



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 12:08 PM UTC

PDB ID : 8IC1 / pdb_00008ic1
Title : endo-alpha-D-arabinanase EndoMA1 D51N mutant from *Microbacterium arabinogalactanolyticum* in complex with arabinooligosaccharides
Authors : Li, J.; Nakashima, C.; Ishiwata, A.; Fujita, K.; Fushinobu, S.
Deposited on : 2023-02-10
Resolution : 1.80 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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

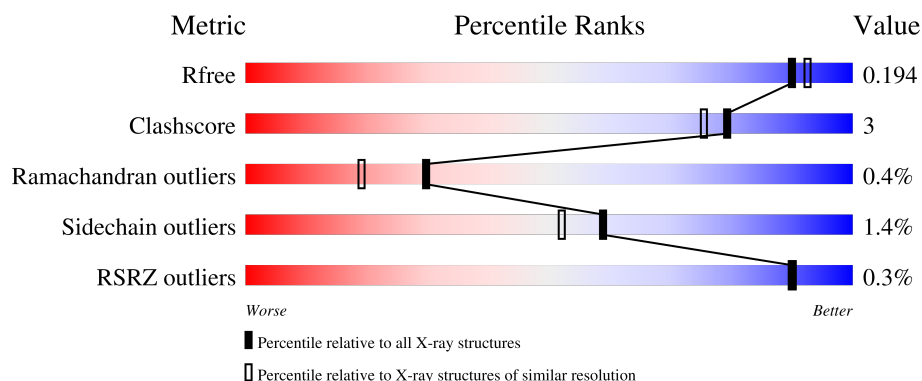
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.80 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.



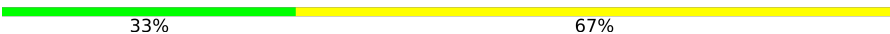
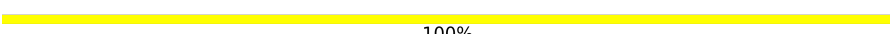
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	7662 (1.80-1.80)
Clashscore	190562	8479 (1.80-1.80)
Ramachandran outliers	187476	8391 (1.80-1.80)
Sidechain outliers	187428	8390 (1.80-1.80)
RSRZ outliers	180081	7663 (1.80-1.80)

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 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	488	90% 7% ..
1	B	488	91% 7% .
1	C	488	91% 7% .
1	D	488	90% 8% .
2	E	4	100%

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Mol	Chain	Length	Quality of chain
2	I	4	 25% 75%
3	F	3	 33% 67%
3	G	3	 33% 67%
3	H	3	 67% 33%
3	J	3	 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
6	PEG	D	505	-	-	X	-

2 Entry composition [i](#)

There are 11 unique types of molecules in this entry. The entry contains 16610 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 endo-alpha-D-arabinanase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	479	Total	C	N	O	S	0	1	0
			3714	2343	646	712	13			
1	B	479	Total	C	N	O	S	0	1	0
			3714	2343	646	712	13			
1	C	479	Total	C	N	O	S	0	3	0
			3727	2353	649	712	13			
1	D	479	Total	C	N	O	S	0	2	0
			3717	2345	646	713	13			

- Molecule 2 is an oligosaccharide called alpha-D-arabinofuranose-(1-5)-alpha-D-arabinofuranose-(1-5)-alpha-D-arabinofuranose-(1-5)-alpha-D-arabinofuranose.



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace
2	E	4	Total	C	O	0	0	0
			37	20	17			
2	I	4	Total	C	O	0	0	0
			37	20	17			

- Molecule 3 is an oligosaccharide called alpha-D-arabinofuranose-(1-5)-alpha-D-arabinofuranose-(1-5)-alpha-D-arabinofuranose.



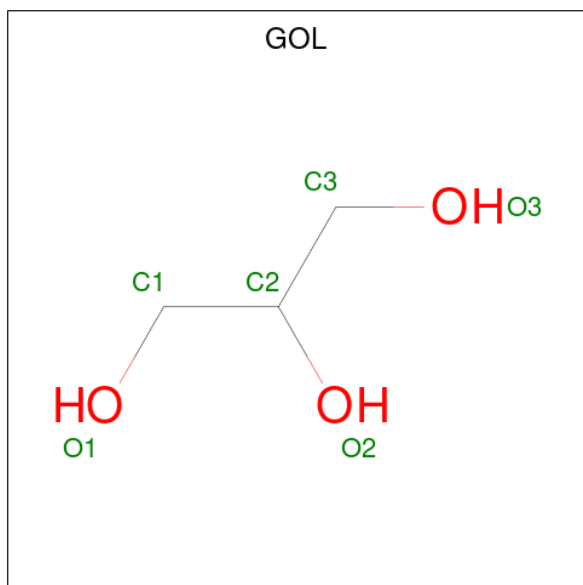
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace
3	F	3	Total	C	O	0	0	0
			27	15	12			

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace
3	G	3	Total	C	O	0	0	0
			27	15	12			
3	H	3	Total	C	O	0	0	0
			27	15	12			
3	J	3	Total	C	O	0	0	0
			27	15	12			

- Molecule 4 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



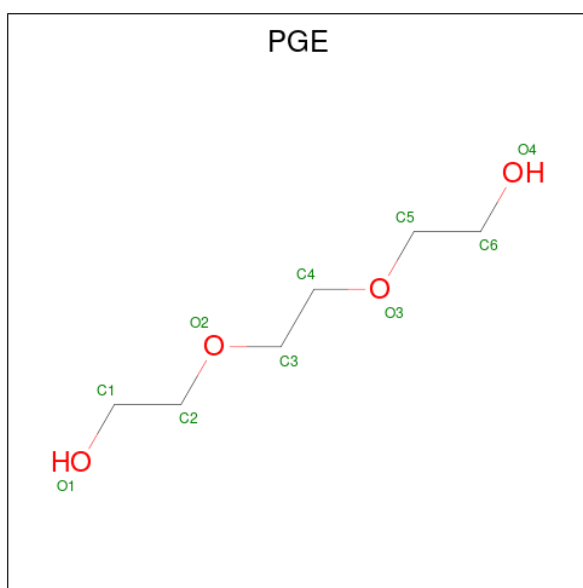
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			6	3	3		
4	A	1	Total	C	O	0	0
			6	3	3		
4	B	1	Total	C	O	0	0
			6	3	3		
4	B	1	Total	C	O	0	0
			6	3	3		
4	B	1	Total	C	O	0	0
			6	3	3		
4	B	1	Total	C	O	0	0
			6	3	3		
4	C	1	Total	C	O	0	0
			6	3	3		

