There are 8 unique types of molecules in this entry. The entry contains 8775 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 Cholinesterase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	527	Total	C	N	O	S	3	3	0
			4204	2713	709	767	15			
1	B	526	Total	C	N	O	S	9	0	0
			4166	2695	702	754	15			

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



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

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



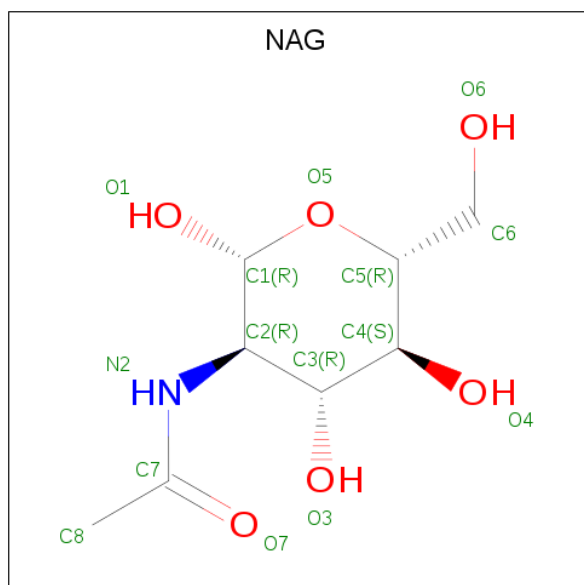
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	D	2	Total	C	N	O	0	0	0
			28	16	2	10			
3	E	2	Total	C	N	O	0	0	0
			28	16	2	10			
3	G	2	Total	C	N	O	0	0	0
			28	16	2	10			

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



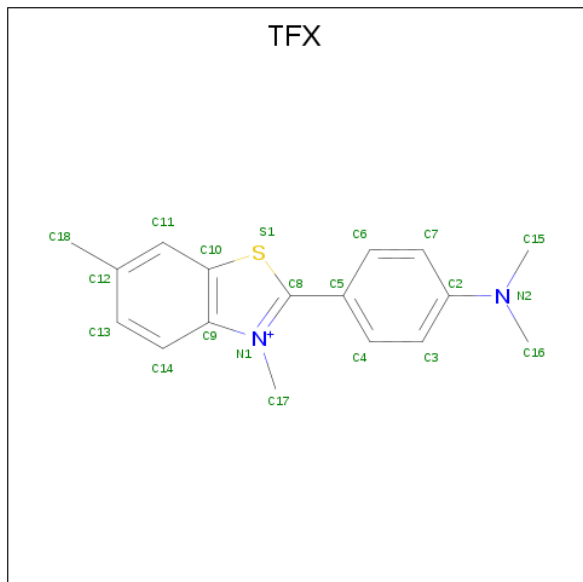
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
4	F	2	Total	C	N	O	0	0	0
			24	14	1	9			

- Molecule 5 is 2-acetamido-2-deoxy-beta-D-glucopyranose (three-letter code: NAG) (formula:  $C_8H_{15}NO_6$ ).



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

- Molecule 6 is 2-[4-(dimethylamino)phenyl]-3,6-dimethyl-1,3-benzothiazol-3-ium (three-letter code: TFX) (formula: C<sub>17</sub>H<sub>19</sub>N<sub>2</sub>S) (labeled as "Ligand of Interest" by author).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
6	A	1	Total	C	N	S	0	0
			20	17	2	1		
6	A	1	Total	C	N	S	0	0
			20	17	2	1		
6	B	1	Total	C	N	S	0	0
			20	17	2	1		
6	B	1	Total	C	N	S	0	0
			20	17	2	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		

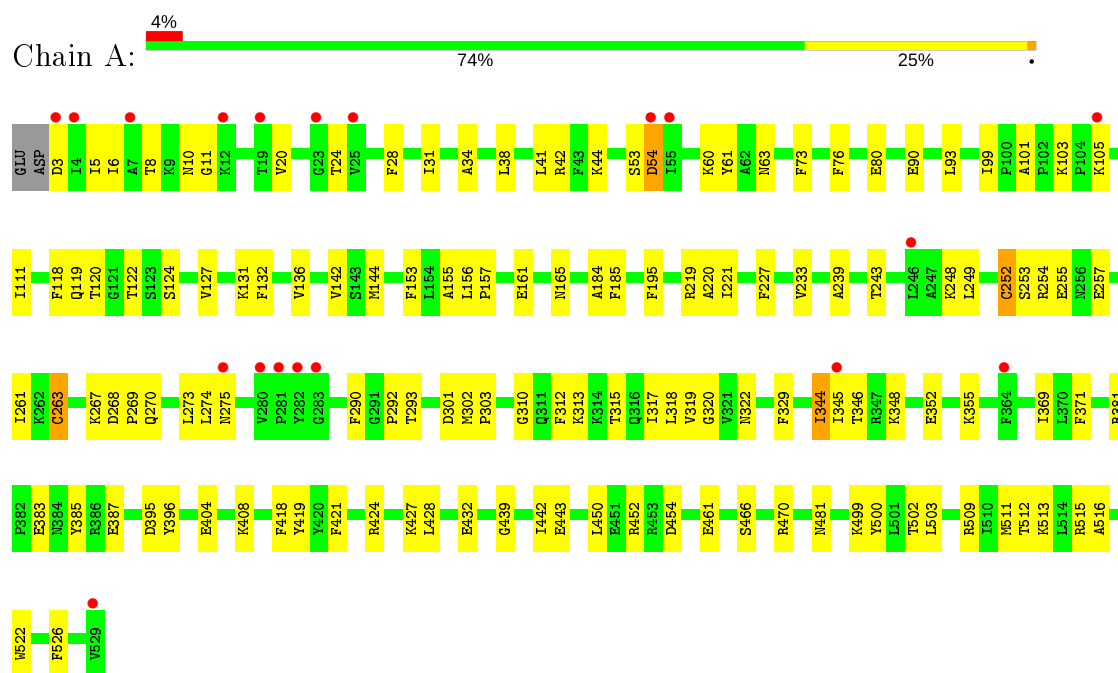
- Molecule 8 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	17	Total	O	0	0
			17	17		
8	B	24	Total	O	0	0
			24	24		

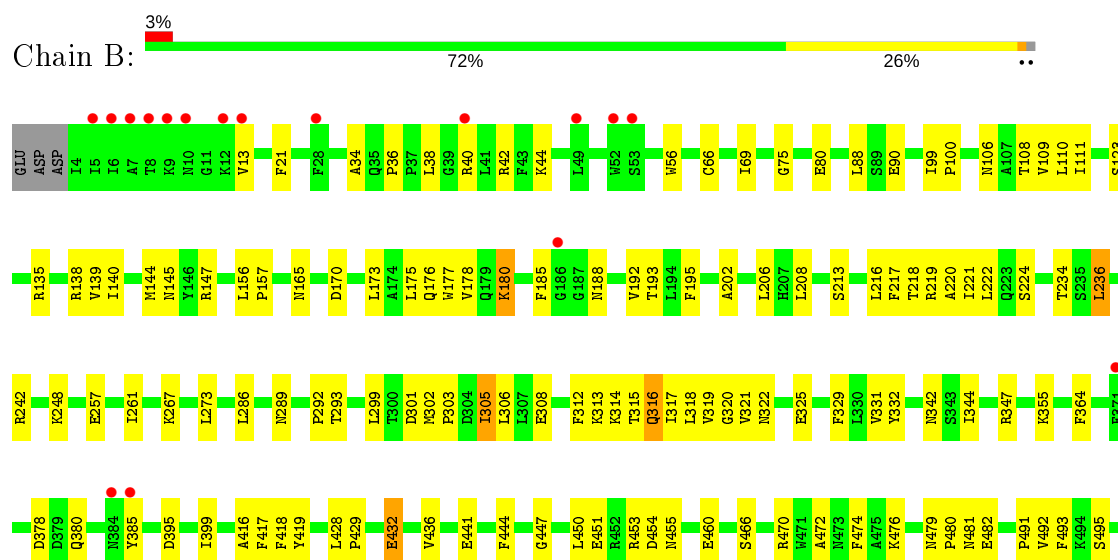
### 3 Residue-property plots

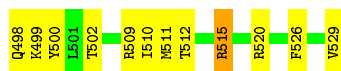
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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: Cholinesterase



#### • Molecule 1: Cholinesterase





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

Chain C: 33% 33% 33%



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

Chain H: 33% 33% 33%



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

Chain D: 50% 50%



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

Chain E: 50% 50%



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

Chain G: 100%



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

Chain F: 100%





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	74.27Å 79.25Å 228.92Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	39.35 – 2.80 48.98 – 2.80	Depositor EDS
% Data completeness (in resolution range)	98.6 (39.35-2.80) 92.5 (48.98-2.80)	Depositor EDS
$R_{merge}$	0.08	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.11 (at 2.81Å)	Xtriage
Refinement program	PHENIX (1.12_2829: ???)	Depositor
R, $R_{free}$	0.206 , 0.262 0.206 , 0.261	Depositor DCC
$R_{free}$ test set	1635 reflections (4.85%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	68.2	Xtriage
Anisotropy	0.404	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.32 , 58.6	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.46$ , $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	8775	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	83.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.57% 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 ⓘ

### 5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: CL, NAG, FUC, TFX

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	0.31	0/4333	0.52	1/5885 (0.0%)
1	B	0.34	1/4285 (0.0%)	0.56	5/5820 (0.1%)
All	All	0.33	1/8618 (0.0%)	0.54	6/11705 (0.1%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	455	ASN	CB-CG	7.69	1.68	1.51

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	106	ASN	N-CA-CB	7.64	124.36	110.60
1	B	106	ASN	CB-CA-C	-7.61	95.17	110.40
1	B	236	LEU	CB-CG-CD1	-6.91	99.26	111.00
1	B	106	ASN	CA-CB-CG	5.99	126.59	113.40
1	B	222	LEU	CB-CG-CD2	-5.81	101.12	111.00
1	A	344	ILE	CB-CG1-CD1	-5.30	99.06	113.90

There are no chirality outliers.

There are no planarity outliers.

### 5.2 Too-close contacts ⓘ

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	4204	0	4094	101	1
1	B	4166	0	4044	100	1
2	C	38	0	34	1	0
2	H	38	0	34	1	0
3	D	28	0	25	0	0
3	E	28	0	25	0	0
3	G	28	0	25	0	0
4	F	24	0	22	0	0
5	A	28	0	26	1	0
5	B	70	0	65	4	0
6	A	40	0	38	6	0
6	B	40	0	38	9	0
7	A	1	0	0	0	0
7	B	1	0	0	0	0
8	A	17	0	0	1	0
8	B	24	0	0	1	0
All	All	8775	0	8470	215	1

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 13.

