



Full wwPDB X-ray Structure Validation Report ⓘ

Oct 2, 2021 – 11:59 AM EDT

PDB ID : 3HS6
Title : X-ray crystal structure of eicosapentaenoic acid bound to the cyclooxygenase channel of cyclooxygenase-2
Authors : Vecchio, A.J.; Simmons, D.M.; Malkowski, M.G.
Deposited on : 2009-06-10
Resolution : 2.40 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

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

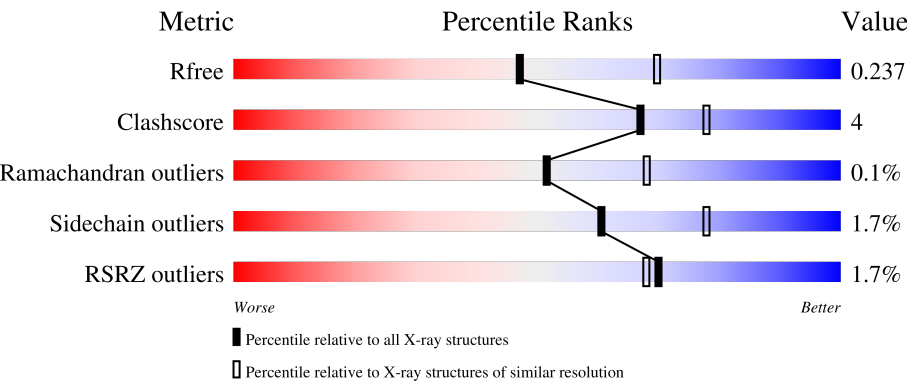
MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.23.2
buster-report : 1.1.7 (2018)
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.23.2

1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.40 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	3907 (2.40-2.40)
Clashscore	141614	4398 (2.40-2.40)
Ramachandran outliers	138981	4318 (2.40-2.40)
Sidechain outliers	138945	4319 (2.40-2.40)
RSRZ outliers	127900	3811 (2.40-2.40)

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 of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	591	
1	B	591	
2	C	2	
2	F	2	
3	D	3	

Continued on next page...

Continued from previous page...

Mol	Chain	Length	Quality of chain
4	E	2	

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	1	X	-	-	-
2	NAG	C	2	-	-	-	X
4	NDG	E	2	-	-	-	X
5	EPA	B	1	-	-	X	-
6	AKR	A	2	-	-	X	-
6	AKR	B	2	-	-	X	-

2 Entry composition [i](#)

There are 11 unique types of molecules in this entry. The entry contains 9765 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 Prostaglandin G/H synthase 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	551	Total	C	N	O	S	0	3	0
			4463	2882	743	813	25			
1	B	551	Total	C	N	O	S	0	3	0
			4442	2868	742	807	25			

There are 14 discrepancies between the modelled and reference sequences:

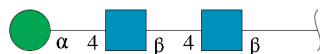
Chain	Residue	Modelled	Actual	Comment	Reference
A	29	HIS	-	expression tag	UNP Q05769
A	30	HIS	-	expression tag	UNP Q05769
A	31	HIS	-	expression tag	UNP Q05769
A	32	HIS	-	expression tag	UNP Q05769
A	33	HIS	-	expression tag	UNP Q05769
A	34	HIS	-	expression tag	UNP Q05769
A	594	ALA	ASN	engineered mutation	UNP Q05769
B	29	HIS	-	expression tag	UNP Q05769
B	30	HIS	-	expression tag	UNP Q05769
B	31	HIS	-	expression tag	UNP Q05769
B	32	HIS	-	expression tag	UNP Q05769
B	33	HIS	-	expression tag	UNP Q05769
B	34	HIS	-	expression tag	UNP Q05769
B	594	ALA	ASN	engineered mutation	UNP Q05769

- 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	C	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			

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



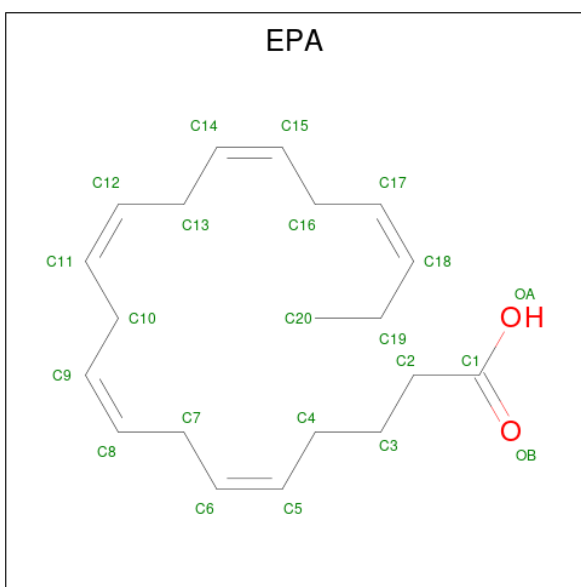
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	D	3	Total	C	N	O	0	0	0
			39	22	2	15			

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



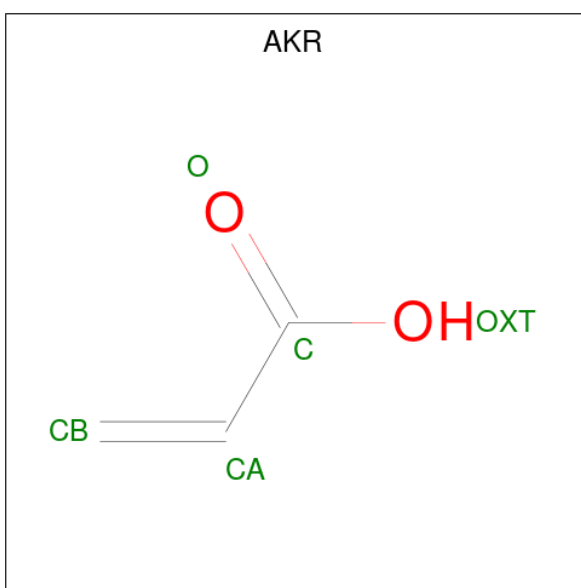
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
4	E	2	Total	C	N	O	0	0	0
			28	16	2	10			

- Molecule 5 is 5,8,11,14,17-EICOSAPENTAENOIC ACID (three-letter code: EPA) (formula: C₂₀H₃₀O₂).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			22	20	2		
5	B	1	Total	C	O	0	0
			22	20	2		

- Molecule 6 is ACRYLIC ACID (three-letter code: AKR) (formula: $C_3H_4O_2$).



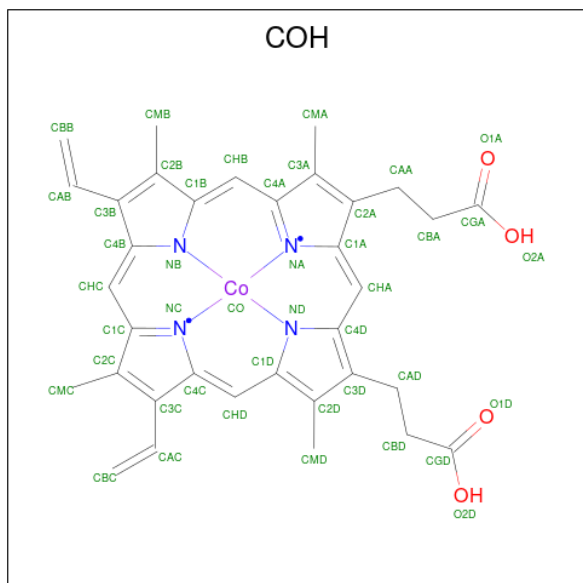
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	A	1	Total	C	O	0	0
			5	3	2		
6	A	1	Total	C	O	0	0
			5	3	2		

Continued on next page...

Continued from previous page...

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	B	1	Total	C	O	0	0
			5	3	2		
6	B	1	Total	C	O	0	0
			5	3	2		

- Molecule 7 is PROTOPORPHYRIN IX CONTAINING CO (three-letter code: COH) (formula: $C_{34}H_{32}CoN_4O_4$).



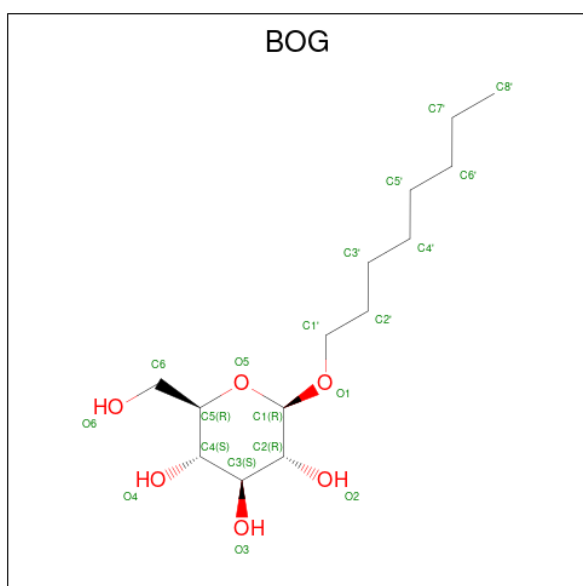
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
7	A	1	Total	C	Co	N	O	0	0
			43	34	1	4	4		
7	B	1	Total	C	Co	N	O	0	0
			43	34	1	4	4		

- Molecule 8 is 2-acetamido-2-deoxy-beta-D-glucopyranose (three-letter code: NAG) (formula: $C_8H_{15}NO_6$).



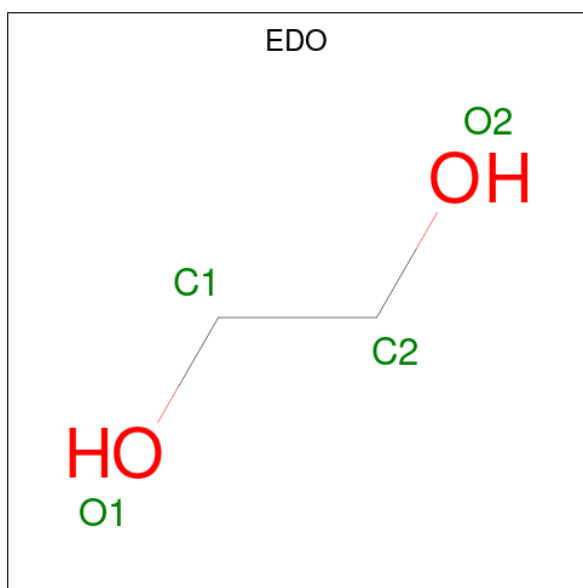
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
8	A	1	Total	C	N	O	0	0
			14	8	1	5		
8	B	1	Total	C	N	O	0	0
			14	8	1	5		

- Molecule 9 is octyl beta-D-glucopyranoside (three-letter code: BOG) (formula: $C_{14}H_{28}O_6$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
9	A	1	Total	C	O	0	0
			20	14	6		
9	B	1	Total	C	O	0	0
			20	14	6		

- Molecule 10 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: $C_2H_6O_2$).