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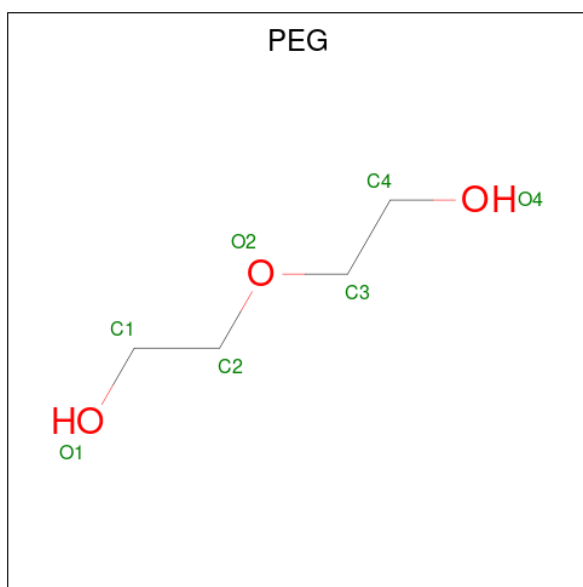
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	C	1	Total	C	O	0	0
			6	3	3		
4	C	1	Total	C	O	0	0
			6	3	3		
4	D	1	Total	C	O	0	0
			6	3	3		
4	D	1	Total	C	O	0	0
			6	3	3		
4	D	1	Total	C	O	0	0
			6	3	3		

- Molecule 5 is TRIETHYLENE GLYCOL (CCD ID: PGE) (formula: $C_6H_{14}O_4$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			10	6	4		
5	A	1	Total	C	O	0	0
			10	6	4		
5	B	1	Total	C	O	0	0
			10	6	4		
5	C	1	Total	C	O	0	0
			10	6	4		
5	D	1	Total	C	O	0	0
			10	6	4		

- Molecule 6 is DI(HYDROXYETHYL)ETHER (CCD ID: PEG) (formula: $C_4H_{10}O_3$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	A	1	Total	C	O	0	0
			7	4	3		
6	B	1	Total	C	O	0	0
			7	4	3		
6	C	1	Total	C	O	0	0
			7	4	3		
6	C	1	Total	C	O	0	0
			7	4	3		
6	D	1	Total	C	O	0	0
			7	4	3		
6	D	1	Total	C	O	0	0
			7	4	3		

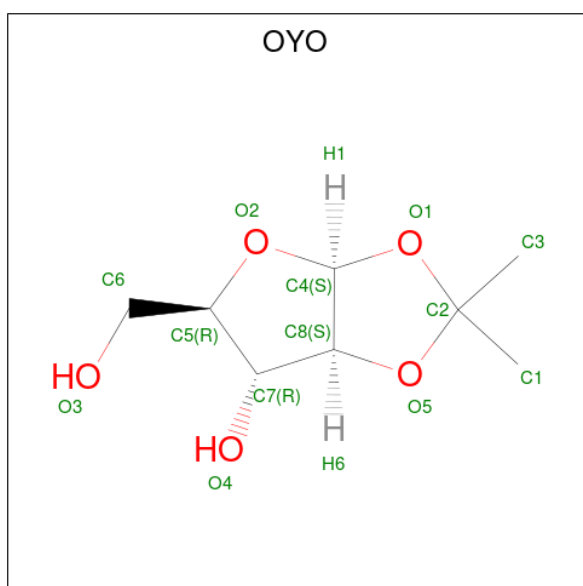
- Molecule 7 is SODIUM ION (CCD ID: NA) (formula: Na).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	1	Total	Na	0	0
			1	1		
7	B	1	Total	Na	0	0
			1	1		
7	C	1	Total	Na	0	0
			1	1		
7	D	1	Total	Na	0	0
			1	1		

- Molecule 8 is CALCIUM ION (CCD ID: CA) (formula: Ca).

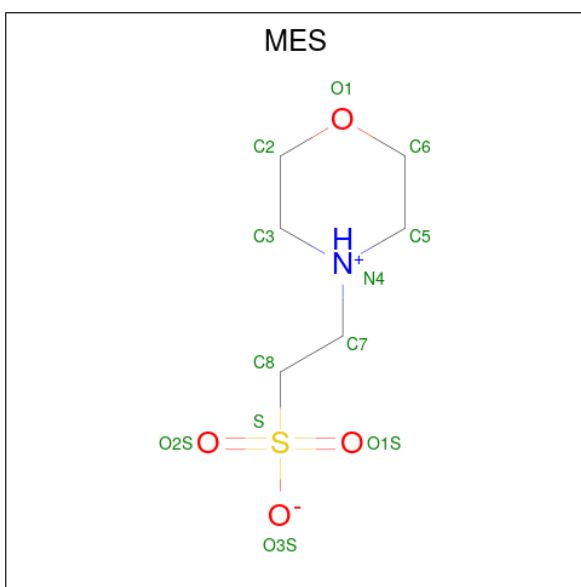
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
8	A	1	Total Ca 1 1	0	0
8	B	1	Total Ca 1 1	0	0
8	C	1	Total Ca 1 1	0	0
8	D	1	Total Ca 1 1	0	0

- Molecule 9 is (3 {a} {S},5 {R},6 {R},6 {a} {S})-5-(hydroxymethyl)-2,2-dimethyl-3 {a},5,6,6 {a}-tetrahydrofuro[2,3-d][1,3]dioxol-6-ol (CCD ID: OYO) (formula: C₈H₁₄O₅) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
9	A	1	Total C O 13 8 5	0	0
9	B	1	Total C O 13 8 5	0	0
9	C	1	Total C O 13 8 5	0	0
9	D	1	Total C O 13 8 5	0	0

- Molecule 10 is 2-(N-MORPHOLINO)-ETHANESULFONIC ACID (CCD ID: MES) (formula: C₆H₁₃NO₄S).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
10	B	1	Total	C	N	O	S	0	0
			12	6	1	4	1		