All (215) 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:8:THR:HG22	1:A:10:ASN:H	1.22	0.98
1:A:381:ARG:NH2	1:A:383:GLU:OE1	1.97	0.97
1:A:249:LEU:HB3	1:A:275:ASN:HD22	1.33	0.92
1:A:381:ARG:HH21	1:A:383:GLU:CD	1.75	0.88
6:A:612:TFX:H17B	6:A:613:TFX:H6	1.55	0.87
1:A:165:ASN:HD21	1:A:293:THR:H	1.20	0.85
1:A:127:VAL:HB	1:A:442:ILE:HD13	1.58	0.83
1:B:40:ARG:O	1:B:44:LYS:NZ	2.15	0.79
1:A:263:CYS:SG	1:A:267:LYS:NZ	2.54	0.78
1:B:417:PHE:CE1	1:B:480:PRO:HD2	2.24	0.72
1:A:345:ILE:HD11	1:A:385:TYR:HB3	1.71	0.72
1:A:99:ILE:HD11	1:A:185:PHE:HB3	1.72	0.72
1:B:502:THR:HG21	1:B:509:ARG:HH11	1.55	0.71
1:A:243:THR:HG22	1:A:290:PHE:HE2	1.57	0.69
1:B:317:ILE:HD13	1:B:319:VAL:HG23	1.76	0.68
1:A:249:LEU:HB3	1:A:275:ASN:ND2	2.09	0.66
1:A:500:TYR:CZ	1:A:511:MET:HG3	2.31	0.65
1:B:110:LEU:HD23	1:B:193:THR:HB	1.78	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:344:ILE:HD12	1:B:344:ILE:H	1.60	0.64
1:B:176:GLN:HG2	1:B:180:LYS:HE2	1.80	0.64
6:B:611:TFX:H17B	6:B:612:TFX:H4	1.79	0.64
1:A:270:GLN:HG3	1:A:274:LEU:HD21	1.80	0.63
1:B:108:THR:HG21	1:B:476:LYS:HG2	1.80	0.63
1:B:417:PHE:CE2	1:B:492:VAL:HG22	2.32	0.63
1:B:474:PHE:HB2	1:B:480:PRO:HD3	1.79	0.63
1:B:470:ARG:HB3	1:B:480:PRO:HB2	1.81	0.62
1:B:109:VAL:HG23	1:B:140:ILE:HB	1.81	0.62
1:A:381:ARG:NH2	1:A:383:GLU:CG	2.63	0.61
6:A:612:TFX:H17B	6:A:613:TFX:C6	2.28	0.61
1:B:99:ILE:HD11	1:B:185:PHE:HB3	1.82	0.61
1:B:176:GLN:O	1:B:180:LYS:HG2	2.01	0.60
2:C:2:NAG:H83	2:C:2:NAG:H3	1.82	0.60
1:B:66:CYS:HB3	1:B:273:LEU:HD11	1.83	0.60
1:A:165:ASN:HD21	1:A:293:THR:N	1.96	0.60
1:A:99:ILE:CD1	1:A:185:PHE:HB3	2.32	0.60
1:A:60:LYS:HD3	1:A:61:TYR:N	2.17	0.60
1:B:378:ASP:OD1	1:B:380:GLN:N	2.35	0.59
1:B:417:PHE:CD1	1:B:480:PRO:HD2	2.37	0.59
1:B:466:SER:O	1:B:470:ARG:HG3	2.02	0.59
1:B:13:VAL:HG12	1:B:56:TRP:HB3	1.85	0.59
1:B:109:VAL:HG12	1:B:192:VAL:HG22	1.85	0.58
1:A:255:GLU:N	1:A:255:GLU:OE1	2.34	0.58
1:A:381:ARG:NH2	1:A:383:GLU:HG3	2.18	0.58
1:B:218:THR:O	1:B:315:THR:HG21	2.04	0.58
5:B:607:NAG:O3	5:B:607:NAG:H82	2.02	0.58
1:B:213:SER:HA	1:B:216:LEU:HD12	1.84	0.58
1:B:202:ALA:O	1:B:206:LEU:HD12	2.03	0.58
1:A:253:SER:O	1:A:254:ARG:HD3	2.04	0.57
1:B:331:VAL:HG23	1:B:332:TYR:CD1	2.39	0.57
1:B:416:ALA:C	1:B:417:PHE:HD2	2.08	0.57
1:A:161:GLU:HB2	1:A:261:ILE:HD13	1.86	0.57
1:B:495:SER:O	1:B:498:GLN:NE2	2.28	0.57
1:B:75:GLY:N	1:B:80:GLU:OE1	2.34	0.57
1:A:269:PRO:O	1:A:273:LEU:HD13	2.04	0.57
1:A:381:ARG:HH21	1:A:383:GLU:CG	2.17	0.57
1:A:352:GLU:HA	1:A:355:LYS:HD2	1.86	0.56
2:H:1:NAG:H83	2:H:1:NAG:H3	1.87	0.56
1:B:416:ALA:O	1:B:493:PHE:N	2.27	0.55
1:A:233:VAL:HG21	1:A:303:PRO:HG2	1.88	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:312:PHE:O	1:B:314:LYS:HD3	2.07	0.55
1:A:404:GLU:HG3	1:A:408:LYS:HE3	1.87	0.55
1:B:218:THR:O	1:B:313:LYS:NZ	2.40	0.55
1:A:442:ILE:HD12	1:A:443:GLU:N	2.21	0.55
1:B:69:ILE:HD11	1:B:88:LEU:HD11	1.89	0.55
1:A:344:ILE:HD13	1:A:383:GLU:HA	1.88	0.54
1:A:381:ARG:NH2	1:A:383:GLU:CD	2.55	0.54
1:B:491:PRO:HG3	1:B:510:ILE:HD11	1.89	0.54
6:B:611:TFX:H17B	6:B:612:TFX:C4	2.37	0.54
1:B:302:MET:HB3	1:B:305:ILE:HD12	1.89	0.54
1:B:286:LEU:HD23	6:B:611:TFX:H16B	1.88	0.54
1:A:270:GLN:O	1:A:273:LEU:N	2.40	0.54
1:A:127:VAL:HB	1:A:442:ILE:CD1	2.35	0.54
1:B:474:PHE:CD1	1:B:480:PRO:HD3	2.42	0.53
5:B:606:NAG:O3	5:B:606:NAG:H82	2.08	0.53
1:A:301:ASP:OD1	1:A:302:MET:N	2.42	0.53
1:A:221:ILE:HG12	1:A:318:LEU:HB3	1.90	0.53
1:B:135:ARG:HD2	1:B:138:ARG:NH2	2.24	0.53
1:A:8:THR:HB	1:A:11:GLY:O	2.09	0.52
1:A:60:LYS:HE2	1:A:60:LYS:HA	1.90	0.52
6:B:611:TFX:H18A	6:B:612:TFX:H18A	1.90	0.52
1:A:239:ALA:O	1:A:243:THR:HG23	2.10	0.52
1:A:439:GLY:O	1:A:442:ILE:HG13	2.10	0.52
1:B:491:PRO:HG3	1:B:510:ILE:CD1	2.40	0.52
1:B:135:ARG:HD2	1:B:138:ARG:HH21	1.74	0.52
1:B:42:ARG:HD2	1:B:90:GLU:OE1	2.09	0.52
1:B:108:THR:HG23	1:B:139:VAL:HG13	1.91	0.52
1:B:21:PHE:HE1	1:B:451:GLU:HB2	1.74	0.52
1:B:177:TRP:HA	1:B:180:LYS:HE3	1.91	0.51
1:B:321:VAL:HG11	1:B:399:ILE:HA	1.92	0.51
1:B:432:GLU:H	1:B:432:GLU:CD	2.14	0.51
1:A:155:ALA:H	1:A:243:THR:HG21	1.76	0.51
1:A:53:SER:HB2	1:A:54:ASP:OD1	2.11	0.51
6:B:611:TFX:C17	6:B:612:TFX:H4	2.40	0.51
1:A:132:PHE:O	1:A:136:VAL:HG23	2.11	0.51
1:A:119:GLN:NE2	1:A:290:PHE:O	2.42	0.51
1:B:395:ASP:OD1	1:B:515:ARG:NH2	2.38	0.50
1:A:20:VAL:HG11	1:A:131:LYS:CG	2.42	0.50
1:A:252:CYS:SG	1:A:267:LYS:NZ	2.84	0.50
6:A:612:TFX:C17	6:A:613:TFX:H6	2.36	0.50
1:A:103:LYS:NZ	1:A:184:ALA:O	2.43	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:157:PRO:HB3	1:B:236:LEU:HD11	1.94	0.49
1:B:221:ILE:HG23	1:B:318:LEU:HB3	1.95	0.49
1:A:344:ILE:CD1	1:A:383:GLU:HA	2.41	0.49
1:A:383:GLU:O	1:A:387:GLU:HG3	2.12	0.49
1:A:395:ASP:OD2	1:A:515:ARG:NH2	2.29	0.49
1:B:322:ASN:HB3	1:B:436:VAL:HG12	1.93	0.49
1:A:345:ILE:HD12	1:A:346:THR:O	2.13	0.48
1:A:313:LYS:NZ	8:A:702:HOH:O	2.45	0.48
1:A:424:ARG:HG3	1:A:428:LEU:HD23	1.95	0.48
1:A:466:SER:O	1:A:470:ARG:HG3	2.14	0.48
1:B:123:SER:HB3	1:B:145:ASN:OD1	2.13	0.48
1:B:165:ASN:OD1	1:B:292:PRO:HA	2.13	0.48
1:A:322:ASN:HD22	1:A:421:PHE:HB3	1.78	0.48
1:A:424:ARG:NH1	1:A:432:GLU:HA	2.29	0.48
1:B:319:VAL:O	1:B:418:PHE:HA	2.14	0.48
1:B:417:PHE:HE1	1:B:480:PRO:HD2	1.74	0.48
1:B:499:LYS:HA	1:B:511:MET:O	2.14	0.48
1:B:470:ARG:HG2	1:B:482:GLU:HB2	1.96	0.48
1:A:24:THR:HG21	5:A:601:NAG:H61	1.96	0.47
1:B:111:ILE:HD11	1:B:178:VAL:HG11	1.96	0.47
1:B:100:PRO:HB3	1:B:138:ARG:HD3	1.95	0.47
1:A:111:ILE:HA	1:A:142:VAL:O	2.14	0.47
1:A:344:ILE:HA	1:A:344:ILE:HD12	1.41	0.47
1:B:147:ARG:NH2	1:B:170:ASP:OD1	2.46	0.47
1:A:369:ILE:HD11	1:A:526:PHE:CE1	2.50	0.47
1:A:120:THR:OG1	6:A:612:TFX:H17A	2.14	0.47
1:B:219:ARG:HA	1:B:315:THR:HG21	1.96	0.47
1:B:502:THR:HG21	1:B:509:ARG:NH1	2.27	0.47
1:B:42:ARG:NH2	1:B:267:LYS:O	2.36	0.47
1:B:474:PHE:HD1	1:B:480:PRO:HD3	1.78	0.47
1:A:509:ARG:CZ	1:A:511:MET:HE1	2.45	0.47
1:A:227:PHE:CD1	1:A:303:PRO:HB2	2.50	0.47
1:A:156:LEU:HD13	1:A:257:GLU:HB3	1.96	0.46
1:B:441:GLU:OE2	1:B:441:GLU:N	2.38	0.46
1:A:6:ILE:HD11	1:A:184:ALA:C	2.36	0.46
1:A:319:VAL:O	1:A:418:PHE:HA	2.16	0.46
1:A:60:LYS:HD3	1:A:61:TYR:H	1.79	0.46
1:B:302:MET:O	1:B:306:LEU:HD12	2.16	0.46
1:B:342:ASN:CG	1:B:344:ILE:HD11	2.36	0.46
1:A:450:LEU:HA	1:A:461:GLU:HG3	1.98	0.45
1:B:248:LYS:HD2	1:B:248:LYS:HA	1.86	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:34:ALA:HB2	1:B:173:LEU:HD23	1.98	0.45
1:A:320:GLY:HA3	1:A:419:TYR:CD2	2.52	0.45
1:A:76:PHE:O	1:A:80:GLU:HG3	2.17	0.45
1:B:320:GLY:HA3	1:B:419:TYR:CE2	2.51	0.45
1:A:270:GLN:HG3	1:A:274:LEU:CD2	2.46	0.45
1:A:499:LYS:HG3	1:A:512:THR:HG22	1.98	0.45
1:A:381:ARG:NE	1:A:383:GLU:HG2	2.32	0.45
1:A:320:GLY:HA3	1:A:419:TYR:CE2	2.52	0.45
1:B:444:PHE:CZ	1:B:460:GLU:HG2	2.51	0.45
6:B:611:TFX:H17B	6:B:612:TFX:C3	2.46	0.45
1:B:316:GLN:CD	1:B:316:GLN:H	2.20	0.45
1:B:472:ALA:O	1:B:476:LYS:HG3	2.17	0.45
6:B:611:TFX:H17B	6:B:612:TFX:H3	1.99	0.45
1:A:322:ASN:ND2	1:A:421:PHE:HB3	2.32	0.44
1:A:371:PHE:O	1:B:520:ARG:NH2	2.51	0.44
1:A:220:ALA:HB3	1:A:317:ILE:HG22	1.98	0.44
1:A:329:PHE:HD1	6:A:613:TFX:C17	2.30	0.44
1:A:481:ASN:N	1:A:481:ASN:OD1	2.51	0.44
1:B:234:THR:HG23	1:B:293:THR:HG22	2.00	0.44
1:B:236:LEU:HA	1:B:236:LEU:HD12	1.79	0.44
1:B:99:ILE:CD1	1:B:185:PHE:HB3	2.48	0.44
1:B:329:PHE:HD1	6:B:612:TFX:H17A	1.83	0.44
1:A:348:LYS:O	1:A:352:GLU:HG2	2.18	0.43
1:A:502:THR:O	1:A:503:LEU:HD23	2.18	0.43
1:B:220:ALA:HB3	1:B:317:ILE:HG22	2.01	0.43
1:A:122:THR:HG23	1:A:124:SER:H	1.82	0.43
1:A:329:PHE:HD1	6:A:613:TFX:H17A	1.83	0.43
1:B:428:LEU:HA	1:B:429:PRO:HD3	1.89	0.43
1:A:395:ASP:CG	1:A:515:ARG:HH21	2.18	0.43
1:A:34:ALA:HB3	1:A:93:LEU:HD12	2.01	0.43
1:B:21:PHE:CE1	1:B:451:GLU:HB2	2.54	0.43
1:B:479:ASN:O	1:B:481:ASN:N	2.46	0.43
1:A:42:ARG:NE	1:A:90:GLU:OE2	2.51	0.43
1:B:208:LEU:HD23	1:B:217:PHE:HZ	1.83	0.43
1:A:156:LEU:HD12	1:A:261:ILE:HD11	2.01	0.43
1:B:193:THR:HG23	1:B:219:ARG:HB2	2.01	0.43
1:B:499:LYS:HG2	1:B:512:THR:HG22	2.01	0.43
1:A:38:LEU:HA	1:A:38:LEU:HD23	1.76	0.42
1:A:427:LYS:NZ	1:A:454:ASP:OD1	2.52	0.42
1:B:36:PRO:HB2	1:B:38:LEU:HD13	2.01	0.42
1:A:28:PHE:HB3	1:A:31:ILE:HD11	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:224:SER:HB3	1:B:325:GLU:OE2	2.19	0.42
1:A:253:SER:C	1:A:254:ARG:HD3	2.39	0.42
1:A:513:LYS:HB3	1:A:516:ALA:HB2	2.01	0.42
1:B:156:LEU:HD13	1:B:257:GLU:HB3	2.00	0.42
1:B:261:ILE:HA	1:B:261:ILE:HD13	1.90	0.42
1:B:347:ARG:HB2	1:B:385:TYR:CZ	2.53	0.42
1:A:153:PHE:HD1	1:A:165:ASN:HA	1.85	0.42
1:B:526:PHE:O	1:B:529:VAL:HG22	2.19	0.42
1:A:165:ASN:ND2	1:A:292:PRO:HA	2.34	0.42
1:B:500:TYR:CZ	1:B:511:MET:HB2	2.55	0.42
6:B:612:TFX:H15B	8:B:724:HOH:O	2.18	0.42
1:A:118:PHE:O	1:A:119:GLN:NE2	2.53	0.42
1:A:3:ASP:CG	1:A:5:ILE:HD11	2.40	0.42
1:A:254:ARG:HH11	1:A:254:ARG:HG3	1.84	0.41
1:B:416:ALA:HB3	1:B:493:PHE:HB3	2.02	0.41
1:A:500:TYR:CE2	1:A:511:MET:HG3	2.55	0.41
1:B:378:ASP:OD1	1:B:380:GLN:HB3	2.19	0.41
1:A:396:TYR:HB2	1:A:522:TRP:CE2	2.55	0.41
1:A:268:ASP:OD1	1:A:268:ASP:N	2.53	0.41
1:B:175:LEU:O	1:B:178:VAL:HG12	2.20	0.41
1:B:188:ASN:HD21	5:B:602:NAG:C6	2.34	0.41
5:B:606:NAG:H2	5:B:606:NAG:H83	1.91	0.41
1:A:219:ARG:HA	1:A:315:THR:CG2	2.50	0.41
1:A:24:THR:OG1	1:A:101:ALA:HB3	2.20	0.41
1:A:310:GLY:HA2	1:A:312:PHE:CE2	2.56	0.41
1:B:242:ARG:HD3	1:B:289:ASN:O	2.20	0.41
1:B:447:GLY:O	1:B:450:LEU:HB2	2.20	0.41
1:B:417:PHE:N	1:B:417:PHE:CD2	2.89	0.41
1:B:299:LEU:CD2	1:B:303:PRO:HG3	2.51	0.41
1:A:41:LEU:O	1:A:44:LYS:HB2	2.21	0.40
1:B:135:ARG:HH11	1:B:138:ARG:HH21	1.67	0.40
1:B:157:PRO:CB	1:B:236:LEU:HD11	2.51	0.40
1:B:302:MET:O	1:B:305:ILE:HB	2.22	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:157:PRO:CB	1:B:453:ARG:NH2[3_544]	2.13	0.07