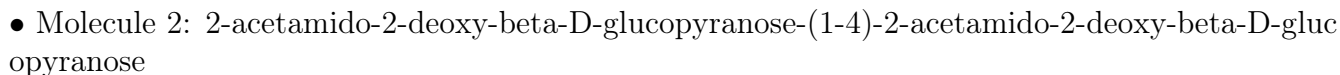
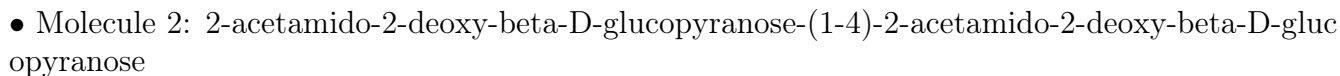


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
10	A	1	Total	C	O	0	0
			4	2	2		
10	A	1	Total	C	O	0	0
			4	2	2		
10	A	1	Total	C	O	0	0
			4	2	2		
10	A	1	Total	C	O	0	0
			4	2	2		
10	A	1	Total	C	O	0	0
			4	2	2		
10	B	1	Total	C	O	0	0
			4	2	2		
10	B	1	Total	C	O	0	0
			4	2	2		
10	B	1	Total	C	O	0	0
			4	2	2		

- Molecule 11 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
11	A	258	Total	O	0	0
			258	258		
11	B	229	Total	O	0	0
			229	229		

- Molecule 1: Prostaglandin G/H synthase 2



MAG1
MAG2

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

Chain D:  100%

MAG1
MAG2
MAN3

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

Chain E:  50% 50%

MAG1
RDG2

4 Data and refinement statistics

Property	Value	Source
Space group	I 2 2 2	Depositor
Cell constants a, b, c, α , β , γ	121.85Å 132.34Å 180.29Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.95 – 2.40 19.95 – 2.40	Depositor EDS
% Data completeness (in resolution range)	99.9 (19.95-2.40) 99.9 (19.95-2.40)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.35 (at 2.41Å)	Xtriage
Refinement program	REFMAC 5.5.0088	Depositor
R, R_{free}	0.176 , 0.223 0.189 , 0.237	Depositor DCC
R_{free} test set	2889 reflections (5.07%)	wwPDB-VP
Wilson B-factor (Å ²)	34.8	Xtriage
Anisotropy	0.045	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 43.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	9765	wwPDB-VP
Average B, all atoms (Å ²)	35.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.23% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: AKR, NAG, EDO, NDG, COH, BOG, EPA, MAN

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	0.44	0/4600	0.53	0/6249
1	B	0.44	0/4578	0.54	0/6218
All	All	0.44	0/9178	0.53	0/12467

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

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	4463	0	4295	28	0
1	B	4442	0	4263	42	0
2	C	28	0	25	0	0
2	F	28	0	25	1	0
3	D	39	0	34	3	0
4	E	28	0	24	0	0
5	A	22	0	29	0	0
5	B	22	0	29	9	0
6	A	10	0	6	5	0
6	B	10	0	6	5	0
7	A	43	0	30	0	0

Continued on next page...

Continued from previous page...

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
7	B	43	0	30	0	0
8	A	14	0	13	0	0
8	B	14	0	13	0	0
9	A	20	0	28	1	0
9	B	20	0	28	5	0
10	A	20	0	30	2	0
10	B	12	0	18	2	0
11	A	258	0	0	3	0
11	B	229	0	0	1	0
All	All	9765	0	8926	70	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 4.

All (70) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:533:GLY:HA3	5:B:1:EPA:H181	1.44	0.99
1:A:477:SER:HB2	6:A:2:AKR:HA1	1.51	0.93
1:B:477:SER:HB2	6:B:2:AKR:HA1	1.51	0.90
1:A:216:ARG:HH11	3:D:2:NAG:H83	1.43	0.81
1:B:185:ARG:CZ	9:B:703:BOG:H4'2	2.11	0.81
1:B:36:CYS:HG	1:B:47:CYS:HG	0.91	0.79
1:B:161:THR:HG21	1:B:166:LYS:O	1.84	0.77
1:B:534:LEU:HG	5:B:1:EPA:H171	1.70	0.72
1:B:533:GLY:HA3	5:B:1:EPA:C18	2.19	0.72
1:A:492:LYS:HD3	6:A:2:AKR:HB2	1.72	0.72
1:A:161:THR:HG21	1:A:166:LYS:O	1.91	0.70
1:A:216:ARG:NH1	3:D:2:NAG:H83	2.07	0.70
1:A:532:LYS:HE2	11:A:10:HOH:O	1.93	0.69
1:B:272:GLU:HG3	6:B:3:AKR:HB2	1.76	0.67
1:B:533:GLY:CA	5:B:1:EPA:H181	2.24	0.66
1:A:216:ARG:HH11	3:D:2:NAG:C8	2.10	0.65
1:A:414:LEU:HD11	1:A:419:LEU:HD12	1.78	0.65
1:A:289:GLN:HG2	1:A:292:PHE:CE1	2.33	0.63
1:A:294:LEU:HG	1:A:295:VAL:HG23	1.81	0.63
1:A:477:SER:CB	6:A:2:AKR:HA1	2.29	0.61
1:B:216:ARG:HG2	2:F:2:NAG:H81	1.83	0.61
1:B:492:LYS:HE2	6:B:2:AKR:HB2	1.81	0.60
1:A:543:GLN:O	1:B:137:LYS:HE2	2.01	0.60
1:B:386:HIS:HB3	1:B:388[B]:HIS:CE1	2.37	0.60

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:276:PRO:HD2	1:A:279:ILE:HD12	1.86	0.58
1:A:230:LEU:HD13	1:A:233:ILE:HD12	1.86	0.58
1:B:162:PRO:O	10:B:7:EDO:H12	2.06	0.56
1:B:459:LYS:HA	10:B:7:EDO:H11	1.86	0.56
1:B:180:LYS:HB3	1:B:490:GLU:HG2	1.88	0.56
1:B:454:GLN:HA	1:B:457:GLU:HG2	1.88	0.55
1:B:185:ARG:HG3	9:B:703:BOG:H7'2	1.89	0.54
1:A:459:LYS:HA	10:A:4:EDO:H11	1.88	0.54
1:B:185:ARG:NE	9:B:703:BOG:H5'1	2.24	0.53
1:A:33:HIS:N	11:A:748:HOH:O	2.42	0.53
1:B:124:ILE:HD11	1:B:528:PRO:HB2	1.92	0.52
1:B:483:GLY:HA3	1:B:511:LYS:HD3	1.91	0.52
1:B:382:ASN:O	1:B:386:HIS:HD2	1.93	0.52
1:B:477:SER:HB2	6:B:2:AKR:CA	2.33	0.52
1:B:377:ILE:CD1	5:B:1:EPA:H201	2.40	0.51
11:A:668:HOH:O	1:B:543:GLN:HB2	2.11	0.50
1:B:36:CYS:CB	1:B:47:CYS:HG	2.24	0.49
1:A:386:HIS:HB3	1:A:388[B]:HIS:CE1	2.49	0.48
1:B:206:THR:HG21	1:B:385:TYR:CE2	2.48	0.48
1:A:161:THR:HG22	1:A:164:GLY:H	1.78	0.47
1:B:185:ARG:NH2	9:B:703:BOG:H4'2	2.29	0.47
1:A:230:LEU:HD22	1:A:232:HIS:HE1	1.81	0.46
1:A:161:THR:HG22	1:A:164:GLY:N	2.31	0.46
1:A:380:GLU:HG2	1:A:466:TYR:CE2	2.52	0.45
1:A:240:ARG:HH11	6:A:3:AKR:HA1	1.81	0.45
1:B:381:PHE:HE2	5:B:1:EPA:H192	1.81	0.45
1:B:344:VAL:O	1:B:348:TYR:HB3	2.17	0.44
1:B:441:PRO:HG2	1:B:444:VAL:HG22	1.99	0.44
1:B:161:THR:HG23	11:B:694:HOH:O	2.17	0.43
1:B:161:THR:CG2	1:B:166:LYS:O	2.60	0.43
1:A:344:VAL:HA	1:A:348:TYR:HB3	2.00	0.43
1:B:377:ILE:HD11	5:B:1:EPA:H201	2.01	0.43
1:A:162:PRO:O	10:A:4:EDO:H12	2.19	0.42
1:B:385:TYR:OH	5:B:1:EPA:H131	2.19	0.42
1:A:398:GLU:HB3	1:A:399:ASP:H	1.72	0.42
1:A:477:SER:HB2	6:A:2:AKR:CA	2.36	0.42
1:B:381:PHE:HE2	5:B:1:EPA:C19	2.33	0.41
1:B:397:ILE:HD11	1:B:426:PHE:CE1	2.55	0.41
1:B:34:HIS:HA	1:B:35:PRO:HD3	1.92	0.41
1:A:398:GLU:OE1	1:A:421:GLN:HG2	2.20	0.41
9:A:703:BOG:H6'1	9:A:703:BOG:H3'1	1.70	0.41

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:240:ARG:NH2	1:B:272:GLU:O	2.33	0.41
1:B:472:LEU:HD21	1:B:524:GLU:HG3	2.02	0.41
1:B:185:ARG:HG3	9:B:703:BOG:H5'1	2.03	0.41
1:B:492:LYS:HD3	6:B:2:AKR:HB2	2.03	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	552/591 (93%)	537 (97%)	13 (2%)	2 (0%)	34	48
1	B	552/591 (93%)	537 (97%)	15 (3%)	0	100	100
All	All	1104/1182 (93%)	1074 (97%)	28 (2%)	2 (0%)	51	62

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	409[A]	TYR
1	A	409[B]	TYR

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.

Continued on next page...

Continued from previous page...

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
1	A	484/529 (92%)	474 (98%)	10 (2%)	53 72
1	B	477/529 (90%)	470 (98%)	7 (2%)	65 80
All	All	961/1058 (91%)	944 (98%)	17 (2%)	60 76

All (17) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	161	THR
1	A	165	VAL
1	A	185	ARG
1	A	289	GLN
1	A	385	TYR
1	A	409[A]	TYR
1	A	409[B]	TYR
1	A	422	PHE
1	A	484	GLU
1	A	556	PHE
1	B	161	THR
1	B	289	GLN
1	B	336	LEU
1	B	368	ASN
1	B	385	TYR
1	B	442	ILE
1	B	484	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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 ⓘ

9 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	0.59	0	17,19,21	1.17	1 (5%)
2	NAG	C	2	2	14,14,15	0.45	0	17,19,21	1.18	2 (11%)
3	NAG	D	1	1,3	14,14,15	0.58	0	17,19,21	1.10	1 (5%)
3	NAG	D	2	3	14,14,15	0.62	0	17,19,21	0.92	0
3	MAN	D	3	3	11,11,12	0.53	0	15,15,17	1.50	2 (13%)
4	NAG	E	1	4,1	14,14,15	0.59	0	17,19,21	1.00	2 (11%)
4	NDG	E	2	4	14,14,15	0.71	0	17,19,21	0.96	0
2	NAG	F	1	1,2	14,14,15	0.62	0	17,19,21	0.97	1 (5%)
2	NAG	F	2	2	14,14,15	0.55	0	17,19,21	0.82	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	C	1	1,2	2/2/5/7	2/6/23/26	0/1/1/1
2	NAG	C	2	2	-	3/6/23/26	0/1/1/1
3	NAG	D	1	1,3	-	0/6/23/26	0/1/1/1
3	NAG	D	2	3	-	4/6/23/26	0/1/1/1
3	MAN	D	3	3	-	2/2/19/22	1/1/1/1
4	NAG	E	1	4,1	-	2/6/23/26	0/1/1/1
4	NDG	E	2	4	-	2/6/23/26	0/1/1/1
2	NAG	F	1	1,2	-	0/6/23/26	0/1/1/1
2	NAG	F	2	2	-	2/6/23/26	0/1/1/1

There are no bond length outliers.