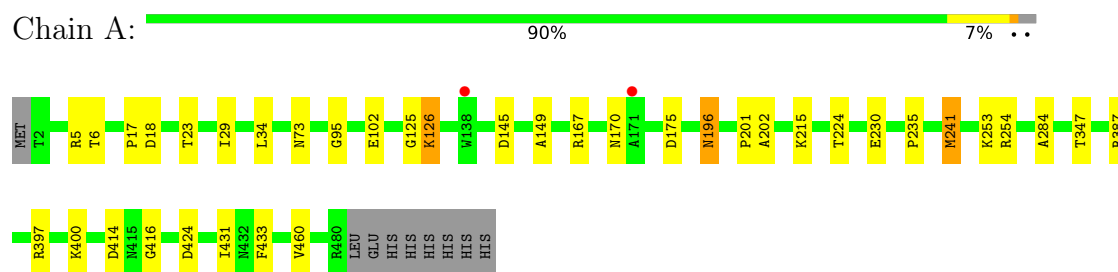
- Molecule 11 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
11	A	354	Total	O	0	0
			354	354		
11	B	330	Total	O	0	0
			330	330		
11	C	313	Total	O	0	0
			313	313		
11	D	317	Total	O	0	0
			317	317		

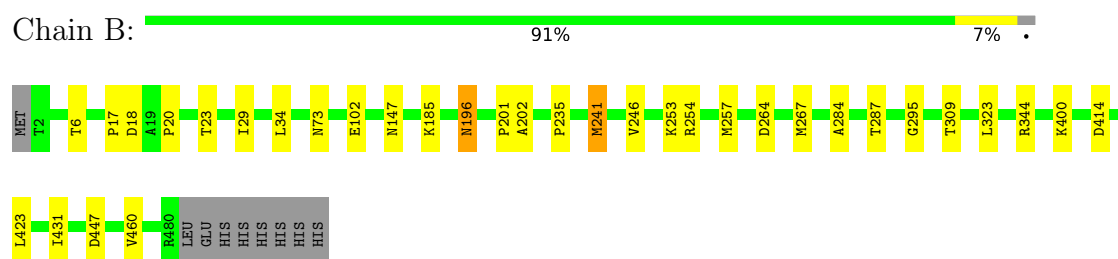
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

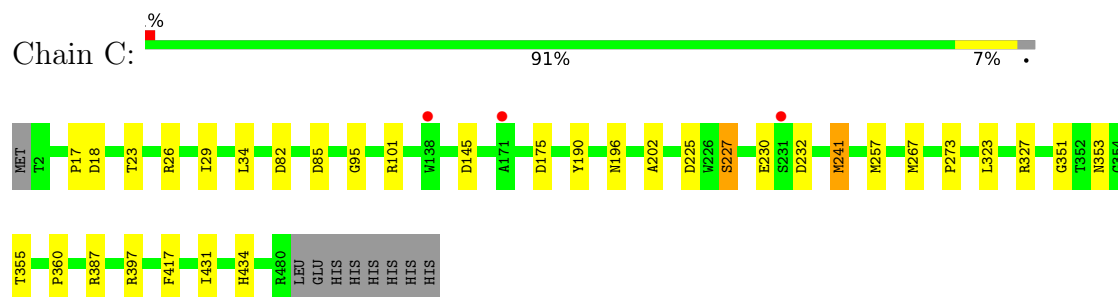
• Molecule 1: endo-alpha-D-arabinanase



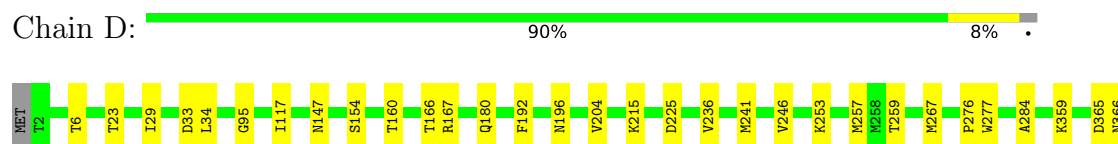
• Molecule 1: endo-alpha-D-arabinanase



• Molecule 1: endo-alpha-D-arabinanase



• Molecule 1: endo-alpha-D-arabinanase





- Molecule 2: alpha-D-arabinofuranose-(1-5)-alpha-D-arabinofuranose-(1-5)-alpha-D-arabinofuranose-(1-5)-alpha-D-arabinofuranose

Chain E: 100%



- Molecule 2: alpha-D-arabinofuranose-(1-5)-alpha-D-arabinofuranose-(1-5)-alpha-D-arabinofuranose-(1-5)-alpha-D-arabinofuranose

Chain I: 25% 75%



- Molecule 3: alpha-D-arabinofuranose-(1-5)-alpha-D-arabinofuranose-(1-5)-alpha-D-arabinofuranose

Chain F: 33% 67%



- Molecule 3: alpha-D-arabinofuranose-(1-5)-alpha-D-arabinofuranose-(1-5)-alpha-D-arabinofuranose

Chain G: 33% 67%



- Molecule 3: alpha-D-arabinofuranose-(1-5)-alpha-D-arabinofuranose-(1-5)-alpha-D-arabinofuranose

Chain H: 67% 33%



- Molecule 3: alpha-D-arabinofuranose-(1-5)-alpha-D-arabinofuranose-(1-5)-alpha-D-arabinofuranose

Chain J: 100%



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	114.71Å 137.73Å 148.59Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	46.65 – 1.80 46.65 – 1.80	Depositor EDS
% Data completeness (in resolution range)	100.0 (46.65-1.80) 100.0 (46.65-1.80)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.22 (at 1.79Å)	Xtriage
Refinement program	REFMAC 5.8.0403	Depositor
R, R_{free}	0.159 , 0.190 (Not available) , 0.194	Depositor DCC
R_{free} test set	10991 reflections (5.06%)	wwPDB-VP
Wilson B-factor (Å ²)	20.9	Xtriage
Anisotropy	0.028	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.39 , 38.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	16610	wwPDB-VP
Average B, all atoms (Å ²)	24.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.55% 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: CA, NA, OYO, PGE, PEG, BXY, MES, GOL

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	0.67	0/3826	1.05	5/5218 (0.1%)
1	B	0.65	0/3826	1.01	3/5218 (0.1%)
1	C	0.64	0/3845	1.01	1/5243 (0.0%)
1	D	0.62	0/3832	1.01	3/5226 (0.1%)
All	All	0.65	0/15329	1.02	12/20905 (0.1%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	3
1	B	0	1
1	C	0	3
1	D	0	2
All	All	0	9

There are no bond length outliers.