## 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	528/529 (100%)	498 (94%)	30 (6%)	0	100	100
1	B	524/529 (99%)	492 (94%)	31 (6%)	1 (0%)	47	78
All	All	1052/1058 (99%)	990 (94%)	61 (6%)	1 (0%)	51	81

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	305	ILE

### 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	451/454 (99%)	441 (98%)	10 (2%)	52	83
1	B	440/454 (97%)	429 (98%)	11 (2%)	47	80
All	All	891/908 (98%)	870 (98%)	21 (2%)	49	81

All (21) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	54	ASP
1	A	63	ASN
1	A	73	PHE
1	A	105	LYS
1	A	144	MET

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Mol	Chain	Res	Type
1	A	195	PHE
1	A	248	LYS
1	A	252	CYS
1	A	263	CYS
1	A	452	ARG
1	B	144	MET
1	B	180	LYS
1	B	195	PHE
1	B	301	ASP
1	B	308	GLU
1	B	316	GLN
1	B	355	LYS
1	B	364	PHE
1	B	432	GLU
1	B	454	ASP
1	B	515	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	165	ASN
1	A	275	ASN
1	A	322	ASN
1	B	188	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 ⓘ

14 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	C	1	1,2	14,14,15	0.97	1 (7%)	17,19,21	2.02	1 (5%)
2	NAG	C	2	2	14,14,15	0.78	1 (7%)	17,19,21	1.60	2 (11%)
2	FUC	C	3	2	10,10,11	0.86	0	14,14,16	0.79	0
3	NAG	D	1	1,3	14,14,15	0.52	0	17,19,21	0.60	0
3	NAG	D	2	3	14,14,15	0.70	1 (7%)	17,19,21	0.68	1 (5%)
3	NAG	E	1	1,3	14,14,15	0.72	1 (7%)	17,19,21	0.86	1 (5%)
3	NAG	E	2	3	14,14,15	0.44	0	17,19,21	0.60	0
4	NAG	F	1	1,4	14,14,15	0.93	1 (7%)	17,19,21	0.73	0
4	FUC	F	2	4	10,10,11	1.85	3 (30%)	14,14,16	1.64	2 (14%)
3	NAG	G	1	1,3	14,14,15	1.83	2 (14%)	17,19,21	1.06	1 (5%)
3	NAG	G	2	3	14,14,15	1.58	1 (7%)	17,19,21	1.39	2 (11%)
2	NAG	H	1	1,2	14,14,15	0.34	0	17,19,21	1.39	2 (11%)
2	NAG	H	2	2	14,14,15	0.23	0	17,19,21	0.50	0
2	FUC	H	3	2	10,10,11	0.68	0	14,14,16	1.36	3 (21%)

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	C	1	1,2	-	2/6/23/26	0/1/1/1
2	NAG	C	2	2	-	4/6/23/26	0/1/1/1
2	FUC	C	3	2	1/1/4/5	-	0/1/1/1
3	NAG	D	1	1,3	-	2/6/23/26	0/1/1/1
3	NAG	D	2	3	-	4/6/23/26	0/1/1/1
3	NAG	E	1	1,3	-	2/6/23/26	0/1/1/1
3	NAG	E	2	3	-	0/6/23/26	0/1/1/1
4	NAG	F	1	1,4	-	2/6/23/26	0/1/1/1
4	FUC	F	2	4	1/1/4/5	-	0/1/1/1
3	NAG	G	1	1,3	-	3/6/23/26	0/1/1/1
3	NAG	G	2	3	-	4/6/23/26	0/1/1/1
2	NAG	H	1	1,2	-	4/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	NAG	H	2	2	-	1/6/23/26	0/1/1/1
2	FUC	H	3	2	1/1/4/5	-	0/1/1/1

All (11) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	G	2	NAG	O5-C1	-5.72	1.34	1.43
3	G	1	NAG	O5-C1	-5.34	1.35	1.43
3	G	1	NAG	C1-C2	-3.97	1.46	1.52
4	F	2	FUC	O5-C1	-3.87	1.37	1.43
2	C	1	NAG	O5-C1	3.50	1.49	1.43
4	F	1	NAG	C1-C2	3.15	1.57	1.52
4	F	2	FUC	C2-C3	2.86	1.56	1.52
3	E	1	NAG	O5-C1	-2.60	1.39	1.43
3	D	2	NAG	O5-C1	-2.40	1.39	1.43
4	F	2	FUC	C1-C2	2.40	1.57	1.52
2	C	2	NAG	C1-C2	2.08	1.55	1.52

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	1	NAG	C1-O5-C5	7.83	122.80	112.19
4	F	2	FUC	C1-C2-C3	4.55	115.26	109.67
2	C	2	NAG	C2-N2-C7	4.53	129.35	122.90
2	H	1	NAG	C2-N2-C7	4.36	129.11	122.90
2	C	2	NAG	C1-O5-C5	3.95	117.54	112.19
3	G	2	NAG	C4-C3-C2	3.78	116.55	111.02
3	G	2	NAG	C3-C4-C5	3.61	116.68	110.24
2	H	3	FUC	C1-C2-C3	3.25	113.66	109.67
3	G	1	NAG	C3-C4-C5	2.95	115.50	110.24
4	F	2	FUC	C2-C3-C4	2.70	115.57	110.89
3	E	1	NAG	C1-O5-C5	2.51	115.60	112.19
2	H	3	FUC	O5-C1-C2	2.41	114.50	110.77
2	H	3	FUC	C1-O5-C5	2.25	117.89	112.78
2	H	1	NAG	C1-C2-N2	2.14	114.14	110.49
3	D	2	NAG	C3-C4-C5	2.07	113.92	110.24

All (3) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
2	H	3	FUC	C1
4	F	2	FUC	C1

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Mol	Chain	Res	Type	Atom
2	C	3	FUC	C1

All (28) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	G	2	NAG	C1-C2-N2-C7
3	D	2	NAG	C4-C5-C6-O6
3	D	2	NAG	O5-C5-C6-O6
3	E	1	NAG	O5-C5-C6-O6
2	C	1	NAG	O5-C5-C6-O6
2	C	1	NAG	C4-C5-C6-O6
3	G	2	NAG	O5-C5-C6-O6
2	C	2	NAG	C8-C7-N2-C2
2	C	2	NAG	O7-C7-N2-C2
3	G	1	NAG	C8-C7-N2-C2
3	G	1	NAG	O7-C7-N2-C2
3	D	2	NAG	C8-C7-N2-C2
3	D	2	NAG	O7-C7-N2-C2
2	H	1	NAG	C8-C7-N2-C2
2	H	1	NAG	O7-C7-N2-C2
3	G	2	NAG	C4-C5-C6-O6
2	C	2	NAG	O5-C5-C6-O6
3	D	1	NAG	O5-C5-C6-O6
4	F	1	NAG	O5-C5-C6-O6
3	E	1	NAG	C4-C5-C6-O6
3	G	1	NAG	O5-C5-C6-O6
4	F	1	NAG	C4-C5-C6-O6
2	C	2	NAG	C3-C2-N2-C7
3	G	2	NAG	C3-C2-N2-C7
2	H	1	NAG	C4-C5-C6-O6
2	H	2	NAG	C4-C5-C6-O6
2	H	1	NAG	C3-C2-N2-C7
3	D	1	NAG	C4-C5-C6-O6