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	3	MAN	C1-O5-C5	4.27	117.97	112.19
3	D	1	NAG	C1-O5-C5	3.84	117.40	112.19
2	F	1	NAG	C1-O5-C5	3.51	116.94	112.19
2	C	1	NAG	C1-O5-C5	3.13	116.43	112.19
3	D	3	MAN	O5-C5-C6	2.80	111.59	107.20
4	E	1	NAG	C4-C3-C2	2.40	114.54	111.02
4	E	1	NAG	C1-O5-C5	2.40	115.44	112.19
2	C	2	NAG	O5-C5-C6	2.31	110.83	107.20
2	C	2	NAG	C1-O5-C5	2.20	115.17	112.19

All (2) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
2	C	1	NAG	C1
2	C	1	NAG	C5

All (17) torsion outliers are listed below:

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

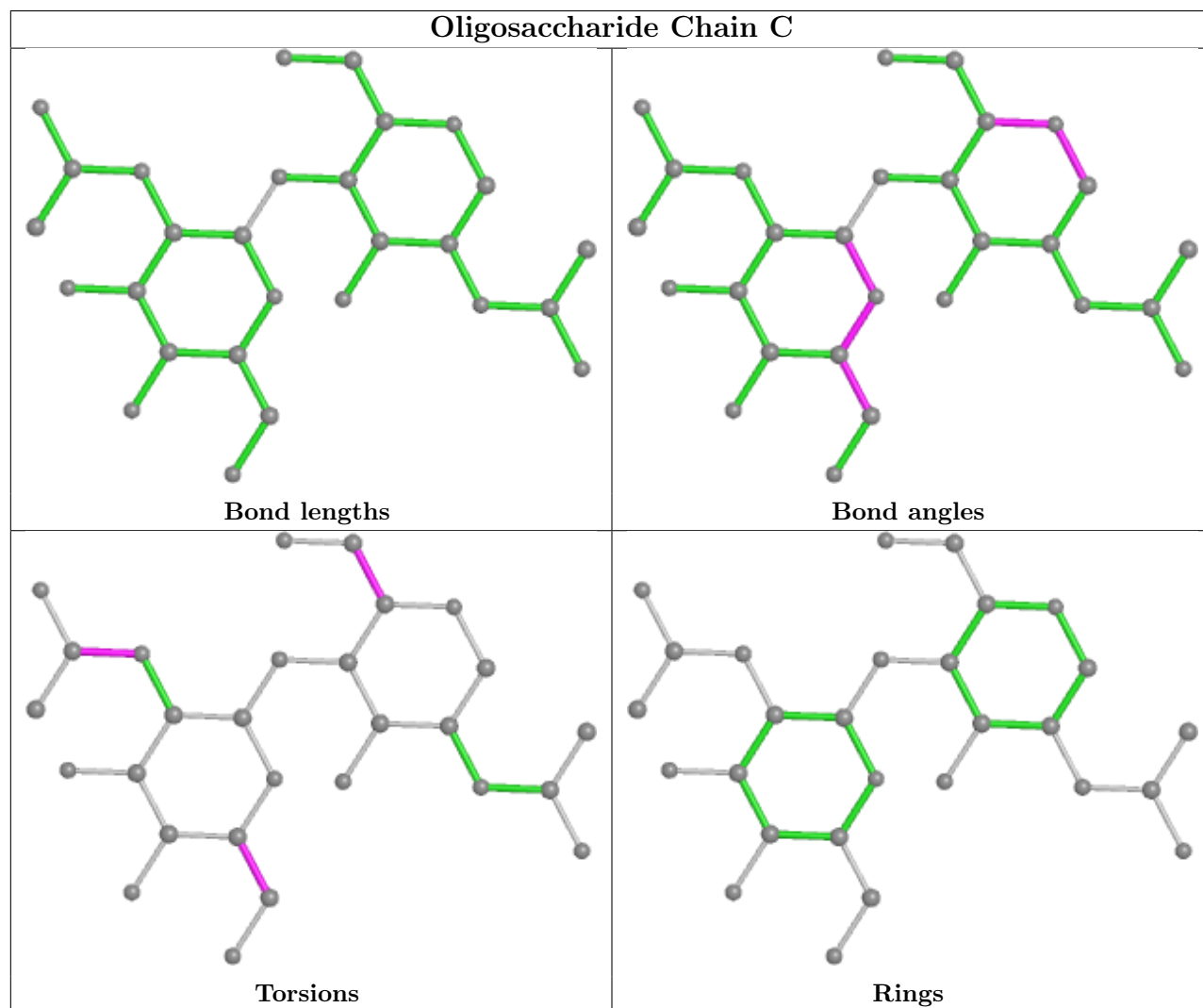
All (1) ring outliers are listed below:

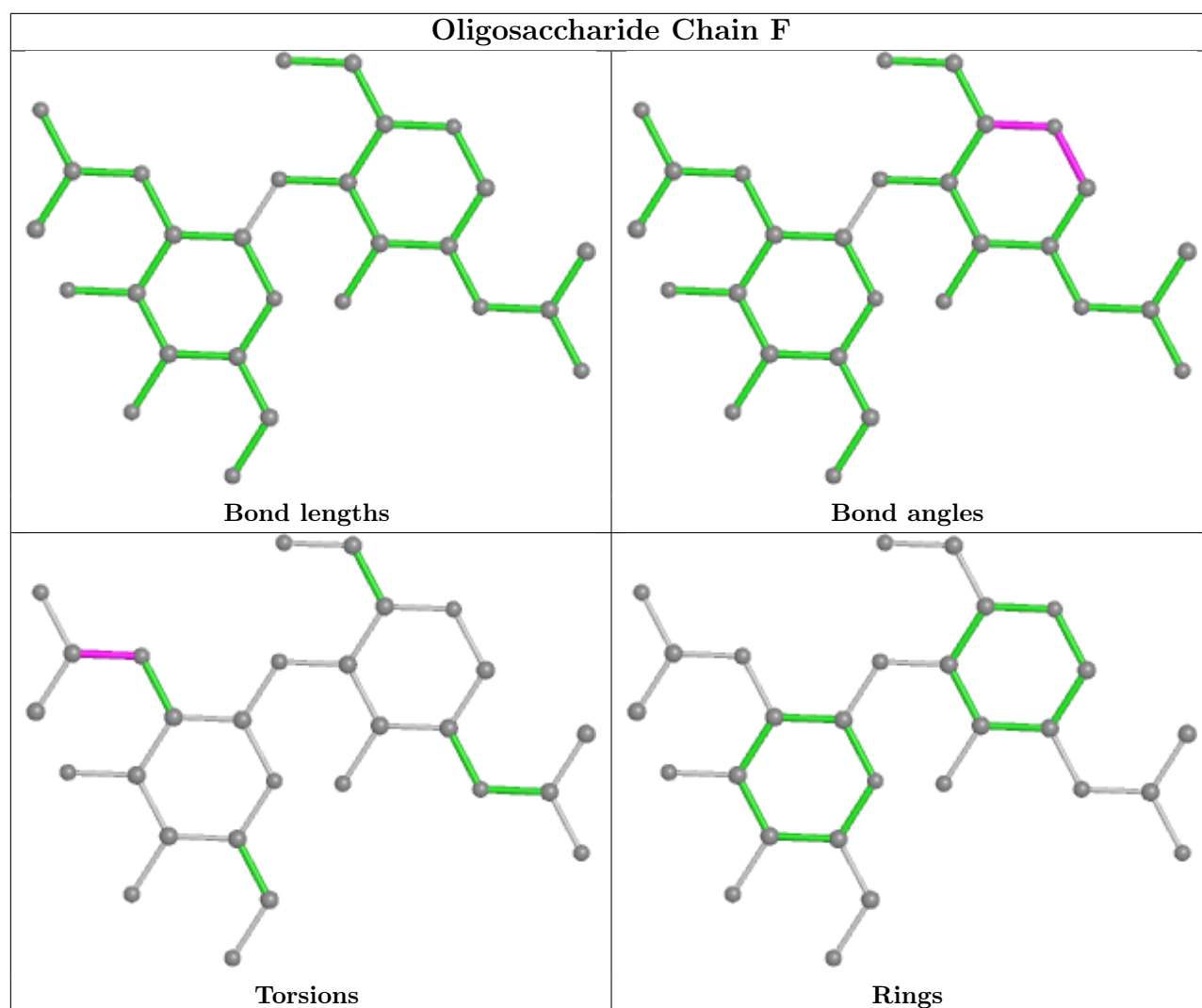
Mol	Chain	Res	Type	Atoms
3	D	3	MAN	C1-C2-C3-C4-C5-O5

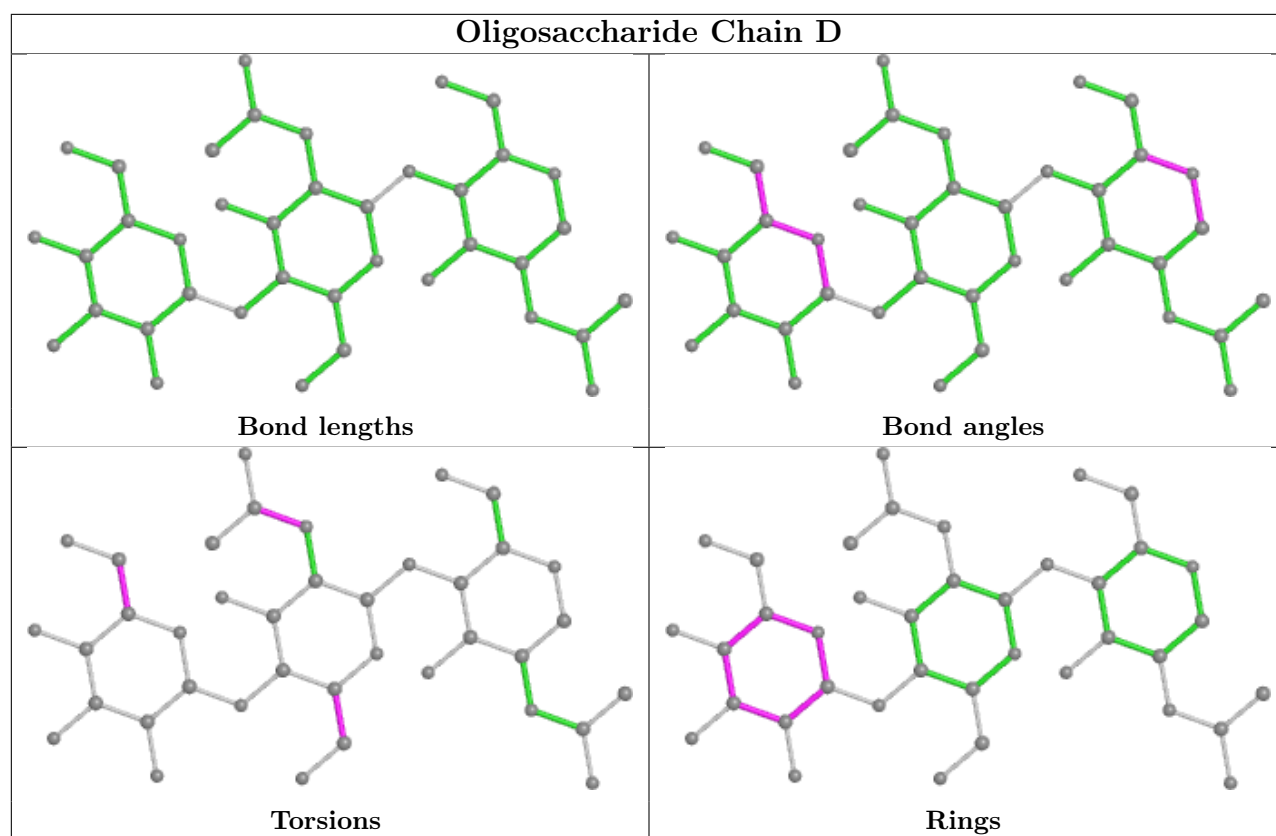
2 monomers are involved in 4 short contacts:

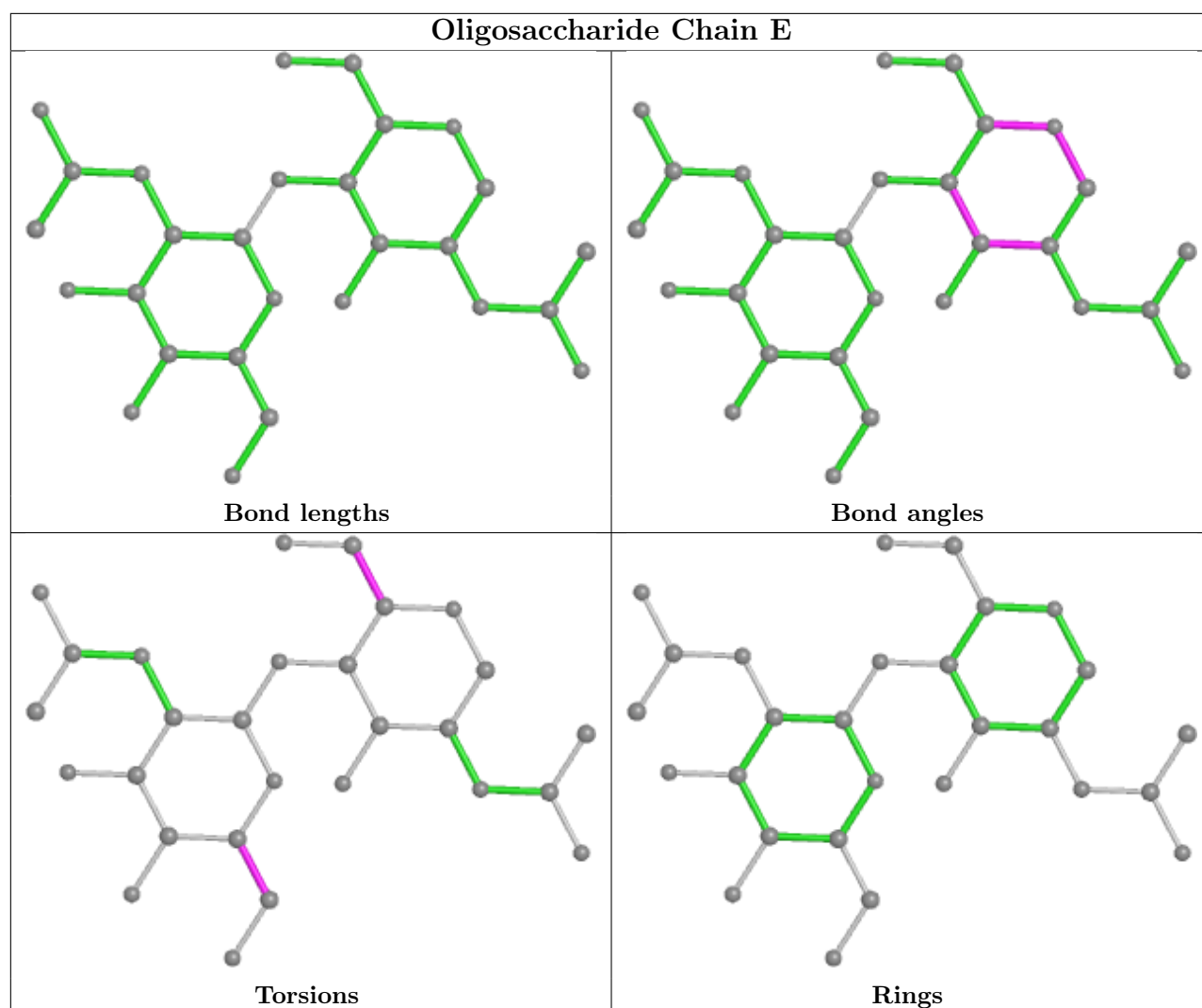
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	F	2	NAG	1	0
3	D	2	NAG	3	0

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









5.6 Ligand geometry [i](#)

20 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	EPA	B	1	-	18,21,21	0.22	0	16,21,21	0.66	0
10	EDO	A	8	-	3,3,3	0.44	0	2,2,2	0.42	0
6	AKR	A	3	-	1,4,4	1.65	0	0,4,4	-	-
10	EDO	B	7	-	3,3,3	0.36	0	2,2,2	0.49	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
10	EDO	B	620	-	3,3,3	0.54	0	2,2,2	0.24	0
7	COH	B	619	1	31,50,50	5.56	22 (70%)	23,82,82	4.58	9 (39%)
6	AKR	B	3	-	1,4,4	1.61	0	0,4,4	-	-
10	EDO	A	4	-	3,3,3	0.44	0	2,2,2	0.43	0
6	AKR	B	2	-	1,4,4	1.61	0	0,4,4	-	-
9	BOG	B	703	-	20,20,20	0.52	0	25,25,25	0.67	0
8	NAG	B	681	1	14,14,15	0.55	0	17,19,21	0.76	0
10	EDO	B	621	-	3,3,3	0.35	0	2,2,2	0.53	0
10	EDO	A	6	-	3,3,3	0.50	0	2,2,2	0.39	0
6	AKR	A	2	-	1,4,4	1.59	0	0,4,4	-	-
9	BOG	A	703	-	20,20,20	0.45	0	25,25,25	0.70	0
10	EDO	A	620	-	3,3,3	0.54	0	2,2,2	0.24	0
8	NAG	A	681	1	14,14,15	0.58	0	17,19,21	1.11	1 (5%)
7	COH	A	619	1	31,50,50	5.52	22 (70%)	23,82,82	4.58	8 (34%)
10	EDO	A	5	-	3,3,3	0.46	0	2,2,2	0.39	0
5	EPA	A	1	-	18,21,21	0.28	0	16,21,21	0.48	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
5	EPA	B	1	-	-	6/17/19/19	-
10	EDO	A	8	-	-	1/1/1/1	-
6	AKR	A	3	-	-	0/0/2/2	-
10	EDO	B	7	-	-	1/1/1/1	-
10	EDO	B	620	-	-	0/1/1/1	-
7	COH	B	619	1	-	4/8/94/94	-
6	AKR	B	3	-	-	0/0/2/2	-
10	EDO	A	4	-	-	1/1/1/1	-
6	AKR	B	2	-	-	0/0/2/2	-
9	BOG	B	703	-	-	6/11/31/31	0/1/1/1
8	NAG	B	681	1	-	3/6/23/26	0/1/1/1
10	EDO	B	621	-	-	1/1/1/1	-
10	EDO	A	6	-	-	1/1/1/1	-
6	AKR	A	2	-	-	0/0/2/2	-
9	BOG	A	703	-	-	5/11/31/31	0/1/1/1
10	EDO	A	620	-	-	0/1/1/1	-
8	NAG	A	681	1	-	0/6/23/26	0/1/1/1
7	COH	A	619	1	-	1/8/94/94	-

Continued on next page...

Continued from previous page...

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
10	EDO	A	5	-	-	0/1/1/1	-
5	EPA	A	1	-	-	5/17/19/19	-

All (44) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
7	B	619	COH	C1D-ND	10.75	1.40	1.34
7	A	619	COH	C1D-ND	10.51	1.39	1.34
7	B	619	COH	C4D-ND	10.15	1.39	1.34
7	A	619	COH	C4D-ND	9.89	1.39	1.34
7	B	619	COH	C4B-NB	9.38	1.39	1.34
7	A	619	COH	C4B-NB	9.16	1.39	1.34
7	A	619	COH	C1B-NB	9.13	1.39	1.34
7	B	619	COH	C1B-NB	9.09	1.39	1.34
7	A	619	COH	CHD-C4C	-8.57	1.38	1.53
7	B	619	COH	CHD-C4C	-8.44	1.38	1.53
7	B	619	COH	CHA-C1A	-8.32	1.39	1.53
7	A	619	COH	CHA-C1A	-8.31	1.39	1.53
7	A	619	COH	CHB-C4A	-8.29	1.39	1.53
7	B	619	COH	CHB-C4A	-8.23	1.39	1.53
7	A	619	COH	CHC-C1C	-7.98	1.39	1.53
7	B	619	COH	CHC-C1C	-7.93	1.39	1.53
7	A	619	COH	C1B-C2B	5.60	1.45	1.38
7	B	619	COH	C3D-C2D	5.59	1.54	1.37
7	B	619	COH	C1B-C2B	5.55	1.45	1.38
7	A	619	COH	C3D-C2D	5.41	1.53	1.37
7	A	619	COH	C4D-C3D	5.33	1.45	1.38
7	B	619	COH	C4D-C3D	5.32	1.45	1.38
7	B	619	COH	CHD-C1D	-5.12	1.38	1.51
7	A	619	COH	CHD-C1D	-5.11	1.38	1.51
7	A	619	COH	C1D-C2D	5.07	1.45	1.38
7	B	619	COH	C1D-C2D	4.99	1.45	1.38
7	A	619	COH	CHA-C4D	-4.99	1.39	1.51
7	B	619	COH	CHB-C1B	-4.98	1.39	1.51
7	A	619	COH	CHB-C1B	-4.98	1.39	1.51
7	B	619	COH	CHA-C4D	-4.97	1.39	1.51
7	A	619	COH	CHC-C4B	-4.65	1.39	1.51
7	B	619	COH	CHC-C4B	-4.59	1.40	1.51
7	A	619	COH	C3B-C2B	-4.23	1.34	1.40
7	B	619	COH	C3B-C2B	-4.10	1.34	1.40
7	A	619	COH	C3B-CAB	3.63	1.55	1.47
7	B	619	COH	C4C-C3C	-3.61	1.45	1.51

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
7	A	619	COH	C4C-C3C	-3.58	1.45	1.51
7	B	619	COH	C3B-CAB	3.55	1.55	1.47
7	B	619	COH	CAA-C2A	3.32	1.55	1.51
7	A	619	COH	CAA-C2A	3.27	1.55	1.51
7	B	619	COH	CMA-C3A	3.23	1.55	1.50
7	A	619	COH	CMA-C3A	3.04	1.55	1.50
7	B	619	COH	CMC-C2C	2.94	1.55	1.50
7	A	619	COH	CMC-C2C	2.83	1.55	1.50

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	A	619	COH	CHB-C4A-NA	10.53	125.61	110.12
7	A	619	COH	CHA-C1A-NA	10.44	125.48	110.12
7	A	619	COH	CHC-C1C-NC	10.38	125.40	110.12
7	A	619	COH	CHD-C4C-NC	10.36	125.37	110.12
7	B	619	COH	CHD-C4C-NC	10.31	125.29	110.12
7	B	619	COH	CHA-C1A-NA	10.29	125.26	110.12
7	B	619	COH	CHC-C1C-NC	10.25	125.21	110.12
7	B	619	COH	CHB-C4A-NA	10.16	125.08	110.12
7	B	619	COH	CBD-CAD-C3D	-3.52	105.99	112.49
8	A	681	NAG	C1-O5-C5	3.47	116.90	112.19
7	B	619	COH	CBA-CAA-C2A	-3.21	108.66	114.35
7	B	619	COH	C4B-C3B-C2B	2.62	108.44	104.41
7	A	619	COH	CMC-C2C-C3C	-2.57	125.36	129.73
7	A	619	COH	C4B-C3B-C2B	2.50	108.25	104.41
7	A	619	COH	CBD-CAD-C3D	-2.42	108.02	112.49
7	B	619	COH	CMC-C2C-C3C	-2.37	125.70	129.73
7	A	619	COH	CHD-C1D-C2D	-2.08	125.86	129.45
7	B	619	COH	CHD-C1D-C2D	-2.01	125.97	129.45

There are no chirality outliers.