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	414	ASP	CA-CB-CG	6.25	118.85	112.60
1	A	196	ASN	CA-CB-CG	6.04	118.64	112.60
1	D	33	ASP	CA-CB-CG	5.98	118.58	112.60
1	D	166	THR	CA-CB-OG1	-5.96	100.65	109.60
1	A	424	ASP	CB-CA-C	-5.94	98.14	109.95
1	B	196	ASN	CA-CB-CG	5.79	118.39	112.60
1	D	117	ILE	CA-C-O	5.51	122.93	119.51
1	B	264	ASP	CA-CB-CG	5.39	117.99	112.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	414	ASP	CA-CB-CG	5.36	117.96	112.60
1	A	170	ASN	CA-C-N	5.08	127.39	120.54
1	A	170	ASN	C-N-CA	5.08	127.39	120.54
1	C	434	HIS	CA-CB-CG	-5.07	108.73	113.80

There are no chirality outliers.

All (9) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	254	ARG	Sidechain
1	A	387	ARG	Sidechain
1	A	5	ARG	Sidechain
1	B	254	ARG	Sidechain
1	C	26	ARG	Sidechain
1	C	327	ARG	Sidechain
1	C	397	ARG	Sidechain
1	D	167	ARG	Sidechain
1	D	451	ARG	Sidechain

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	3714	0	3496	16	0
1	B	3714	0	3496	18	0
1	C	3727	0	3520	18	0
1	D	3717	0	3501	25	0
2	E	37	0	31	0	0
2	I	37	0	31	0	0
3	F	27	0	22	0	0
3	G	27	0	22	0	0
3	H	27	0	22	0	0
3	J	27	0	22	0	0
4	A	12	0	16	1	0
4	B	30	0	40	2	0
4	C	18	0	24	0	0
4	D	18	0	24	3	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	A	20	0	28	2	0
5	B	10	0	14	1	0
5	C	10	0	14	1	0
5	D	10	0	14	1	0
6	A	7	0	10	0	0
6	B	7	0	10	1	0
6	C	14	0	20	0	0
6	D	14	0	20	7	0
7	A	1	0	0	0	0
7	B	1	0	0	0	0
7	C	1	0	0	0	0
7	D	1	0	0	0	0
8	A	1	0	0	0	0
8	B	1	0	0	0	0
8	C	1	0	0	0	0
8	D	1	0	0	0	0
9	A	13	0	0	0	0
9	B	13	0	0	0	0
9	C	13	0	0	0	0
9	D	13	0	0	0	0
10	B	12	0	13	0	0
11	A	354	0	0	2	0
11	B	330	0	0	1	0
11	C	313	0	0	2	0
11	D	317	0	0	4	0
All	All	16610	0	14410	80	0

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

All (80) 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:C:225:ASP:HB2	11:C:826:HOH:O	1.76	0.84
1:D:276:PRO:HA	6:D:505:PEG:H22	1.61	0.83
1:D:277:TRP:H	6:D:505:PEG:H21	1.48	0.79
1:D:277:TRP:H	6:D:505:PEG:C2	1.97	0.78
1:D:225:ASP:HB2	11:D:827:HOH:O	1.83	0.78
4:A:501:GOL:H2	11:A:889:HOH:O	1.89	0.70
1:D:400:LYS:HG2	1:D:447:ASP:OD1	1.95	0.66
1:D:246:VAL:O	4:D:501:GOL:H12	1.96	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:145:ASP:HB3	5:C:502:PGE:H5	1.77	0.64
1:B:400:LYS:HG2	1:B:447:ASP:OD1	1.98	0.63
1:D:463:ASN:HB2	11:D:835:HOH:O	2.01	0.61
1:D:359:LYS:NZ	11:D:602:HOH:O	2.33	0.61
1:D:241:MET:HG2	1:D:259:THR:HG21	1.84	0.58
1:C:227:SER:OG	1:C:232:ASP:OD2	2.19	0.55
1:B:202:ALA:HB2	1:B:241:MET:HE2	1.87	0.55
1:B:6:THR:HG23	1:B:284:ALA:HA	1.89	0.55
1:A:431:ILE:HD12	1:A:431:ILE:C	2.32	0.55
1:C:82:ASP:OD2	1:C:85:ASP:OD2	2.25	0.54
1:A:126:LYS:HE3	1:A:126:LYS:N	2.24	0.53
1:A:145:ASP:HB3	5:A:502:PGE:H4	1.93	0.51
4:D:503:GOL:H12	6:D:504:PEG:H42	1.93	0.50
1:B:447:ASP:HB3	11:B:706:HOH:O	2.11	0.50
1:B:17:PRO:O	1:B:18[B]:ASP:HB2	2.13	0.49
1:A:202:ALA:HB2	1:A:241:MET:HE2	1.95	0.48
1:C:431:ILE:C	1:C:431:ILE:HD12	2.38	0.48
4:B:503:GOL:H32	1:D:366:ASN:H	1.78	0.48
1:A:253:LYS:HD2	5:A:503:PGE:H52	1.94	0.48
1:B:201:PRO:HB2	1:B:235:PRO:HB3	1.96	0.47
1:B:287:THR:CG2	1:B:344:ARG:HG2	2.43	0.47
1:B:73:ASN:O	1:B:102:GLU:HA	2.16	0.46
1:C:82:ASP:OD2	1:C:85:ASP:CG	2.59	0.46
1:A:149:ALA:O	1:A:167:ARG:HA	2.16	0.46
1:B:287:THR:HG22	1:B:344:ARG:HG2	1.98	0.46
1:D:381:ARG:HH22	6:D:504:PEG:H21	1.80	0.46
1:D:147:ASN:OD1	5:D:502:PGE:H1	2.16	0.45
1:A:125:GLY:C	1:A:126:LYS:HE3	2.41	0.45
1:C:101[A]:ARG:NH1	11:C:607:HOH:O	2.45	0.45
1:D:154:SER:HB2	1:D:160:THR:O	2.17	0.45
1:A:6:THR:HG23	1:A:284:ALA:HA	1.99	0.45
1:D:365:ASP:O	1:D:376:HIS:HA	2.17	0.45
1:A:400:LYS:NZ	11:A:614:HOH:O	2.50	0.45
1:A:201:PRO:HB2	1:A:235:PRO:HB3	1.98	0.44
1:D:6:THR:HG23	1:D:284:ALA:HA	2.00	0.44
1:C:241:MET:HE3	1:C:241:MET:HB3	1.81	0.44
1:D:241:MET:HE2	1:D:241:MET:HB3	1.89	0.44
1:C:17:PRO:O	1:C:18[A]:ASP:HB2	2.16	0.44
1:D:257:MET:O	1:D:267:MET:HA	2.17	0.44
1:D:253:LYS:NZ	11:D:615:HOH:O	2.50	0.44
1:B:185:LYS:NZ	6:B:506:PEG:H32	2.32	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:D:503:GOL:H12	6:D:504:PEG:C4	2.48	0.43
1:A:175:ASP:OD1	1:A:230:GLU:OE2	2.36	0.43
1:B:241:MET:HE3	1:B:241:MET:HB3	1.74	0.43
1:D:431:ILE:C	1:D:431:ILE:HD12	2.42	0.43
1:A:416:GLY:HA3	1:A:433:PHE:CE1	2.54	0.43
1:C:323:LEU:HD12	1:C:323:LEU:C	2.44	0.43
1:A:241:MET:HB3	1:A:241:MET:HE3	1.57	0.42
1:D:241:MET:HE3	1:D:259:THR:HG21	2.01	0.42
1:C:190:TYR:CE1	1:C:273:PRO:HG2	2.54	0.42
1:B:257:MET:O	1:B:267:MET:HA	2.20	0.42
1:B:431:ILE:C	1:B:431:ILE:HD12	2.45	0.42
1:B:323:LEU:C	1:B:323:LEU:HD12	2.45	0.42
1:D:277:TRP:N	6:D:505:PEG:H21	2.26	0.42
1:A:23:THR:HB	1:A:29:ILE:HB	2.02	0.41
1:A:73:ASN:O	1:A:102:GLU:HA	2.20	0.41
1:B:147:ASN:HA	5:B:502:PGE:H5	2.02	0.41
1:D:448:ALA:O	1:D:451:ARG:NH1	2.45	0.41
1:D:23:THR:HB	1:D:29:ILE:HB	2.01	0.41
1:D:180:GLN:HA	1:D:192:PHE:O	2.20	0.41
1:A:17:PRO:O	1:A:18[A]:ASP:HB2	2.19	0.41
1:C:175:ASP:OD1	1:C:230:GLU:OE2	2.38	0.41
1:C:202:ALA:HB2	1:C:241:MET:HE2	2.01	0.41
1:C:257:MET:O	1:C:267:MET:HA	2.21	0.41
1:B:246:VAL:O	4:B:501:GOL:C1	2.69	0.41
1:C:23:THR:HB	1:C:29:ILE:HB	2.03	0.41
1:D:204:VAL:HG12	1:D:236:VAL:HG11	2.03	0.41
1:B:295:GLY:HA2	1:B:309:THR:O	2.20	0.41
1:B:23:THR:HB	1:B:29:ILE:HB	2.03	0.40
1:C:351:GLY:HA2	1:C:355:THR:O	2.22	0.40
1:C:360:PRO:O	1:C:387:ARG:HD2	2.21	0.40
1:C:417:PHE:HA	1:C:431:ILE:O	2.22	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	478/488 (98%)	462 (97%)	14 (3%)	2 (0%)	30	19
1	B	478/488 (98%)	459 (96%)	18 (4%)	1 (0%)	43	31
1	C	480/488 (98%)	465 (97%)	13 (3%)	2 (0%)	30	19
1	D	479/488 (98%)	459 (96%)	18 (4%)	2 (0%)	30	19
All	All	1915/1952 (98%)	1845 (96%)	63 (3%)	7 (0%)	30	19