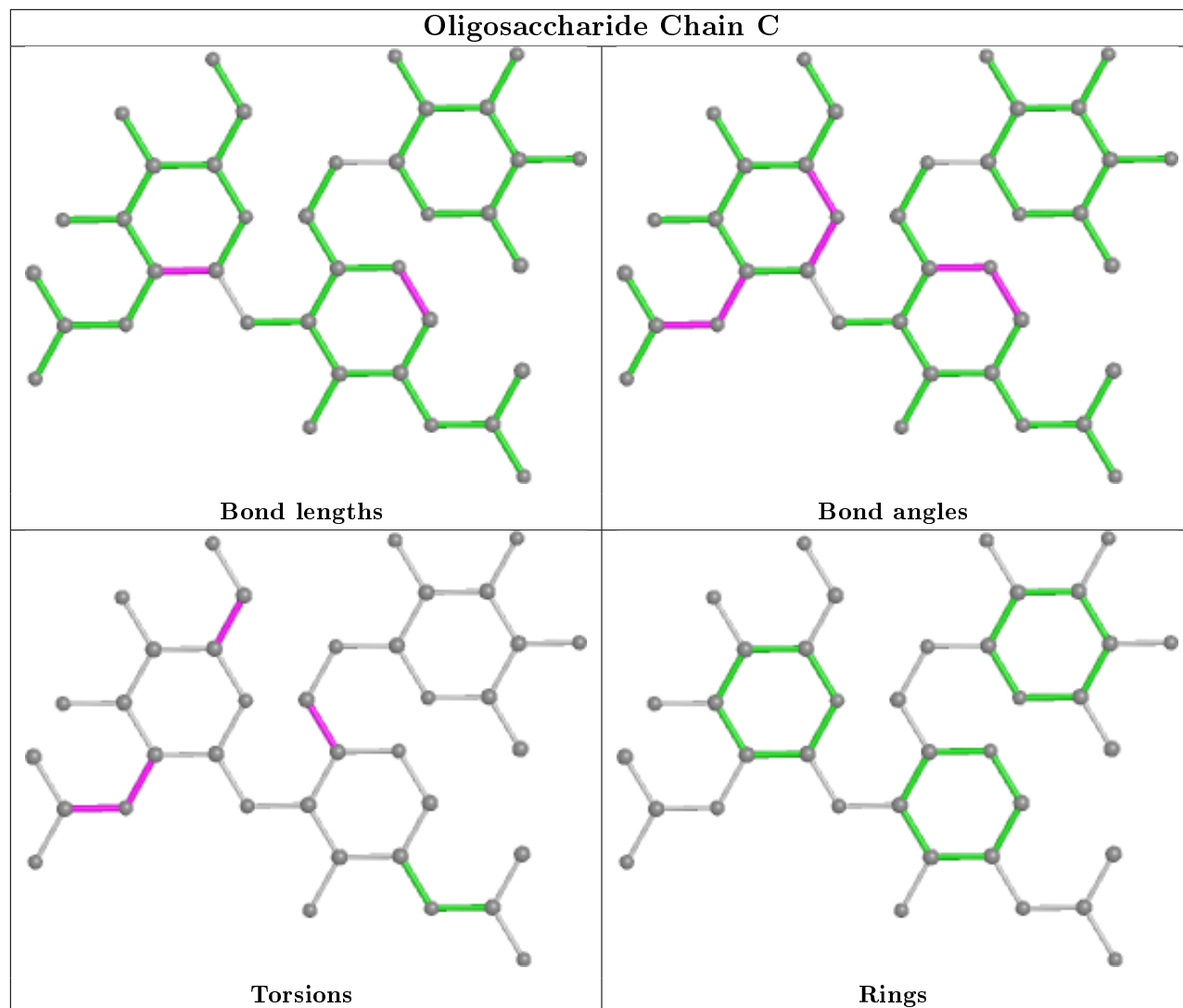
There are no ring outliers.

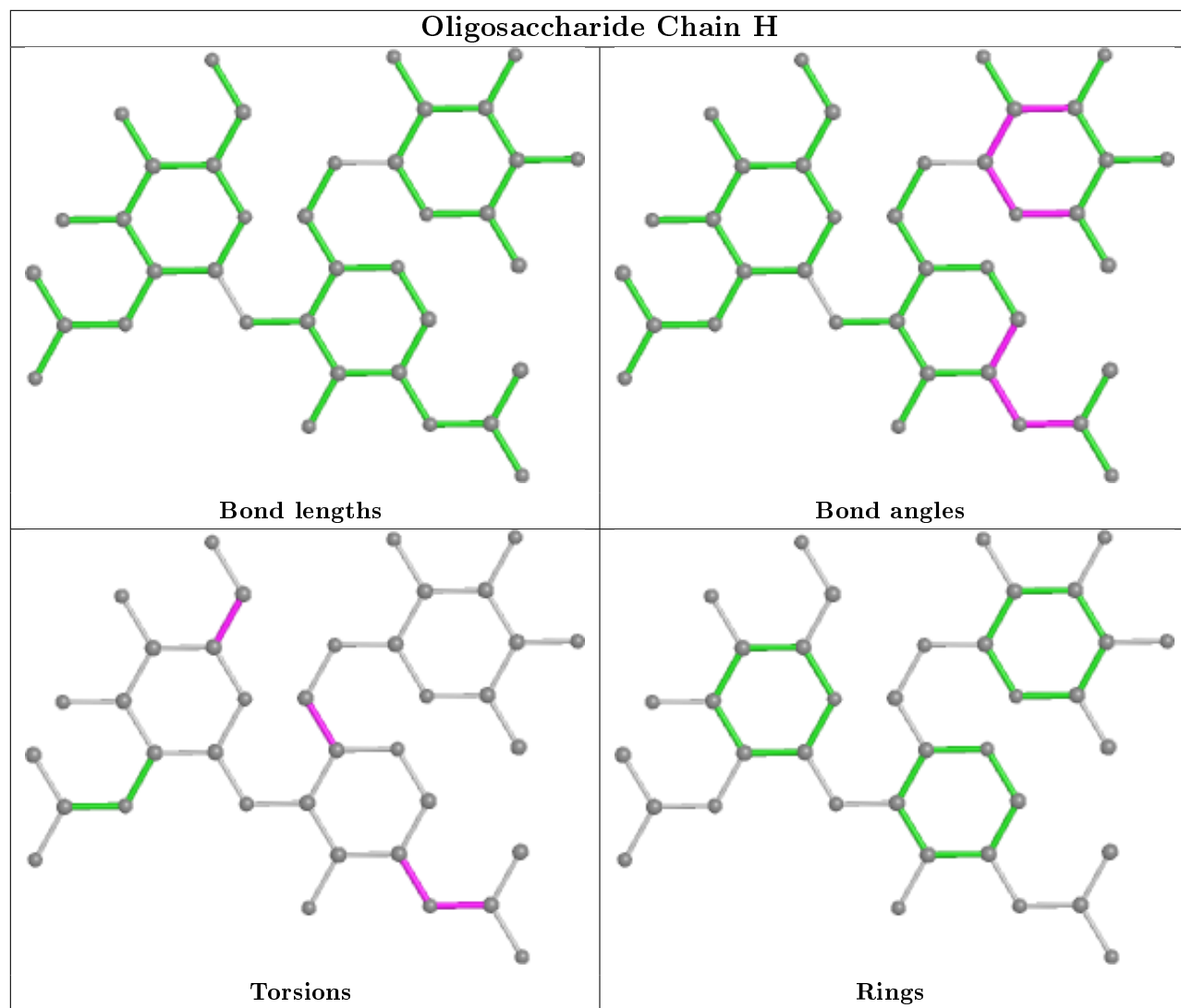
2 monomers are involved in 2 short contacts:

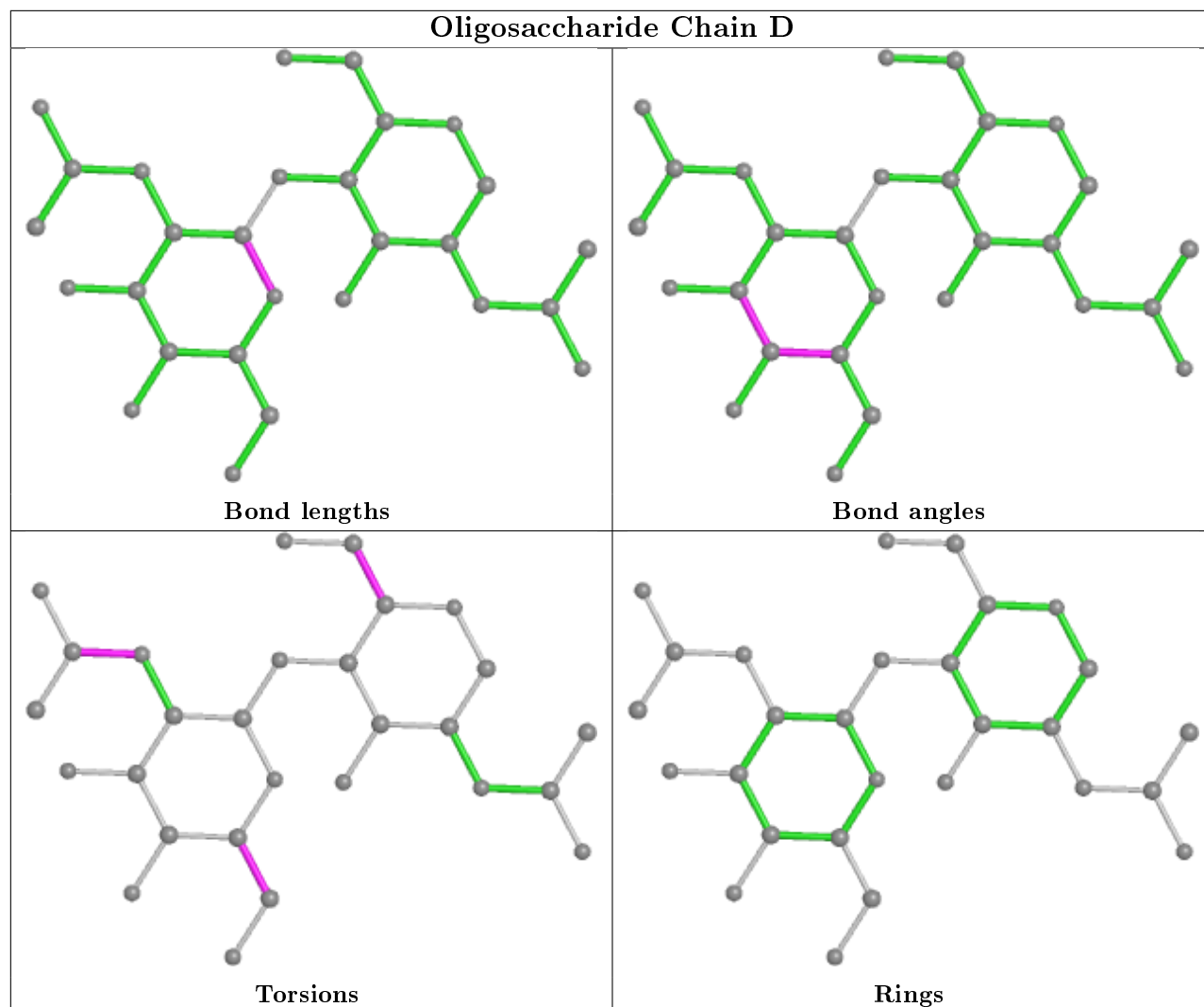
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	2	NAG	1	0
2	H	1	NAG	1	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths,