All (35) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	A	1	EPA	C1-C2-C3-C4
5	B	1	EPA	C12-C13-C14-C15
7	B	619	COH	C2A-CAA-CBA-CGA
9	A	703	BOG	C2'-C1'-O1-C1
9	B	703	BOG	O5-C1-O1-C1'
9	B	703	BOG	C4-C5-C6-O6
9	A	703	BOG	C3'-C4'-C5'-C6'

Continued on next page...

Continued from previous page...

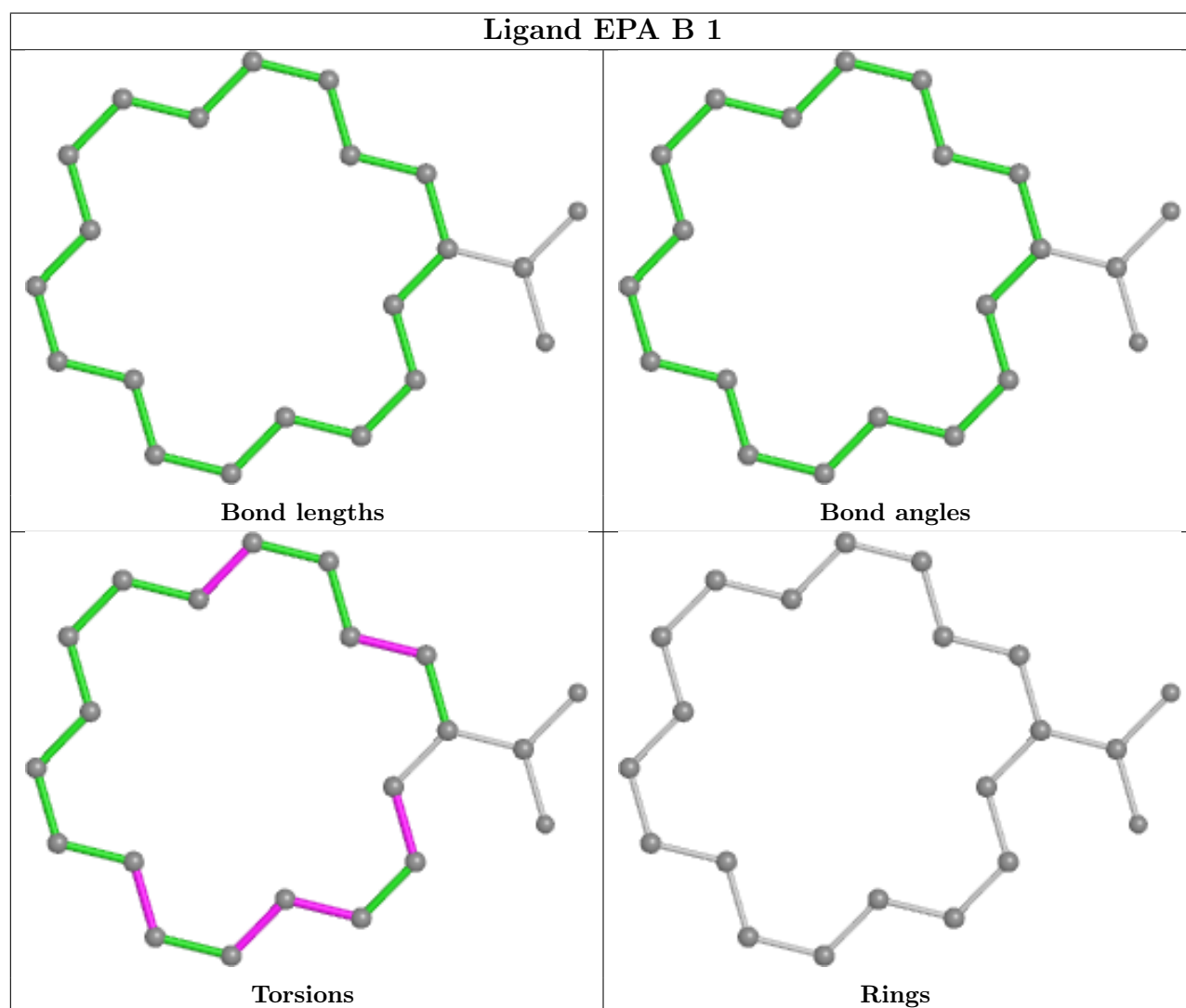
Mol	Chain	Res	Type	Atoms
9	B	703	BOG	O5-C5-C6-O6
9	B	703	BOG	O1-C1'-C2'-C3'
9	A	703	BOG	C2'-C3'-C4'-C5'
9	B	703	BOG	C3'-C4'-C5'-C6'
9	B	703	BOG	C4'-C5'-C6'-C7'
7	B	619	COH	C2C-C3C-CAC-CBC
10	A	4	EDO	O1-C1-C2-O2
10	B	621	EDO	O1-C1-C2-O2
10	B	7	EDO	O1-C1-C2-O2
5	B	1	EPA	C2-C3-C4-C5
7	B	619	COH	C4C-C3C-CAC-CBC
5	A	1	EPA	C2-C3-C4-C5
5	A	1	EPA	C11-C12-C13-C14
5	A	1	EPA	C15-C16-C17-C18
5	B	1	EPA	C5-C6-C7-C8
5	B	1	EPA	C14-C15-C16-C17
5	B	1	EPA	C15-C16-C17-C18
8	B	681	NAG	C8-C7-N2-C2
8	B	681	NAG	O7-C7-N2-C2
9	A	703	BOG	C1'-C2'-C3'-C4'
7	A	619	COH	C2A-CAA-CBA-CGA
5	A	1	EPA	C11-C10-C9-C8
9	A	703	BOG	O1-C1'-C2'-C3'
5	B	1	EPA	C17-C18-C19-C20
10	A	8	EDO	O1-C1-C2-O2
8	B	681	NAG	C4-C5-C6-O6
7	B	619	COH	C3A-C2A-CAA-CBA
10	A	6	EDO	O1-C1-C2-O2

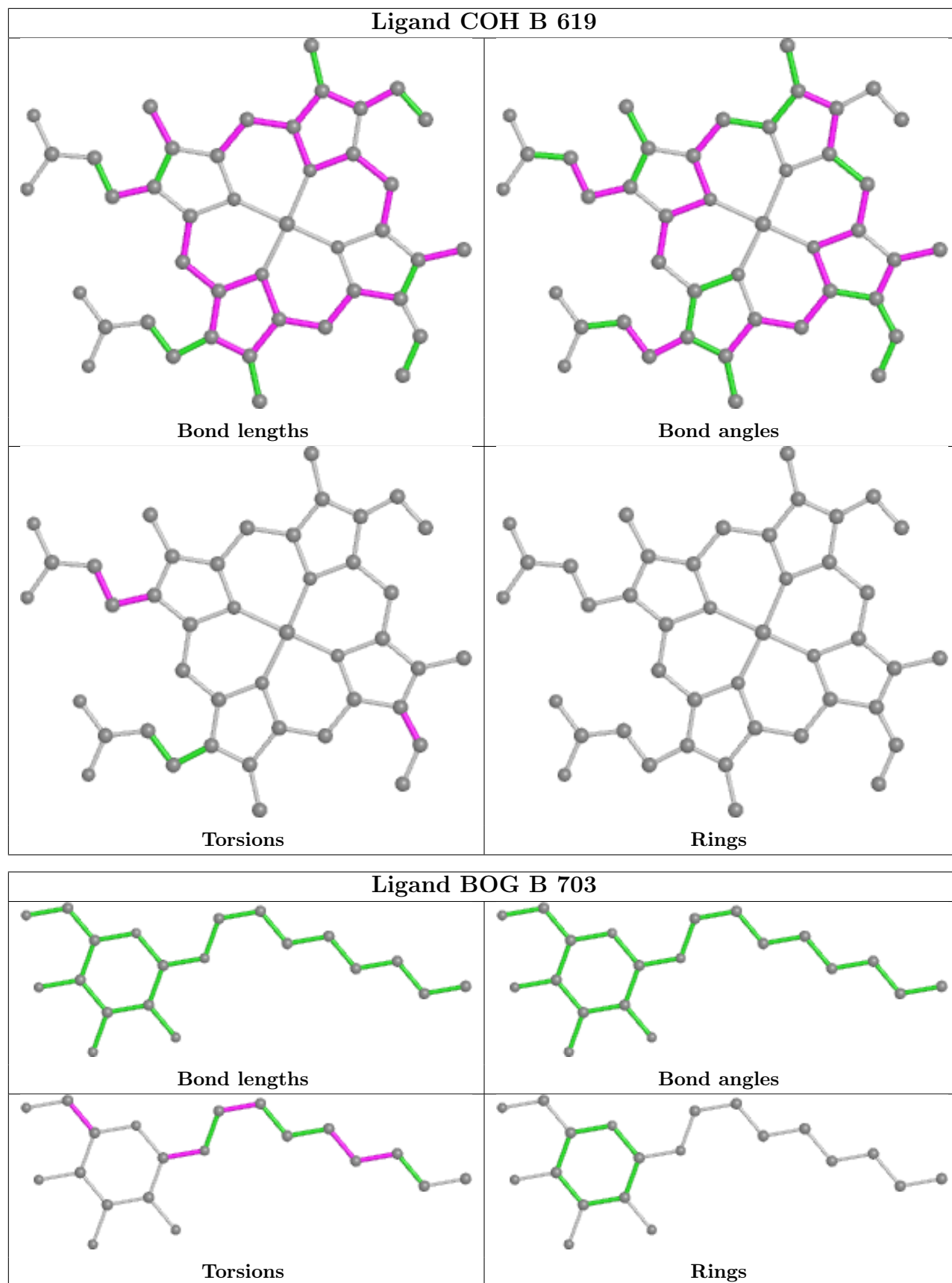
There are no ring outliers.