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	34	LEU
1	C	34	LEU
1	D	34	LEU
1	B	34	LEU
1	D	95	GLY
1	A	95	GLY
1	C	95	GLY

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	384/392 (98%)	376 (98%)	8 (2%)	47	36
1	B	384/392 (98%)	378 (98%)	6 (2%)	55	47
1	C	386/392 (98%)	382 (99%)	4 (1%)	68	64
1	D	385/392 (98%)	382 (99%)	3 (1%)	73	70
All	All	1539/1568 (98%)	1518 (99%)	21 (1%)	59	52

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

Mol	Chain	Res	Type
1	A	126	LYS

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Mol	Chain	Res	Type
1	A	196	ASN
1	A	215	LYS
1	A	224	THR
1	A	241	MET
1	A	347	THR
1	A	397	ARG
1	A	460	VAL
1	B	20	PRO
1	B	196	ASN
1	B	241	MET
1	B	253	LYS
1	B	423	LEU
1	B	460	VAL
1	C	196	ASN
1	C	227	SER
1	C	241	MET
1	C	353	ASN
1	D	196	ASN
1	D	215	LYS
1	D	480	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	147	ASN
1	A	353	ASN
1	B	330	ASN
1	B	353	ASN
1	B	383	ASN
1	B	415	ASN
1	B	432	ASN
1	C	330	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 ⓘ

20 monosaccharides are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	BXY	E	1	2	10,10,10	0.39	0	13,14,14	1.41	2 (15%)
2	BXY	E	2	2	9,9,10	0.68	0	11,12,14	1.70	3 (27%)
2	BXY	E	3	2	9,9,10	0.59	0	11,12,14	1.48	2 (18%)
2	BXY	E	4	2	9,9,10	0.84	0	11,12,14	1.11	1 (9%)
3	BXY	F	1	9,3	9,9,10	0.46	0	11,12,14	0.69	0
3	BXY	F	2	3	9,9,10	1.01	1 (11%)	11,12,14	1.26	2 (18%)
3	BXY	F	3	3	9,9,10	0.83	1 (11%)	11,12,14	1.55	3 (27%)
3	BXY	G	1	9,3	9,9,10	0.58	0	11,12,14	1.12	1 (9%)
3	BXY	G	2	3	9,9,10	0.69	0	11,12,14	1.14	0
3	BXY	G	3	3	9,9,10	0.69	0	11,12,14	1.32	2 (18%)
3	BXY	H	1	9,3	9,9,10	0.78	0	11,12,14	0.67	0
3	BXY	H	2	3	9,9,10	0.61	0	11,12,14	0.73	0
3	BXY	H	3	3	9,9,10	0.59	0	11,12,14	1.00	1 (9%)
2	BXY	I	1	2	10,10,10	0.45	0	13,14,14	0.97	1 (7%)
2	BXY	I	2	2	9,9,10	0.62	0	11,12,14	1.99	3 (27%)
2	BXY	I	3	2	9,9,10	0.55	0	11,12,14	1.03	2 (18%)
2	BXY	I	4	2	9,9,10	0.65	0	11,12,14	0.92	0
3	BXY	J	1	9,3	9,9,10	0.98	1 (11%)	11,12,14	0.82	0
3	BXY	J	2	3	9,9,10	0.72	0	11,12,14	1.12	1 (9%)
3	BXY	J	3	3	9,9,10	0.87	0	11,12,14	1.59	2 (18%)