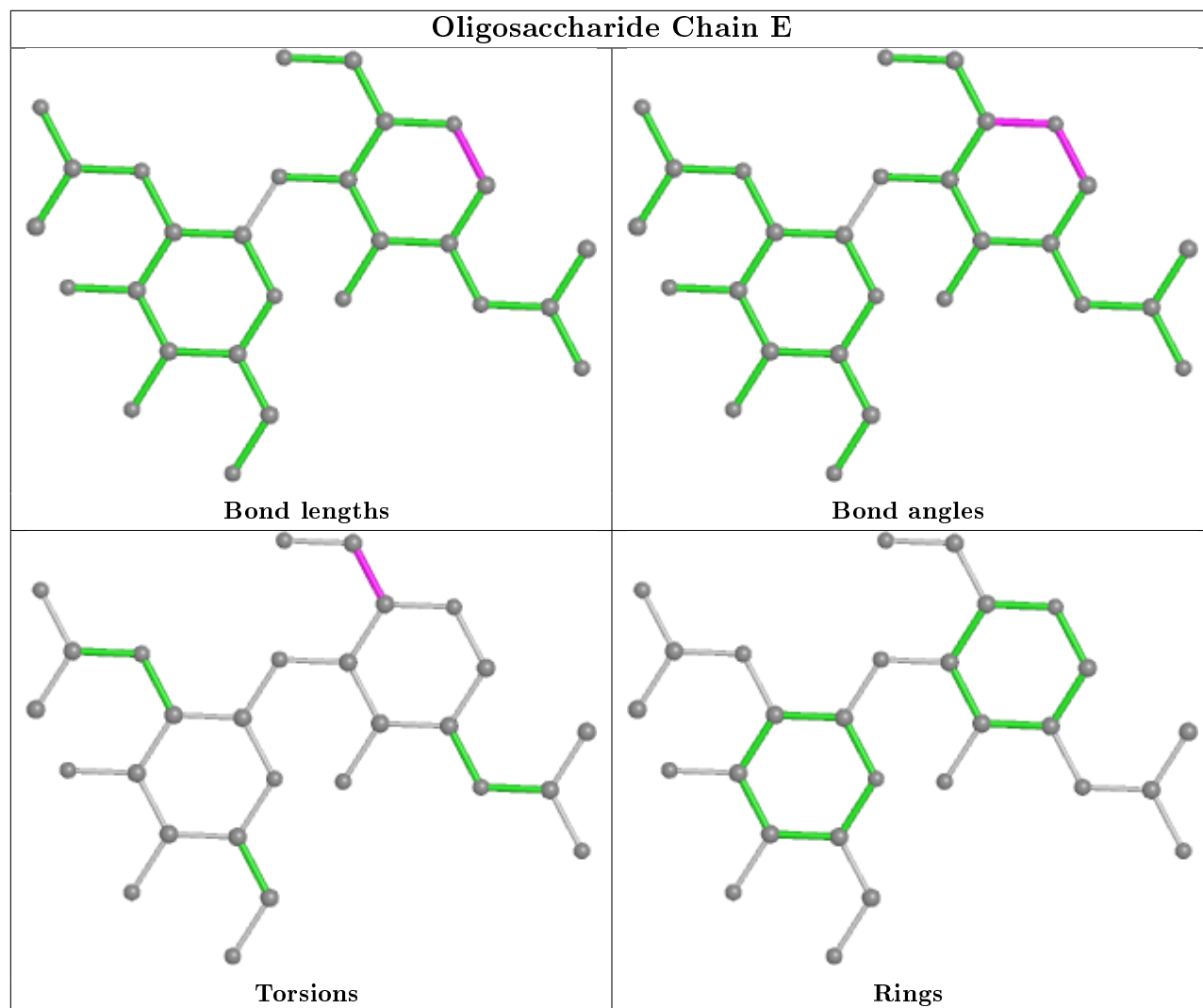
bond angles, torsion angles, and ring geometry for oligosaccharide.

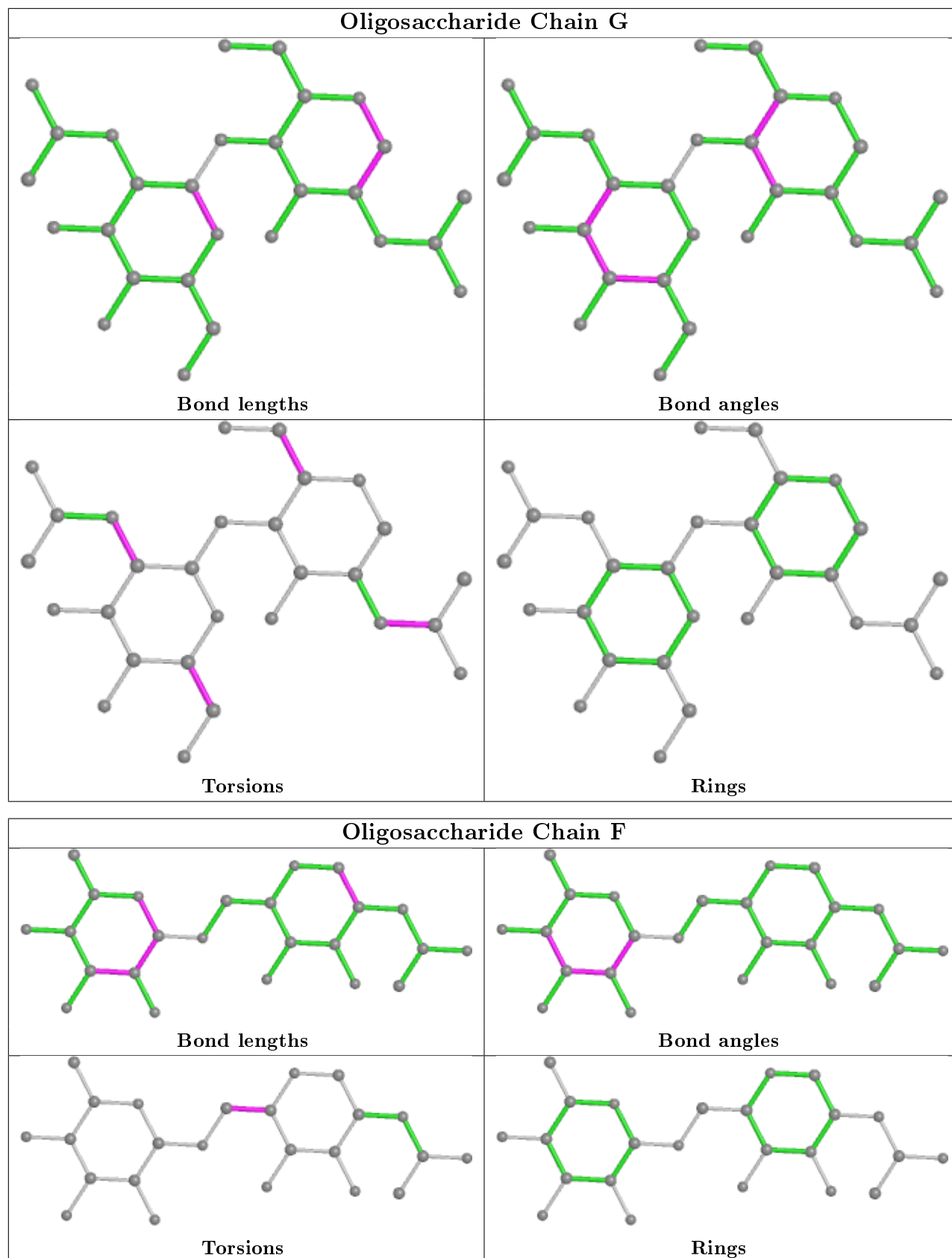












## 5.6 Ligand geometry

Of 13 ligands modelled in this entry, 2 are monoatomic - leaving 11 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
6	TFX	B	612	-	17,22,22	1.45	2 (11%)	23,32,32	2.41	4 (17%)
5	NAG	A	601	1	14,14,15	0.92	1 (7%)	17,19,21	1.36	1 (5%)
5	NAG	B	601	1	14,14,15	1.27	1 (7%)	17,19,21	0.80	0
6	TFX	A	613	-	17,22,22	1.41	2 (11%)	23,32,32	2.75	5 (21%)
6	TFX	B	611	-	17,22,22	1.38	2 (11%)	23,32,32	2.49	4 (17%)
5	NAG	B	605	1	14,14,15	0.92	1 (7%)	17,19,21	0.60	0
5	NAG	B	602	1	14,14,15	2.49	3 (21%)	17,19,21	2.23	4 (23%)
5	NAG	B	606	1	14,14,15	3.02	5 (35%)	17,19,21	2.21	3 (17%)
5	NAG	B	607	1	14,14,15	0.71	1 (7%)	17,19,21	1.42	1 (5%)
6	TFX	A	612	-	17,22,22	1.34	2 (11%)	23,32,32	2.64	4 (17%)
5	NAG	A	607	1	14,14,15	0.34	0	17,19,21	0.41	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
6	TFX	B	612	-	-	1/8/8/8	0/3/3/3
5	NAG	A	601	1	-	3/6/23/26	0/1/1/1
5	NAG	B	601	1	-	3/6/23/26	0/1/1/1
6	TFX	A	613	-	-	2/8/8/8	0/3/3/3
6	TFX	B	611	-	-	6/8/8/8	0/3/3/3
5	NAG	B	605	1	-	4/6/23/26	0/1/1/1
5	NAG	B	602	1	-	5/6/23/26	0/1/1/1
5	NAG	B	606	1	-	4/6/23/26	0/1/1/1
5	NAG	B	607	1	-	4/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
6	TFX	A	612	-	-	2/8/8/8	0/3/3/3
5	NAG	A	607	1	-	1/6/23/26	0/1/1/1

All (20) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	B	606	NAG	O5-C1	9.85	1.59	1.43
5	B	602	NAG	O5-C1	7.65	1.55	1.43
5	B	601	NAG	O5-C1	-4.26	1.36	1.43
5	B	602	NAG	C1-C2	-3.79	1.46	1.52
6	B	612	TFX	C5-C8	-3.61	1.39	1.48
6	B	612	TFX	C2-N2	3.51	1.45	1.37
6	A	612	TFX	C2-N2	3.50	1.45	1.37
6	A	613	TFX	C5-C8	-3.44	1.39	1.48
6	A	613	TFX	C2-N2	3.44	1.45	1.37
6	B	611	TFX	C2-N2	3.41	1.45	1.37
6	B	611	TFX	C5-C8	-3.40	1.39	1.48
5	A	601	NAG	O5-C1	3.30	1.49	1.43
6	A	612	TFX	C5-C8	-3.14	1.40	1.48
5	B	606	NAG	C3-C2	-3.13	1.45	1.52
5	B	605	NAG	O5-C1	-3.10	1.38	1.43
5	B	606	NAG	C1-C2	-2.85	1.48	1.52
5	B	602	NAG	C3-C2	-2.48	1.47	1.52
5	B	607	NAG	O5-C1	-2.43	1.39	1.43
5	B	606	NAG	C2-N2	-2.16	1.42	1.46
5	B	606	NAG	O5-C5	2.04	1.47	1.43

All (26) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	A	613	TFX	C5-C8-N1	11.11	149.79	123.75
6	A	612	TFX	C5-C8-N1	10.91	149.32	123.75
6	B	611	TFX	C5-C8-N1	9.99	147.18	123.75
6	B	612	TFX	C5-C8-N1	9.59	146.24	123.75
5	B	606	NAG	C1-O5-C5	5.91	120.20	112.19
5	B	602	NAG	C2-N2-C7	5.47	130.69	122.90
5	A	601	NAG	C1-O5-C5	5.35	119.44	112.19
5	B	602	NAG	C1-O5-C5	4.79	118.69	112.19
5	B	606	NAG	O6-C6-C5	4.65	127.23	111.29
6	B	611	TFX	C11-C10-S1	4.03	133.16	125.10
6	A	613	TFX	C11-C10-S1	4.02	133.15	125.10
5	B	607	NAG	C2-N2-C7	4.02	128.63	122.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	B	612	TFX	C11-C10-S1	3.97	133.05	125.10
5	B	606	NAG	C4-C3-C2	-3.80	105.44	111.02
6	A	612	TFX	C11-C10-S1	3.79	132.69	125.10
5	B	602	NAG	C3-C4-C5	3.37	116.25	110.24
5	B	602	NAG	O5-C5-C6	-3.02	102.47	107.20
6	A	613	TFX	C7-C2-N2	-2.79	117.85	121.63
6	A	612	TFX	C9-C10-S1	-2.61	108.39	111.85
6	B	612	TFX	C3-C2-N2	-2.59	118.13	121.63
6	B	611	TFX	C9-C10-S1	-2.56	108.45	111.85
6	A	613	TFX	C9-C10-S1	-2.55	108.47	111.85
6	B	612	TFX	C9-C10-S1	-2.51	108.53	111.85
6	A	613	TFX	C13-C12-C11	2.11	121.00	118.40
6	B	611	TFX	C13-C12-C11	2.10	120.99	118.40
6	A	612	TFX	C17-N1-C8	2.06	127.87	125.21

There are no chirality outliers.