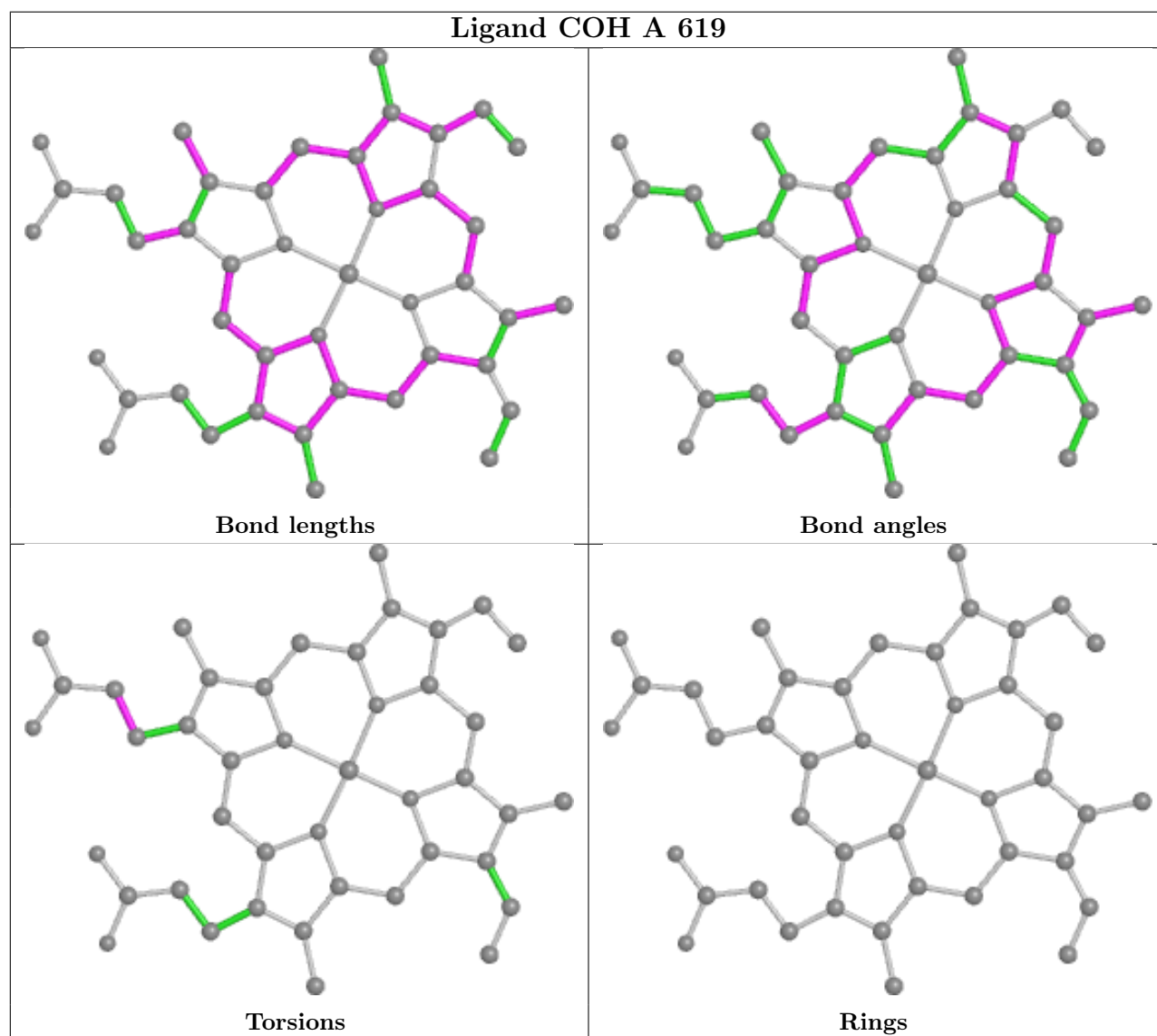
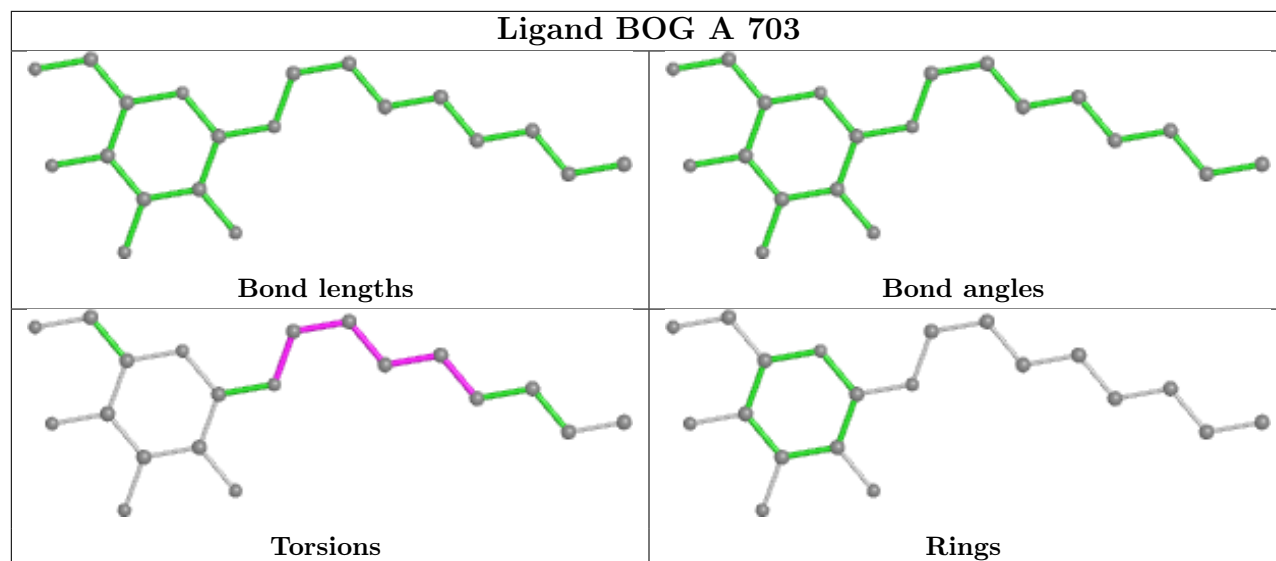
9 monomers are involved in 29 short contacts:

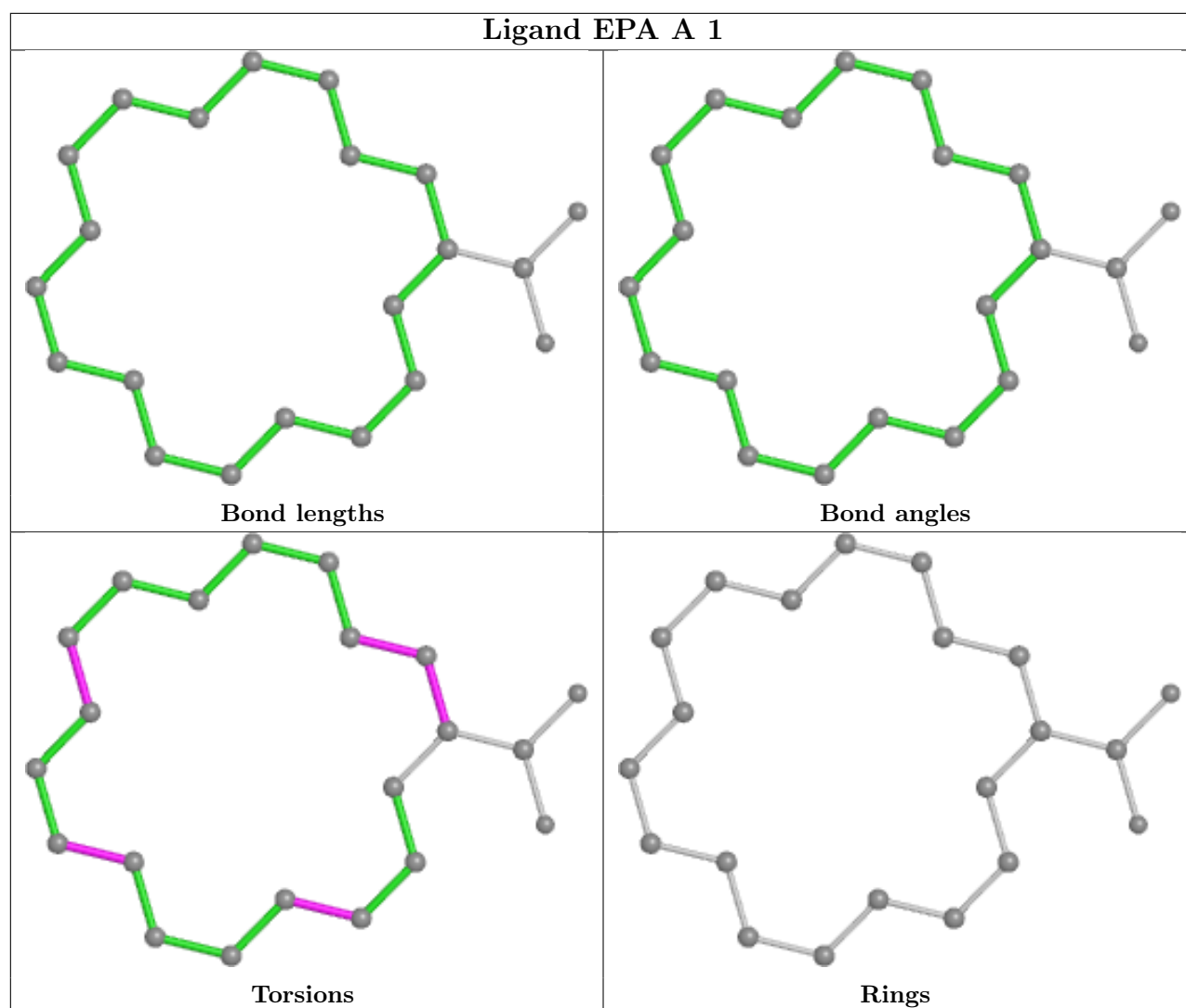
Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	B	1	EPA	9	0
6	A	3	AKR	1	0
10	B	7	EDO	2	0
6	B	3	AKR	1	0
10	A	4	EDO	2	0
6	B	2	AKR	4	0
9	B	703	BOG	5	0
6	A	2	AKR	4	0
9	A	703	BOG	1	0

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









5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	551/591 (93%)	-0.30	7 (1%) 77 75	17, 33, 51, 58	0
1	B	551/591 (93%)	-0.19	12 (2%) 62 60	20, 34, 53, 68	1 (0%)
All	All	1102/1182 (93%)	-0.24	19 (1%) 70 68	17, 34, 52, 68	1 (0%)

All (19) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	277	PRO	3.6
1	A	278	HIS	3.5
1	B	122	TYR	3.4
1	A	81	LEU	3.3
1	B	74	PHE	3.2
1	A	399	ASP	3.0
1	B	368	ASN	2.9
1	A	277	PRO	2.9
1	B	157	ASP	2.8
1	A	368	ASN	2.8
1	B	399	ASP	2.4
1	A	122	TYR	2.3
1	B	105	ASN	2.3
1	B	115	TYR	2.3
1	A	53	ASP	2.1
1	B	278	HIS	2.1
1	B	78	ILE	2.1
1	B	91	TYR	2.0
1	B	65	TYR	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

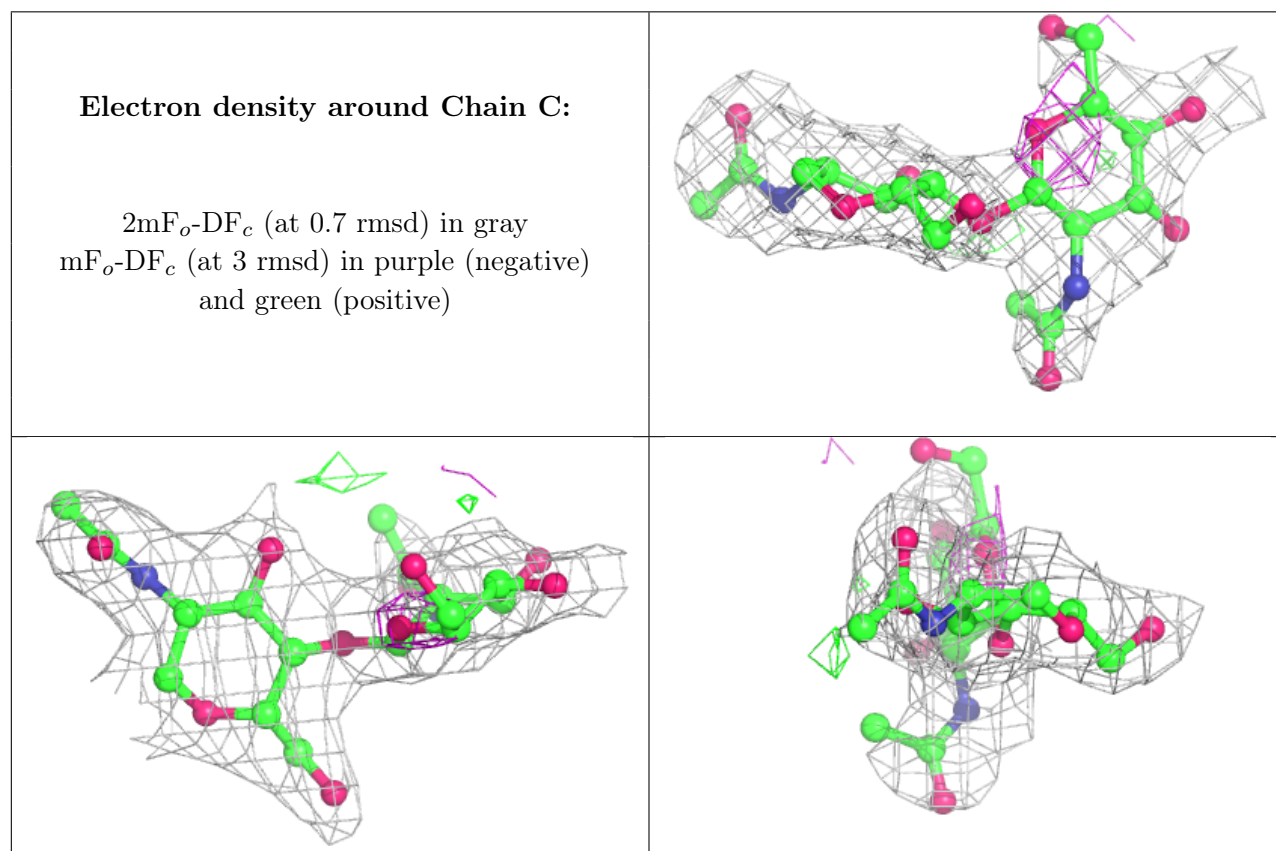
There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates ⓘ