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	BXY	E	1	2	-	0/2/18/18	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	BXY	E	2	2	-	0/2/15/18	0/1/1/1
2	BXY	E	3	2	-	0/2/15/18	0/1/1/1
2	BXY	E	4	2	-	0/2/15/18	0/1/1/1
3	BXY	F	1	9,3	-	0/2/15/18	0/1/1/1
3	BXY	F	2	3	-	0/2/15/18	0/1/1/1
3	BXY	F	3	3	-	2/2/15/18	0/1/1/1
3	BXY	G	1	9,3	-	0/2/15/18	0/1/1/1
3	BXY	G	2	3	-	0/2/15/18	0/1/1/1
3	BXY	G	3	3	-	0/2/15/18	0/1/1/1
3	BXY	H	1	9,3	-	0/2/15/18	0/1/1/1
3	BXY	H	2	3	-	0/2/15/18	0/1/1/1
3	BXY	H	3	3	-	0/2/15/18	0/1/1/1
2	BXY	I	1	2	-	0/2/18/18	0/1/1/1
2	BXY	I	2	2	-	0/2/15/18	0/1/1/1
2	BXY	I	3	2	-	0/2/15/18	0/1/1/1
2	BXY	I	4	2	-	0/2/15/18	0/1/1/1
3	BXY	J	1	9,3	-	0/2/15/18	0/1/1/1
3	BXY	J	2	3	-	0/2/15/18	0/1/1/1
3	BXY	J	3	3	-	0/2/15/18	0/1/1/1

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	F	2	BXY	C1-C2	2.36	1.55	1.51
3	J	1	BXY	O4-C1	2.18	1.48	1.43
3	F	3	BXY	C1-C2	2.16	1.55	1.51

All (26) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	J	3	BXY	O2-C2-C3	4.10	119.50	111.43
2	E	1	BXY	C1-C2-C3	3.73	106.88	102.29
2	E	3	BXY	C1-C2-C3	3.49	107.22	101.63
2	I	2	BXY	O2-C2-C3	-3.38	104.76	111.43
3	G	3	BXY	O4-C4-C3	3.29	107.68	104.63
2	I	2	BXY	O4-C4-C3	3.27	107.65	104.63
2	I	2	BXY	C1-C2-C3	3.27	106.85	101.63
3	F	3	BXY	O3-C3-C2	-3.11	104.61	111.97
2	E	2	BXY	O4-C4-C3	2.83	107.25	104.63
3	J	2	BXY	O2-C2-C3	2.74	116.82	111.43
2	E	4	BXY	C1-C2-C3	2.71	105.96	101.63
3	F	3	BXY	C5-C4-C3	2.65	121.36	115.10

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	J	3	BXY	O4-C4-C3	2.59	107.02	104.63
3	F	2	BXY	O4-C4-C3	-2.58	102.25	104.63
2	E	2	BXY	C1-C2-C3	2.52	105.66	101.63
2	E	1	BXY	C5-C4-C3	-2.50	109.20	115.10
3	G	1	BXY	O2-C2-C3	2.49	116.34	111.43
2	E	3	BXY	O4-C4-C3	2.36	106.82	104.63
2	I	1	BXY	C1-C2-C3	2.36	105.19	102.29
3	G	3	BXY	O2-C2-C3	2.29	115.94	111.43
2	I	3	BXY	O4-C4-C3	2.27	106.73	104.63
3	F	3	BXY	O4-C4-C3	2.15	106.62	104.63
2	E	2	BXY	O3-C3-C2	2.11	116.96	111.97
3	H	3	BXY	O3-C3-C2	-2.07	107.07	111.97
3	F	2	BXY	C1-C2-C3	2.06	104.92	101.63
2	I	3	BXY	C1-C2-C3	2.05	104.90	101.63

There are no chirality outliers.

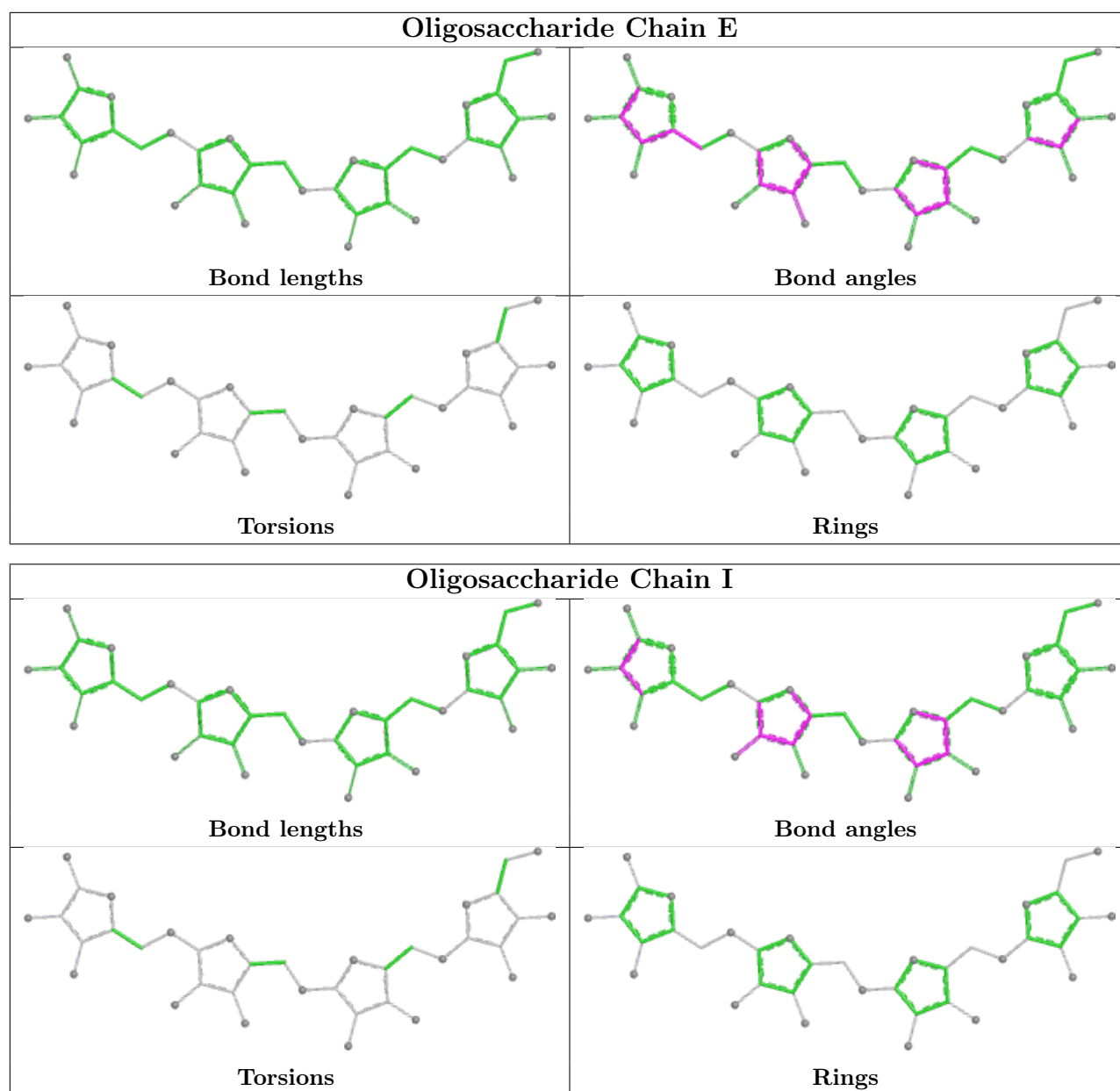
All (2) torsion outliers are listed below:

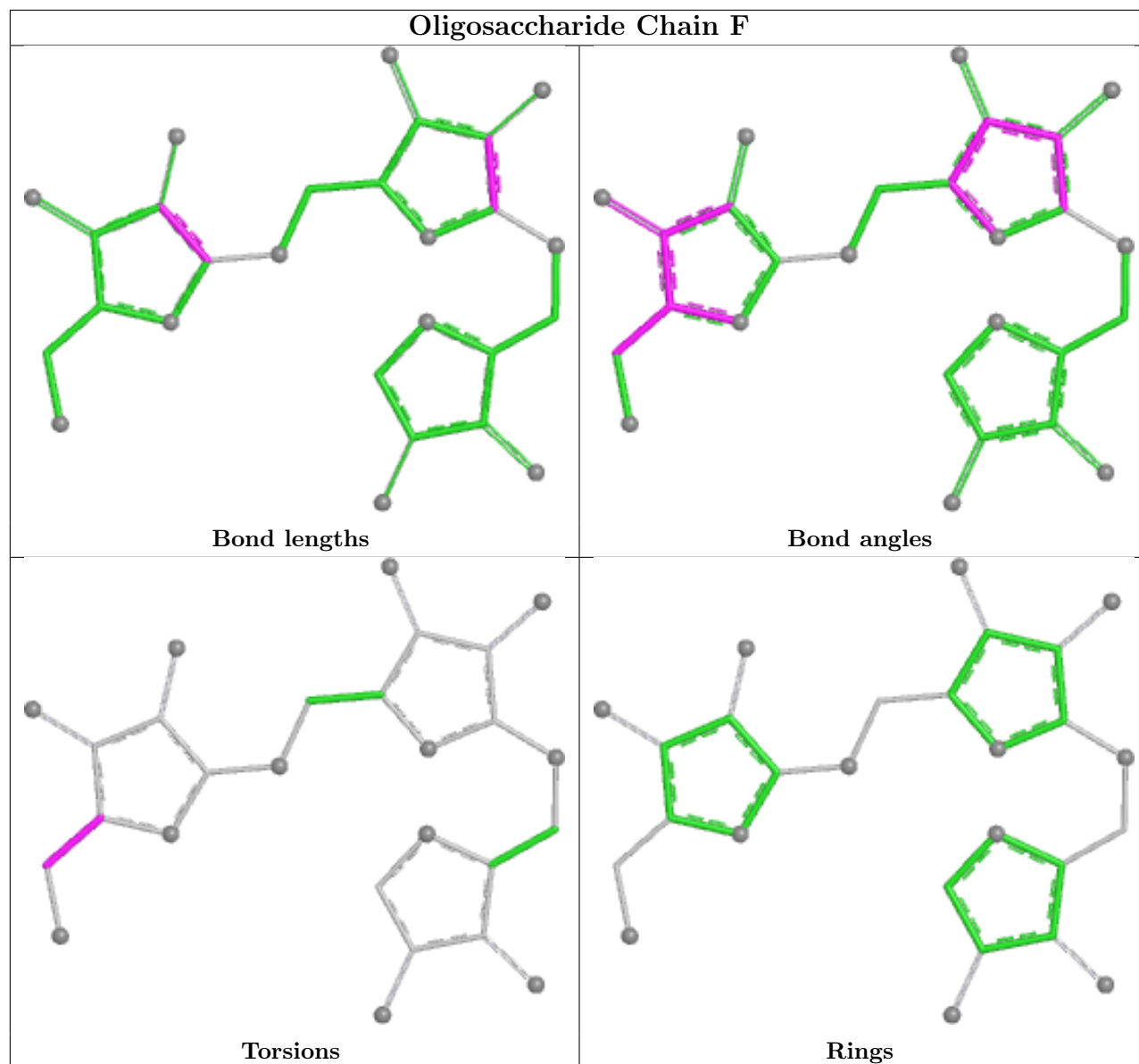
Mol	Chain	Res	Type	Atoms
3	F	3	BXY	O4-C4-C5-O5
3	F	3	BXY	C3-C4-C5-O5