All (35) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
6	B	611	TFX	C4-C5-C8-N1
6	A	612	TFX	C4-C5-C8-N1
6	A	612	TFX	C6-C5-C8-N1
5	B	607	NAG	O5-C5-C6-O6
5	B	606	NAG	O5-C5-C6-O6
5	B	605	NAG	O5-C5-C6-O6
5	B	607	NAG	C4-C5-C6-O6
5	B	606	NAG	C4-C5-C6-O6
5	B	602	NAG	O5-C5-C6-O6
5	B	602	NAG	C4-C5-C6-O6
5	A	601	NAG	C8-C7-N2-C2
5	A	601	NAG	O7-C7-N2-C2
5	B	605	NAG	C8-C7-N2-C2
5	B	605	NAG	O7-C7-N2-C2
5	B	602	NAG	C8-C7-N2-C2
5	B	602	NAG	O7-C7-N2-C2
5	B	606	NAG	C8-C7-N2-C2
5	B	606	NAG	O7-C7-N2-C2
5	B	607	NAG	C8-C7-N2-C2
5	B	607	NAG	O7-C7-N2-C2
5	B	605	NAG	C4-C5-C6-O6
6	B	611	TFX	C7-C2-N2-C15
6	B	611	TFX	C6-C5-C8-N1

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Mol	Chain	Res	Type	Atoms
6	B	611	TFX	C3-C2-N2-C16
5	B	601	NAG	C1-C2-N2-C7
6	B	611	TFX	C3-C2-N2-C15
6	B	611	TFX	C7-C2-N2-C16
5	B	601	NAG	O5-C5-C6-O6
5	A	601	NAG	O5-C5-C6-O6
6	B	612	TFX	C6-C5-C8-S1
6	A	613	TFX	C4-C5-C8-S1
6	A	613	TFX	C6-C5-C8-S1
5	B	601	NAG	C3-C2-N2-C7
5	B	602	NAG	C3-C2-N2-C7
5	A	607	NAG	C4-C5-C6-O6

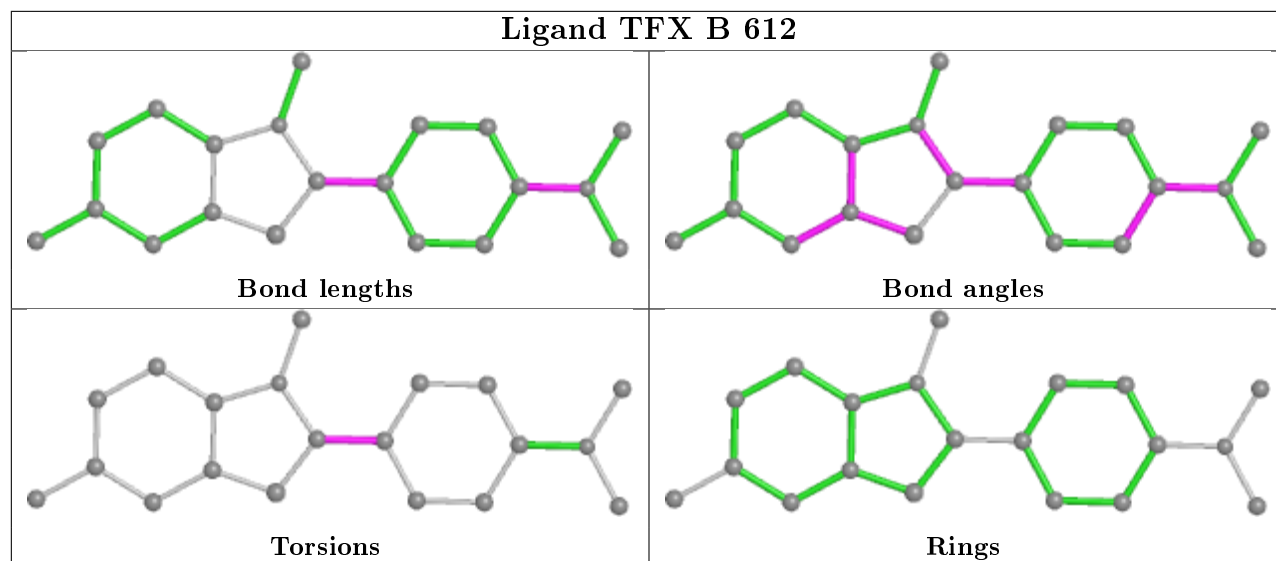
There are no ring outliers.

8 monomers are involved in 20 short contacts:

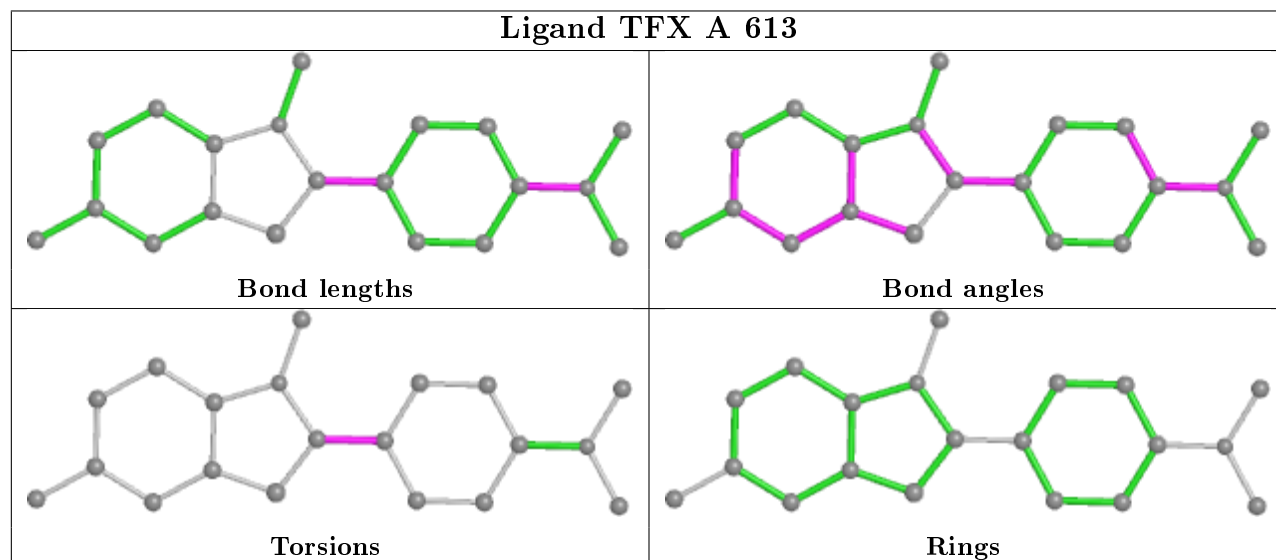
Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	B	612	TFX	8	0
5	A	601	NAG	1	0
6	A	613	TFX	5	0
6	B	611	TFX	7	0
5	B	602	NAG	1	0
5	B	606	NAG	2	0
5	B	607	NAG	1	0
6	A	612	TFX	4	0

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.

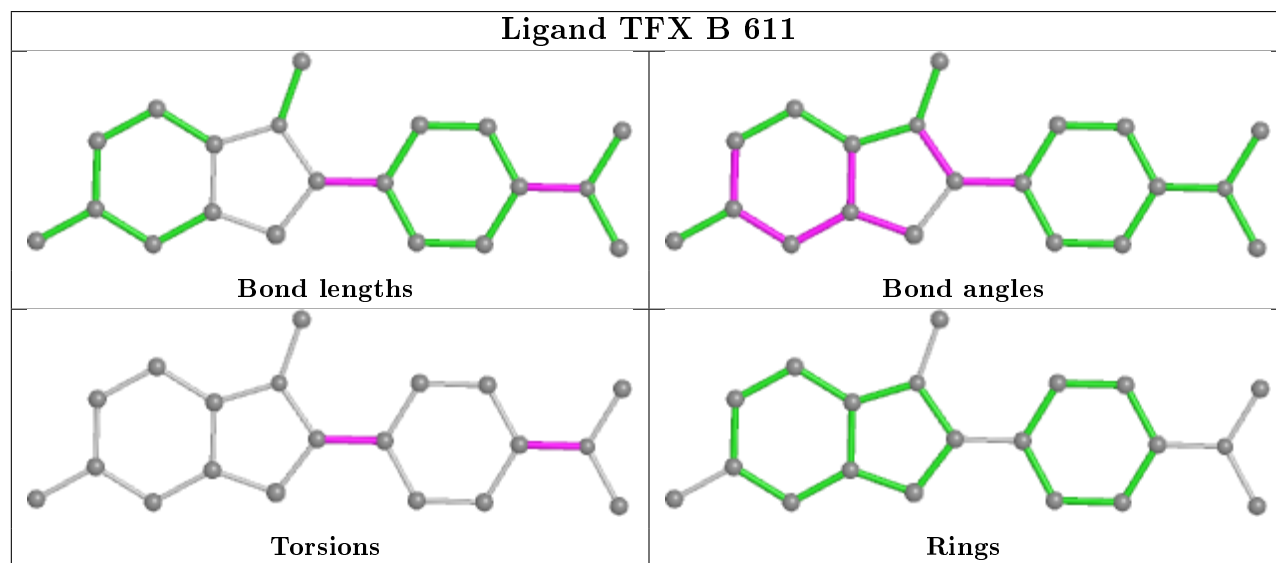
## Ligand TFX B 612

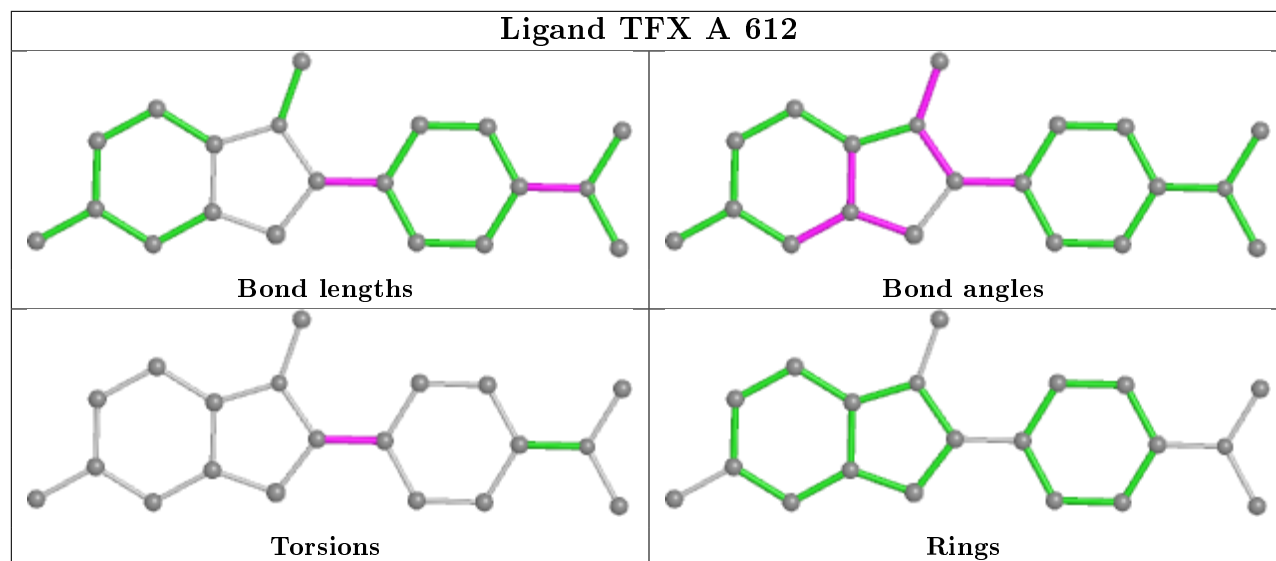


## Ligand TFX A 613



## Ligand TFX B 611





## 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	527/529 (99%)	0.18	19 (3%)	42 32	44, 78, 116, 167	2 (0%)
1	B	526/529 (99%)	0.17	17 (3%)	47 37	42, 78, 124, 192	4 (0%)
All	All	1053/1058 (99%)	0.18	36 (3%)	45 35	42, 78, 121, 192	6 (0%)