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

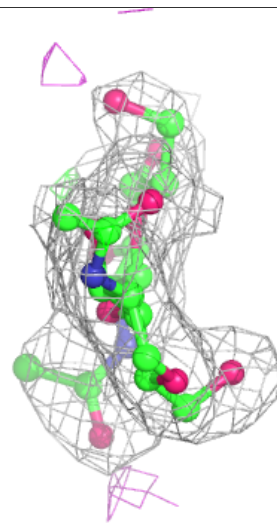
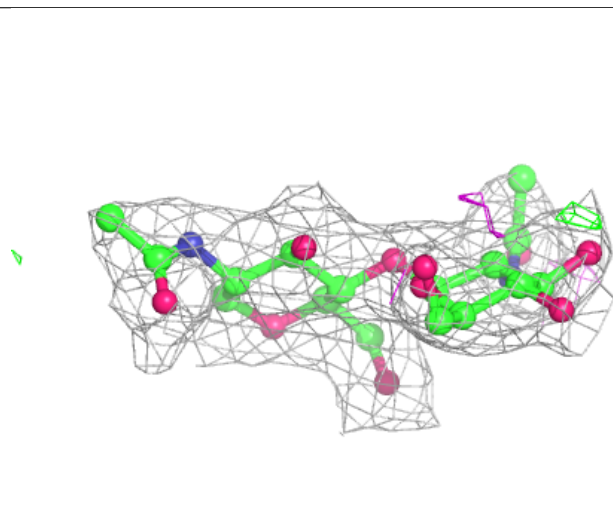
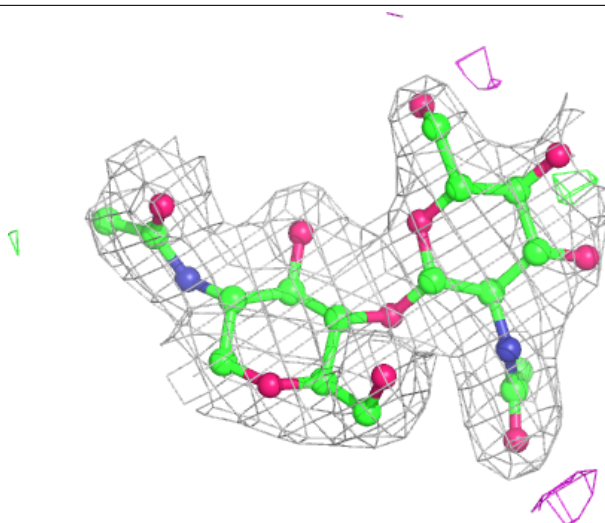
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	NAG	C	2	14/15	0.75	0.56	64,67,68,68	0
4	NDG	E	2	14/15	0.78	0.53	69,72,73,73	0
3	MAN	D	3	11/12	0.81	0.39	60,62,63,63	0
4	NAG	E	1	14/15	0.84	0.32	52,57,60,65	0
3	NAG	D	2	14/15	0.87	0.20	46,48,52,56	0
2	NAG	C	1	14/15	0.89	0.24	45,51,54,59	0
2	NAG	F	2	14/15	0.90	0.28	43,46,48,48	0
2	NAG	F	1	14/15	0.94	0.13	30,34,37,41	0
3	NAG	D	1	14/15	0.95	0.10	30,36,38,42	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.



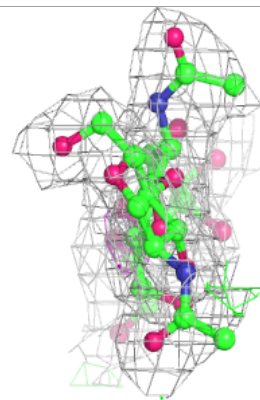
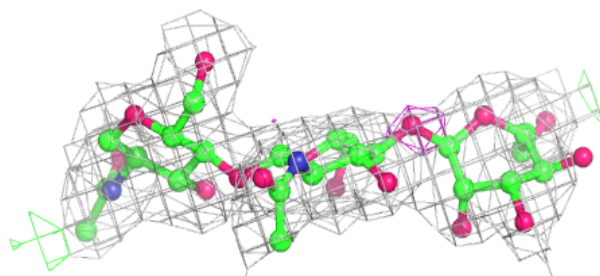
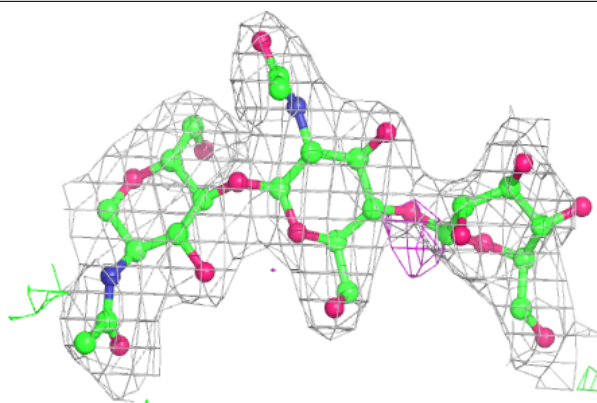
Electron density around Chain F:

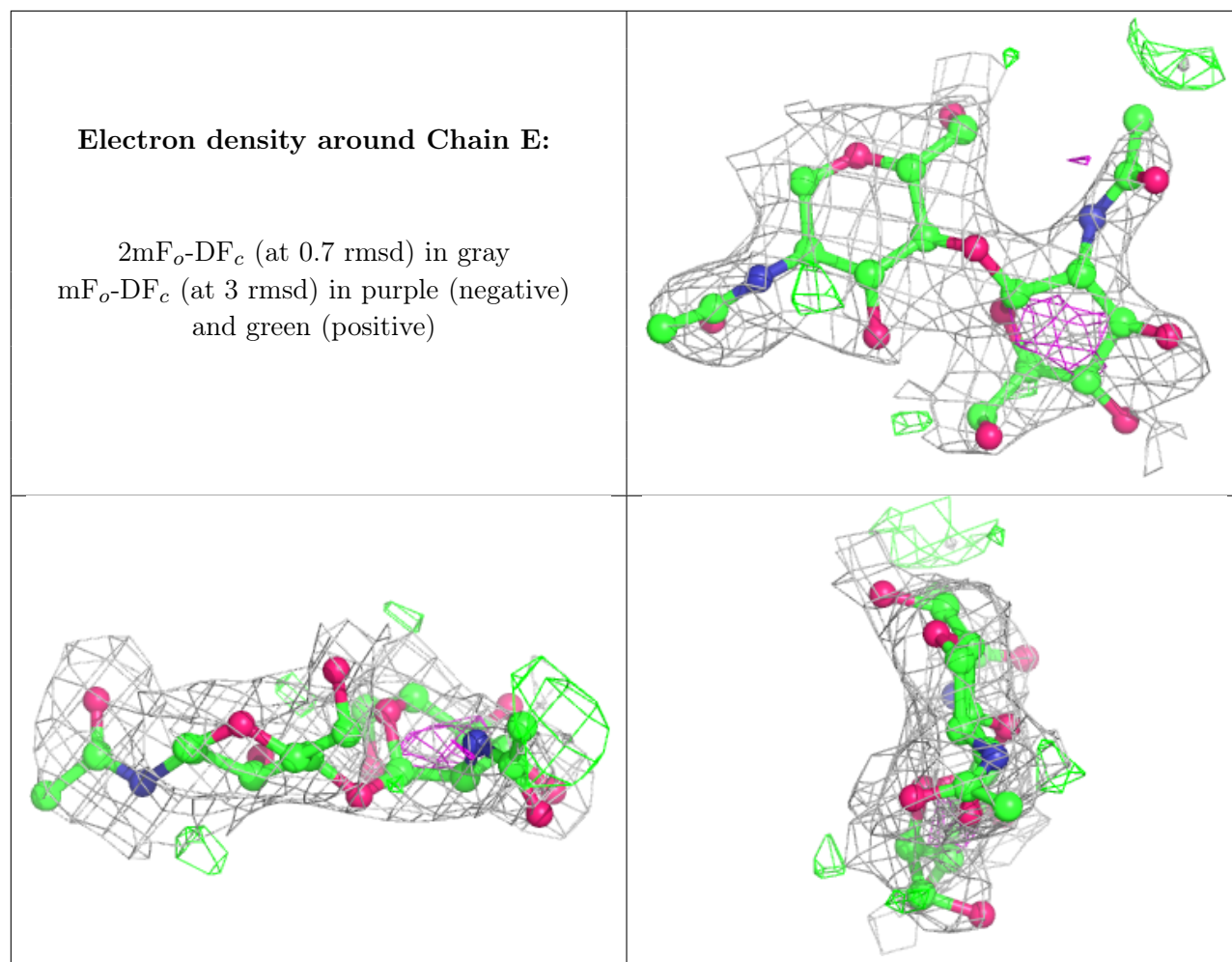
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around Chain D:

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





6.4 Ligands ⓘ

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
7	COH	B	619	43/43	0.64	0.36	80,85,87,88	0
7	COH	A	619	43/43	0.68	0.35	92,94,96,97	0
10	EDO	A	6	4/4	0.75	0.20	58,58,58,58	0
6	AKR	B	3	5/5	0.79	0.31	63,63,63,63	0
10	EDO	A	5	4/4	0.83	0.18	61,62,62,63	0
6	AKR	A	3	5/5	0.85	0.22	54,54,54,54	0
8	NAG	A	681	14/15	0.86	0.37	48,52,54,55	0
8	NAG	B	681	14/15	0.86	0.33	52,56,59,59	0
5	EPA	B	1	22/22	0.87	0.23	42,45,50,52	0
5	EPA	A	1	22/22	0.90	0.21	41,43,48,49	0

Continued on next page...