All (36) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	4	ILE	5.3
1	A	54	ASP	4.5
1	A	282	TYR	4.0
1	A	55	ILE	3.7
1	B	40	ARG	3.6
1	B	53	SER	3.6
1	B	13	VAL	3.1
1	A	246	LEU	3.1
1	A	283	GLY	3.1
1	B	10	ASN	3.0
1	B	371	PHE	2.9
1	A	19	THR	2.9
1	A	3	ASP	2.9
1	A	280	VAL	2.9
1	A	345	ILE	2.9
1	A	105	LYS	2.7
1	B	384	ASN	2.7
1	B	5	ILE	2.6
1	A	275	ASN	2.6
1	A	7	ALA	2.5
1	A	281	PRO	2.4
1	B	28	PHE	2.4
1	A	23	GLY	2.3
1	A	25	VAL	2.3

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Mol	Chain	Res	Type	RSRZ
1	A	364	PHE	2.3
1	B	52	TRP	2.3
1	B	6	ILE	2.3
1	B	186	GLY	2.3
1	B	385	TYR	2.2
1	A	12	LYS	2.2
1	B	12	LYS	2.1
1	A	529	VAL	2.1
1	B	49	LEU	2.1
1	B	7	ALA	2.1
1	B	8	THR	2.1
1	B	9	LYS	2.0

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

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

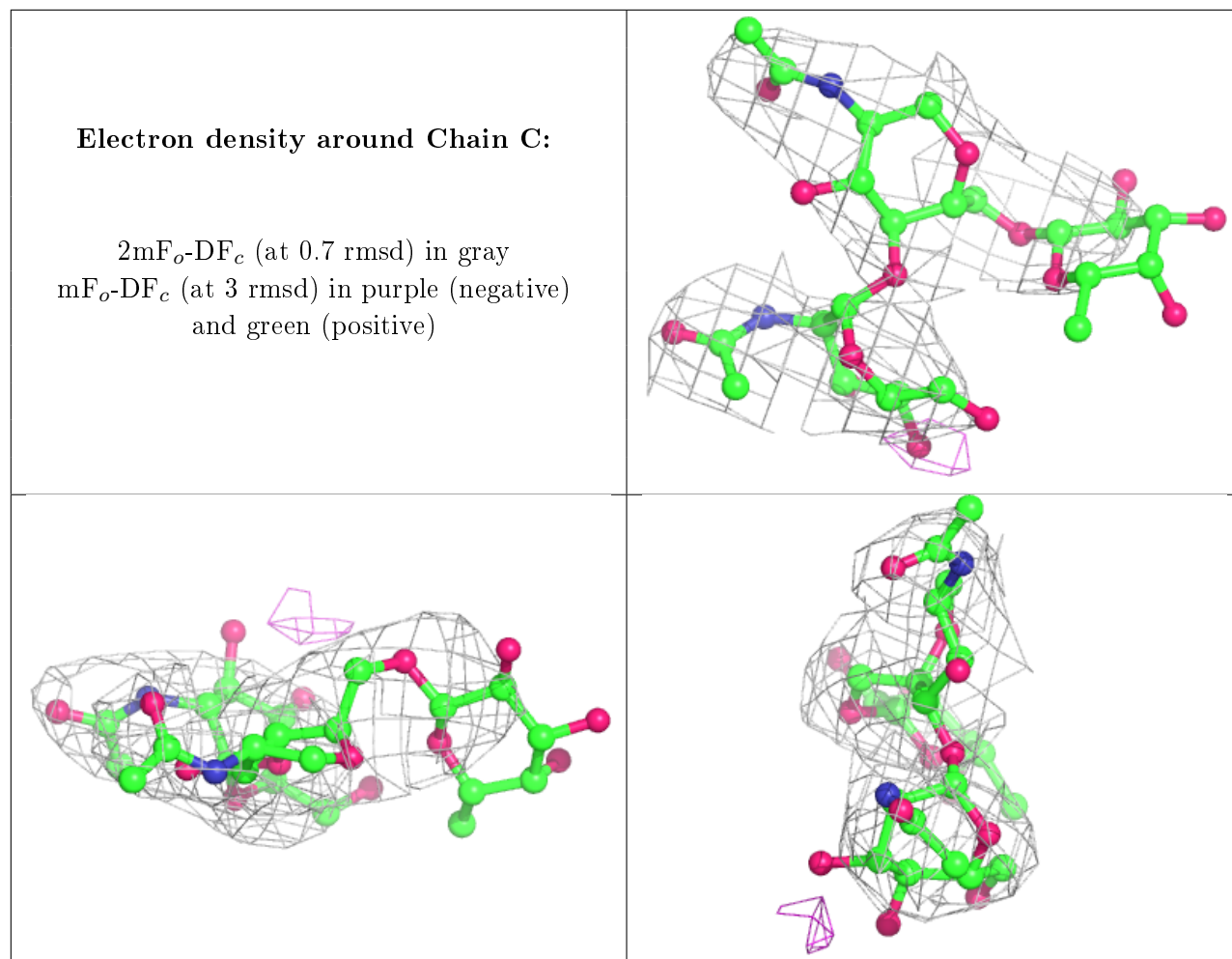
## 6.3 Carbohydrates ⓘ

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
4	NAG	F	1	14/15	0.65	0.21	126,151,168,176	0
4	FUC	F	2	10/11	0.72	0.33	182,192,196,198	0
3	NAG	E	2	14/15	0.74	0.26	114,147,158,163	0
2	FUC	C	3	10/11	0.76	0.79	180,199,205,209	0
3	NAG	G	2	14/15	0.77	0.38	189,200,206,207	0
2	NAG	C	2	14/15	0.77	0.41	153,166,179,182	0
3	NAG	D	2	14/15	0.77	0.40	144,154,162,166	0
3	NAG	G	1	14/15	0.83	0.20	128,143,151,170	0
2	NAG	C	1	14/15	0.86	0.23	112,121,149,158	0
2	NAG	H	1	14/15	0.86	0.18	119,126,155,156	0
3	NAG	D	1	14/15	0.91	0.32	67,91,108,122	0
2	NAG	H	2	14/15	0.91	0.19	109,120,130,134	0
3	NAG	E	1	14/15	0.91	0.23	118,129,143,148	0
2	FUC	H	3	10/11	0.94	0.31	115,122,137,155	0

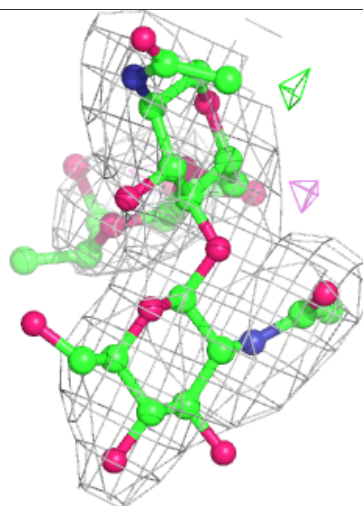
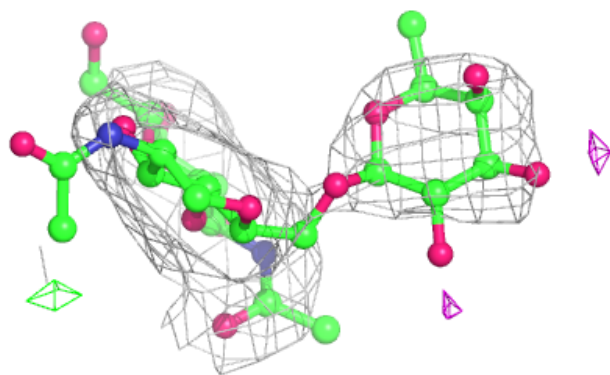
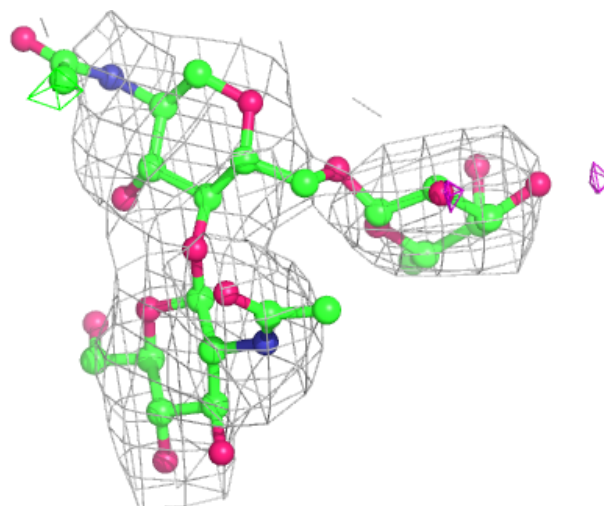
The following is a graphical depiction of the model fit to experimental electron density for oligosac-

charide. Each fit is shown from different orientation to approximate a three-dimensional view.



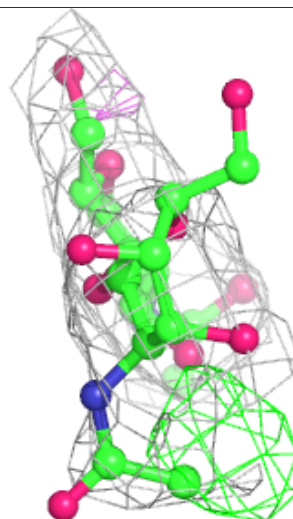
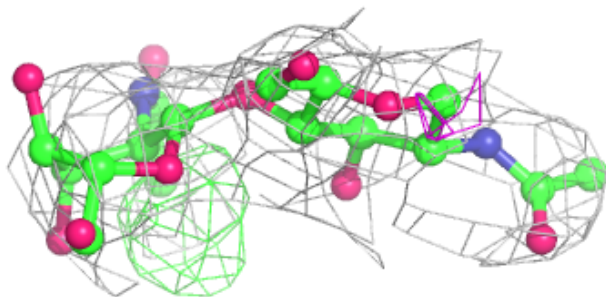
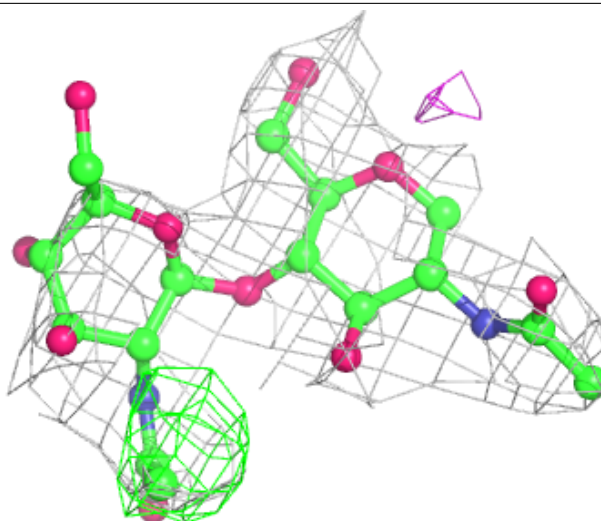
**Electron density around Chain H:**

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



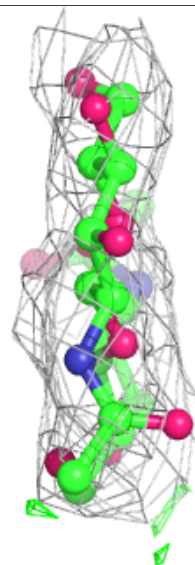
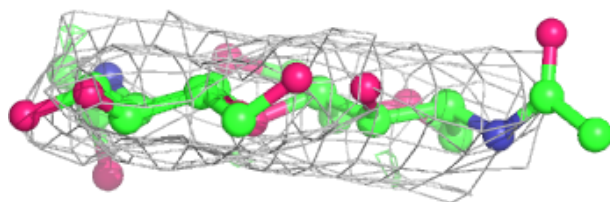
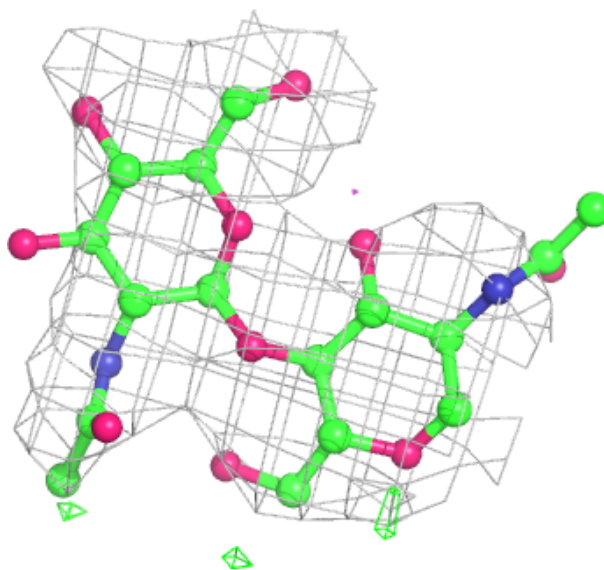
**Electron density around Chain D:**