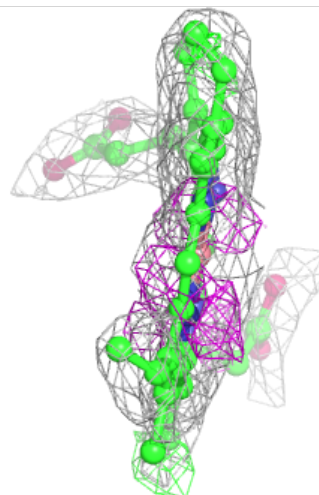
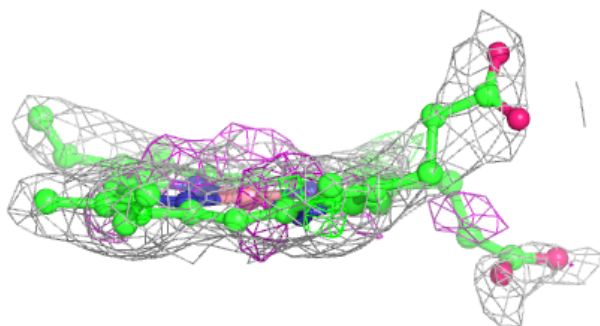
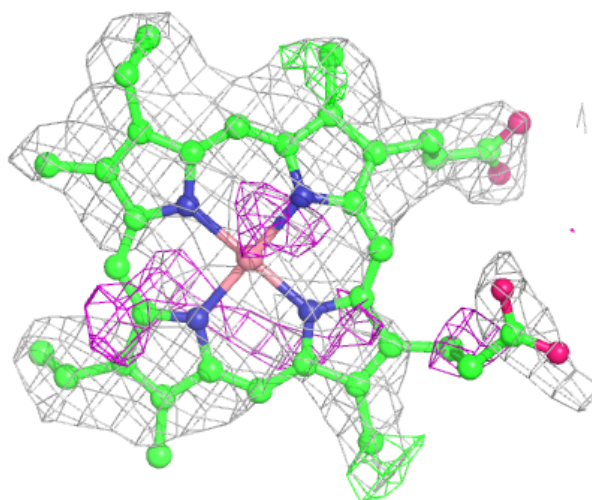
Continued from previous page...

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
6	AKR	B	2	5/5	0.91	0.24	65,65,66,66	0
6	AKR	A	2	5/5	0.92	0.28	52,52,52,52	0
10	EDO	A	4	4/4	0.92	0.18	38,39,41,41	0
10	EDO	A	620	4/4	0.93	0.21	36,36,37,37	0
10	EDO	B	621	4/4	0.93	0.14	33,33,34,34	0
10	EDO	A	8	4/4	0.94	0.17	39,39,41,42	0
10	EDO	B	620	4/4	0.94	0.19	34,35,36,36	0
9	BOG	B	703	20/20	0.94	0.14	36,37,37,37	20
10	EDO	B	7	4/4	0.94	0.12	39,39,39,40	0
9	BOG	A	703	20/20	0.96	0.14	22,23,27,27	20

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.

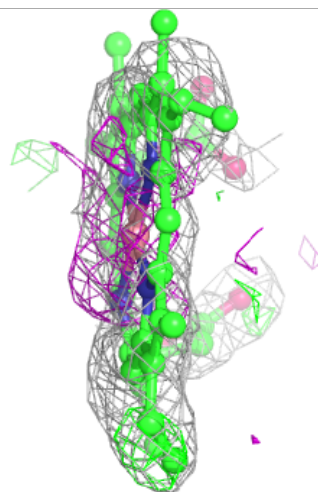
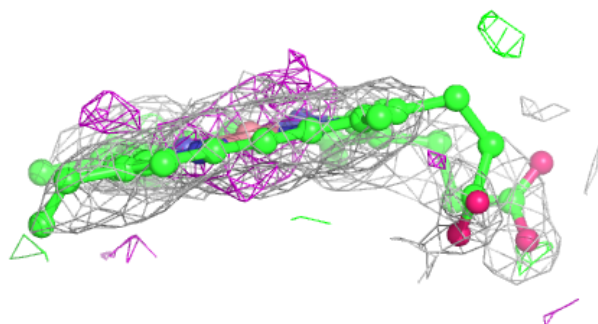
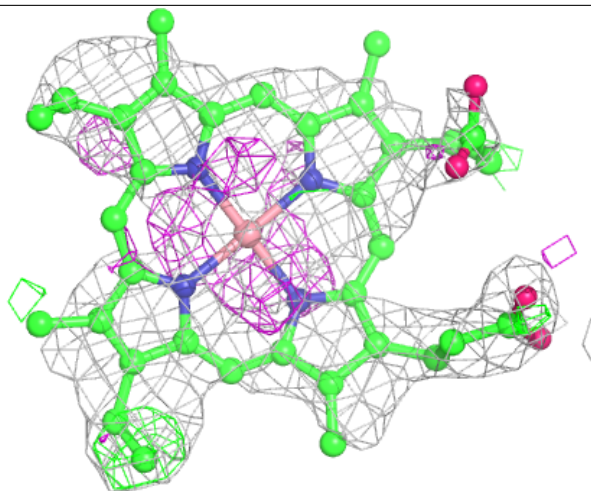
Electron density around COH B 619:

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



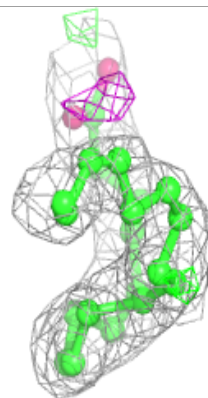
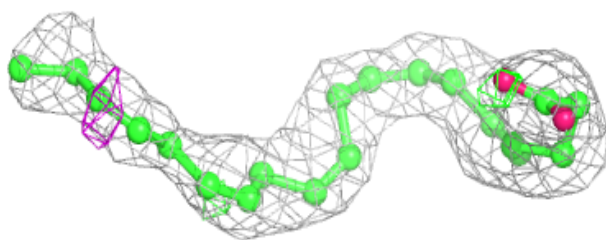
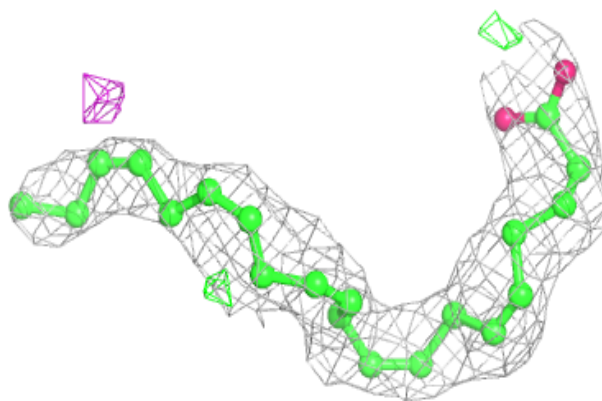
Electron density around COH A 619:

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



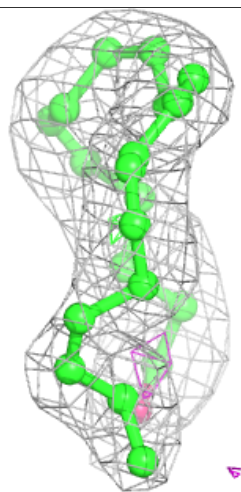
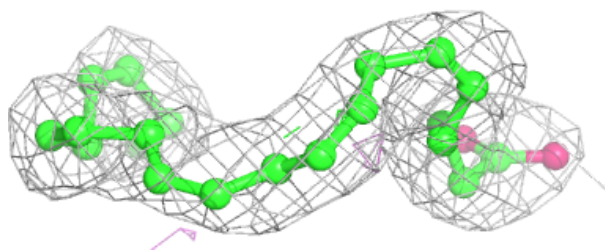
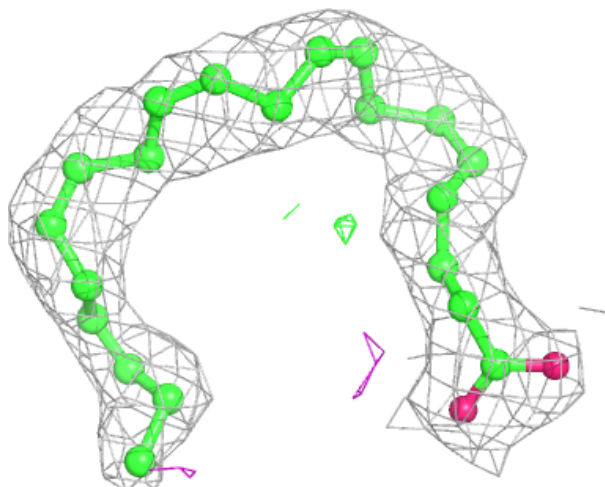
Electron density around EPA B 1:

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



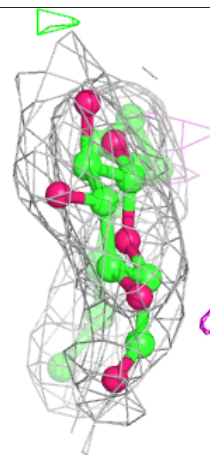
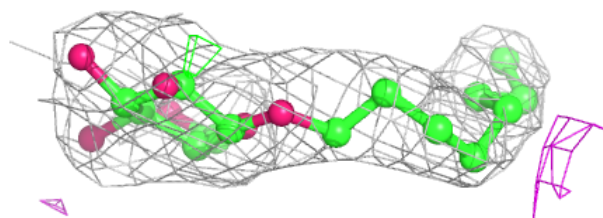
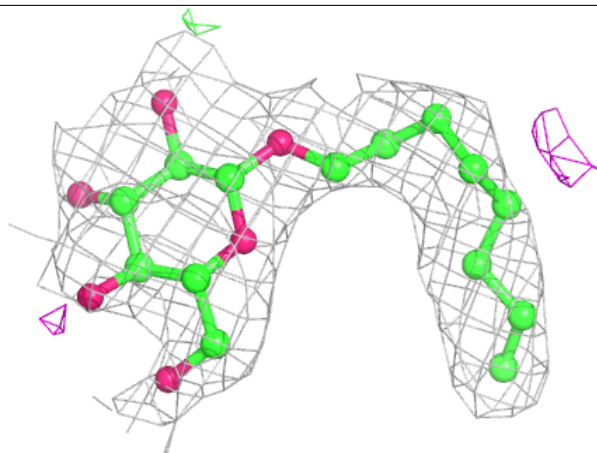
Electron density around EPA A 1:

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



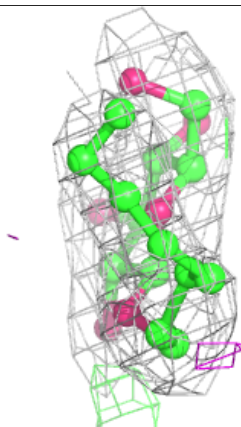
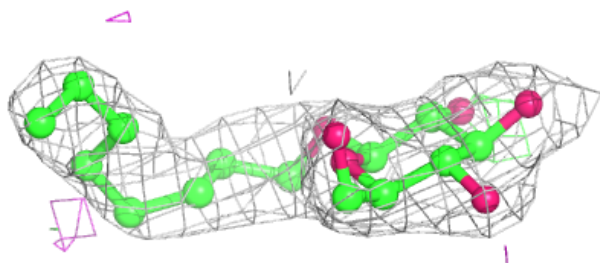
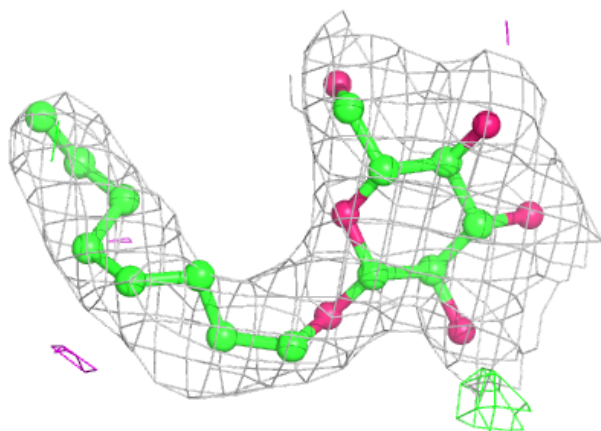
Electron density around BOG B 703:

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



Electron density around BOG A 703:

$2mF_o-DF_c$ (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.