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



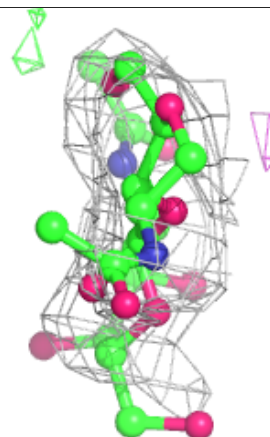
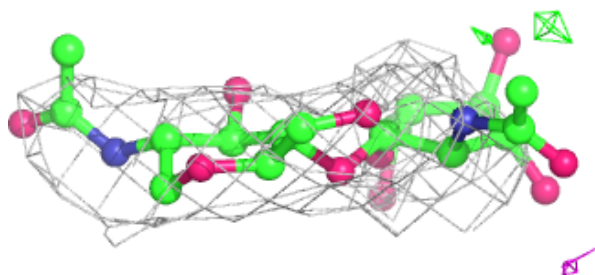
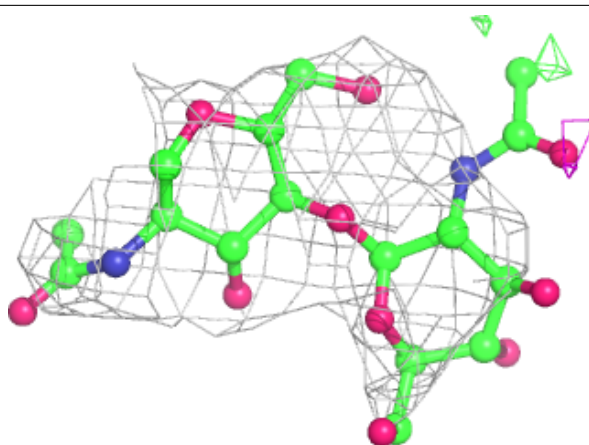
**Electron density around Chain E:**

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

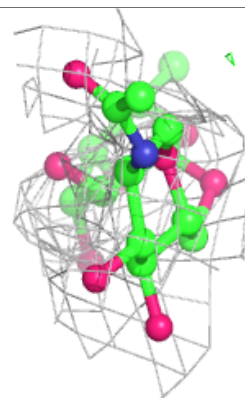
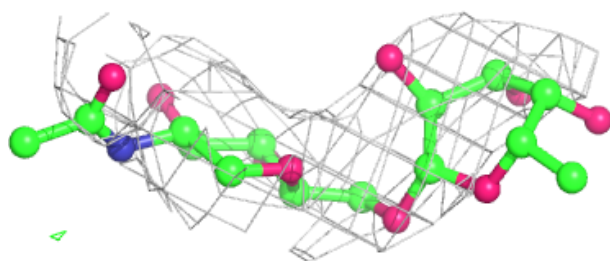
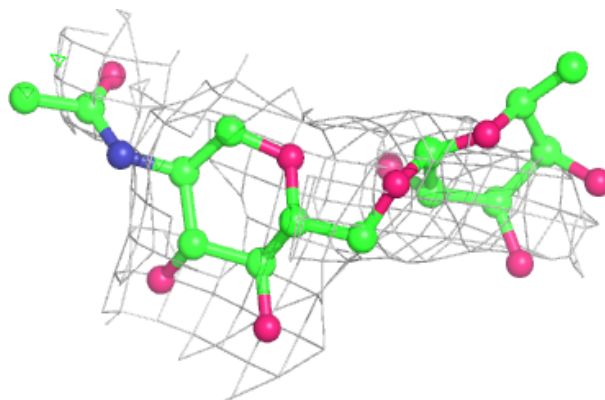


**Electron density around Chain G:**

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

**Electron density around Chain F:**

$2mF_o-DF_c$  (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.

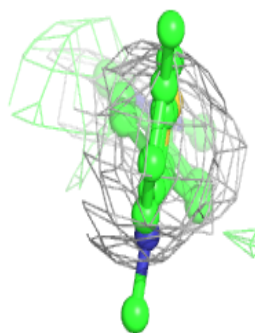
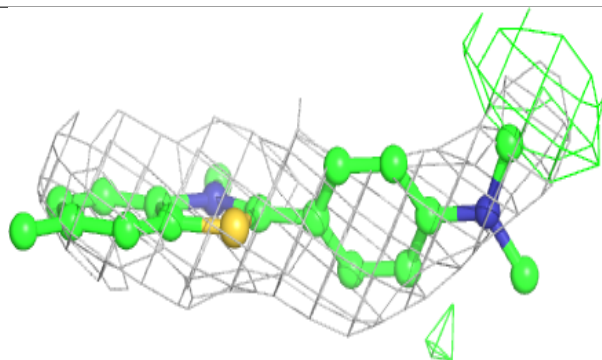
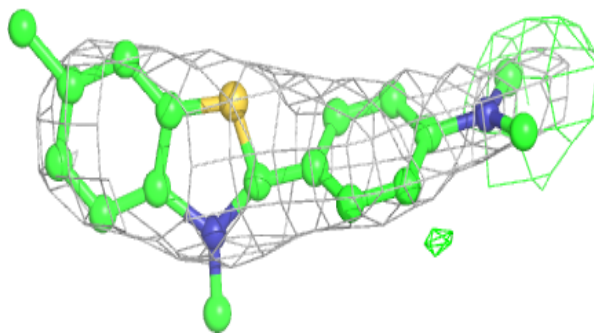
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å <sup>2</sup> )	Q<0.9
7	CL	A	614	1/1	0.25	1.35	222,222,222,222	0
7	CL	B	613	1/1	0.60	0.36	195,195,195,195	0
5	NAG	B	606	14/15	0.68	0.25	115,128,139,141	0
5	NAG	B	601	14/15	0.76	0.24	127,142,152,153	0
5	NAG	B	602	14/15	0.78	0.27	130,137,147,153	0
5	NAG	A	607	14/15	0.80	0.22	112,125,134,134	0
5	NAG	A	601	14/15	0.80	0.37	126,142,158,162	0
6	TFX	A	613	20/20	0.83	0.31	120,137,149,152	0
5	NAG	B	607	14/15	0.83	0.21	158,168,173,174	0
5	NAG	B	605	14/15	0.85	0.14	97,102,114,117	0
6	TFX	B	612	20/20	0.92	0.28	86,104,118,120	0
6	TFX	B	611	20/20	0.94	0.27	56,96,111,120	0
6	TFX	A	612	20/20	0.94	0.25	76,142,164,168	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.

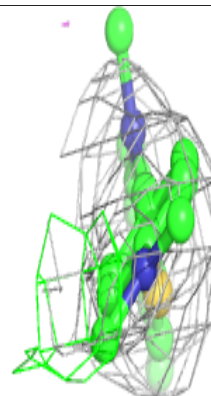
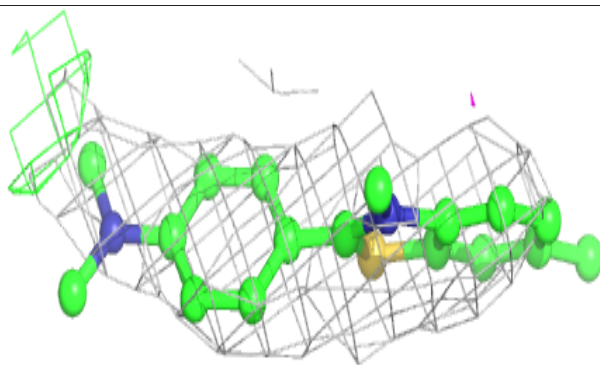
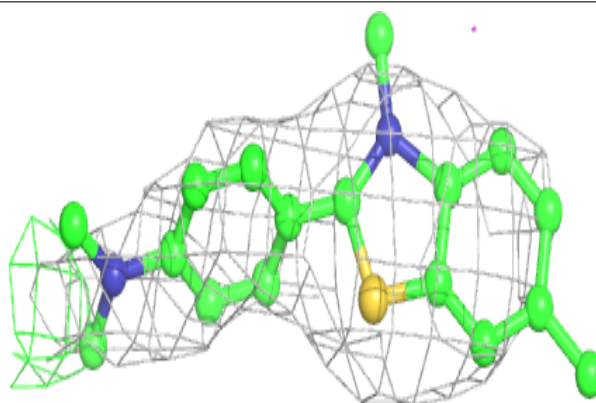


**Electron density around TFX A 613:**

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

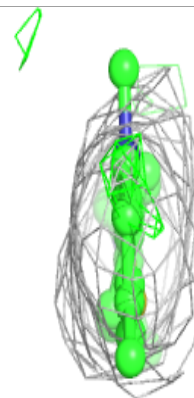
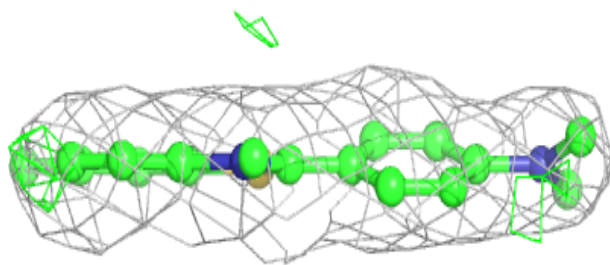
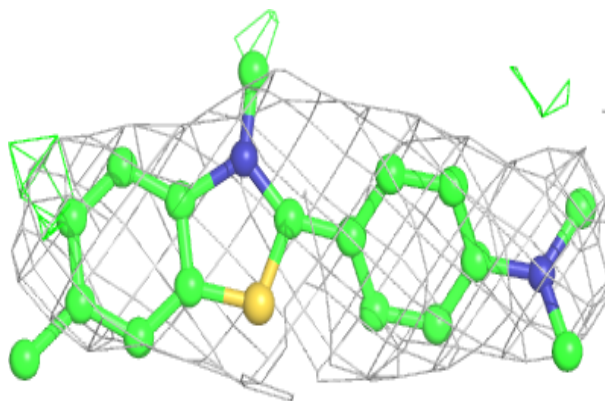
**Electron density around TFX B 612:**

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

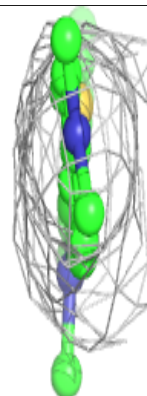
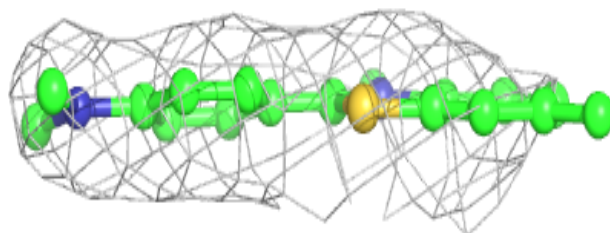
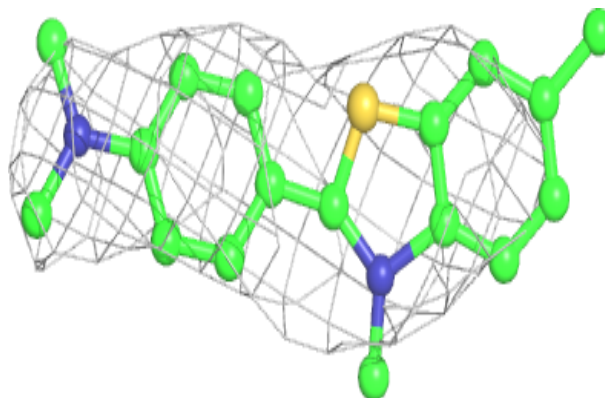


**Electron density around TFX B 611:**

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

**Electron density around TFX A 612:**

$2mF_o-DF_c$  (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.