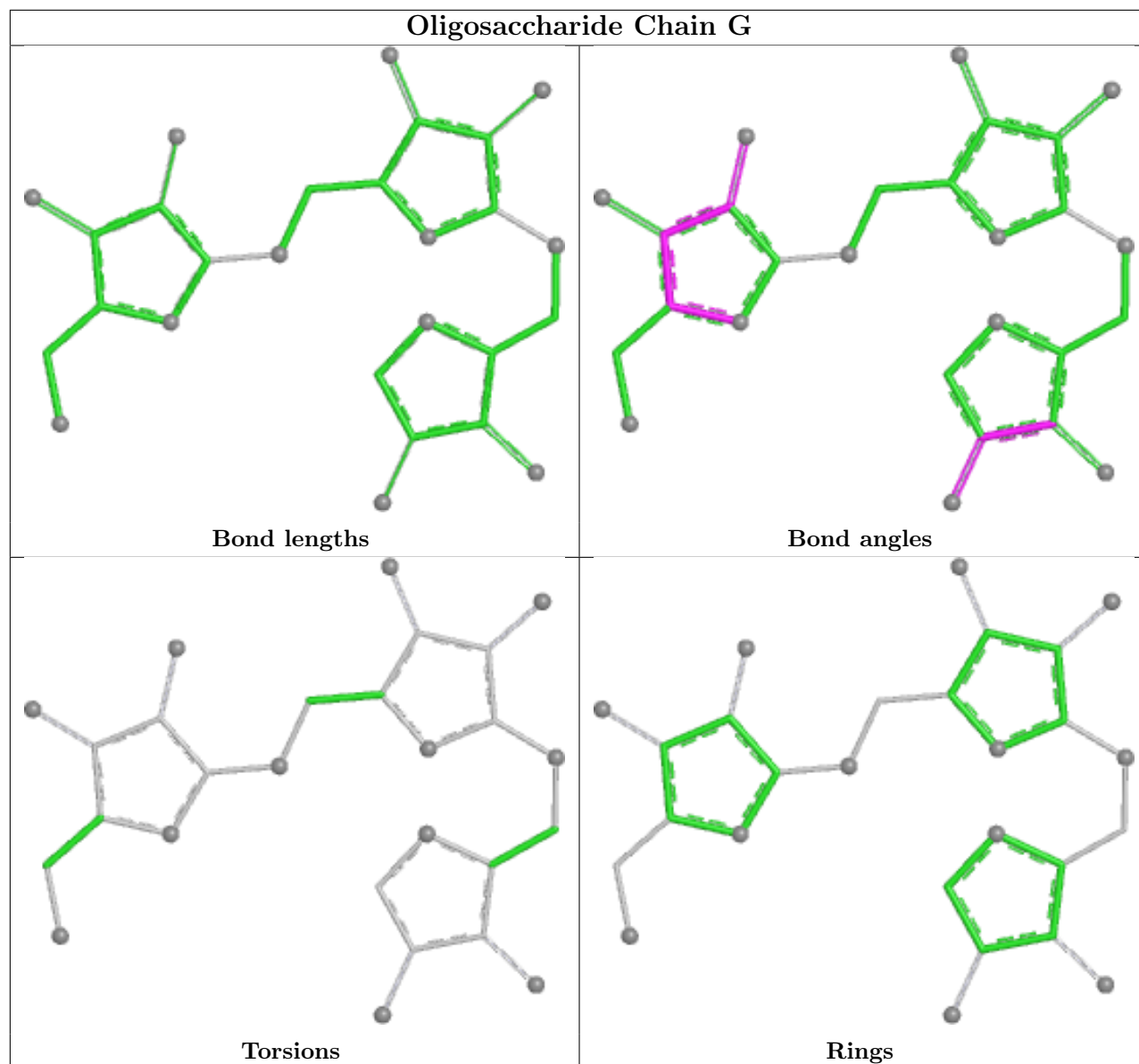
There are no ring outliers.

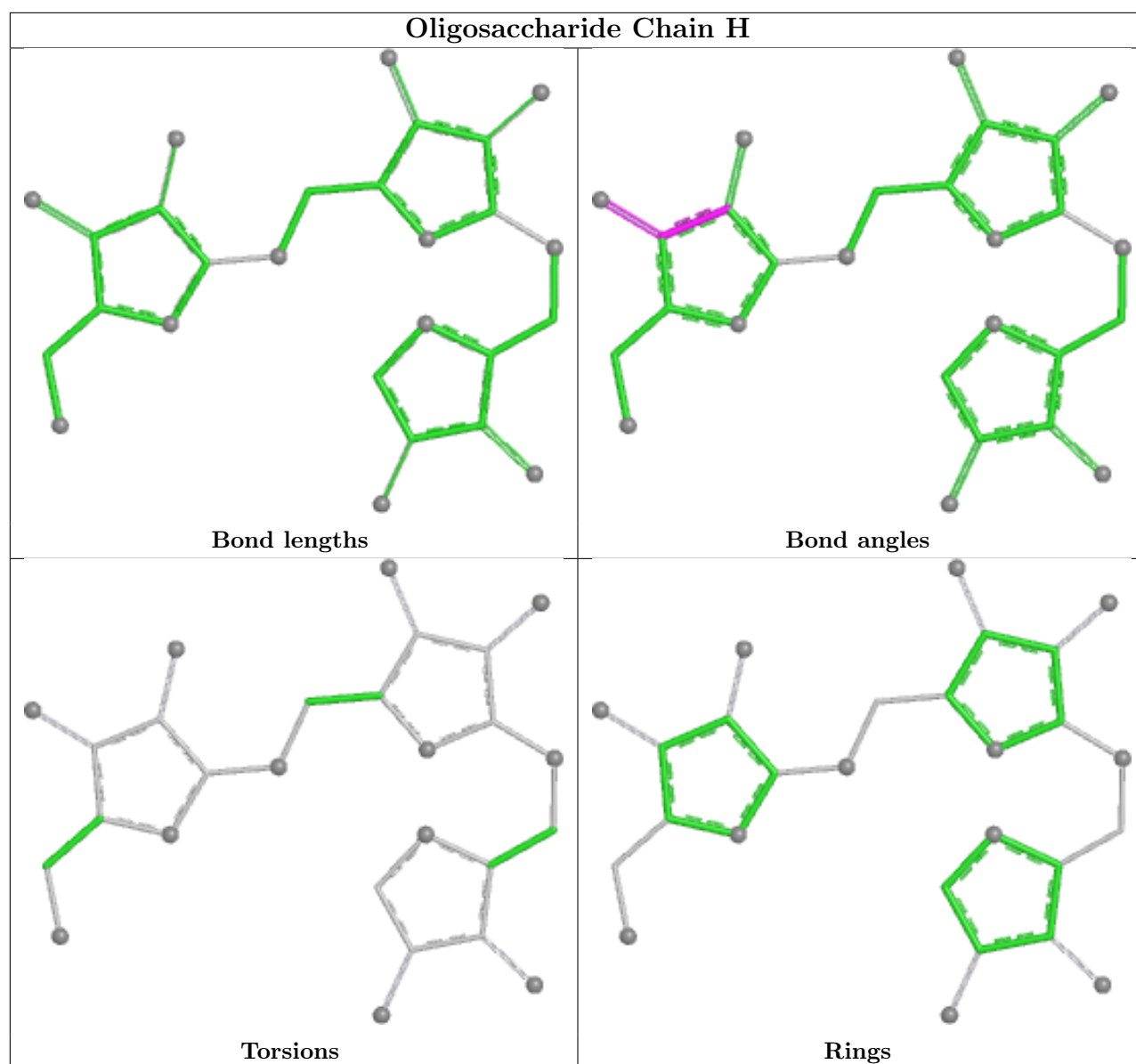
No monomer is involved in short contacts.

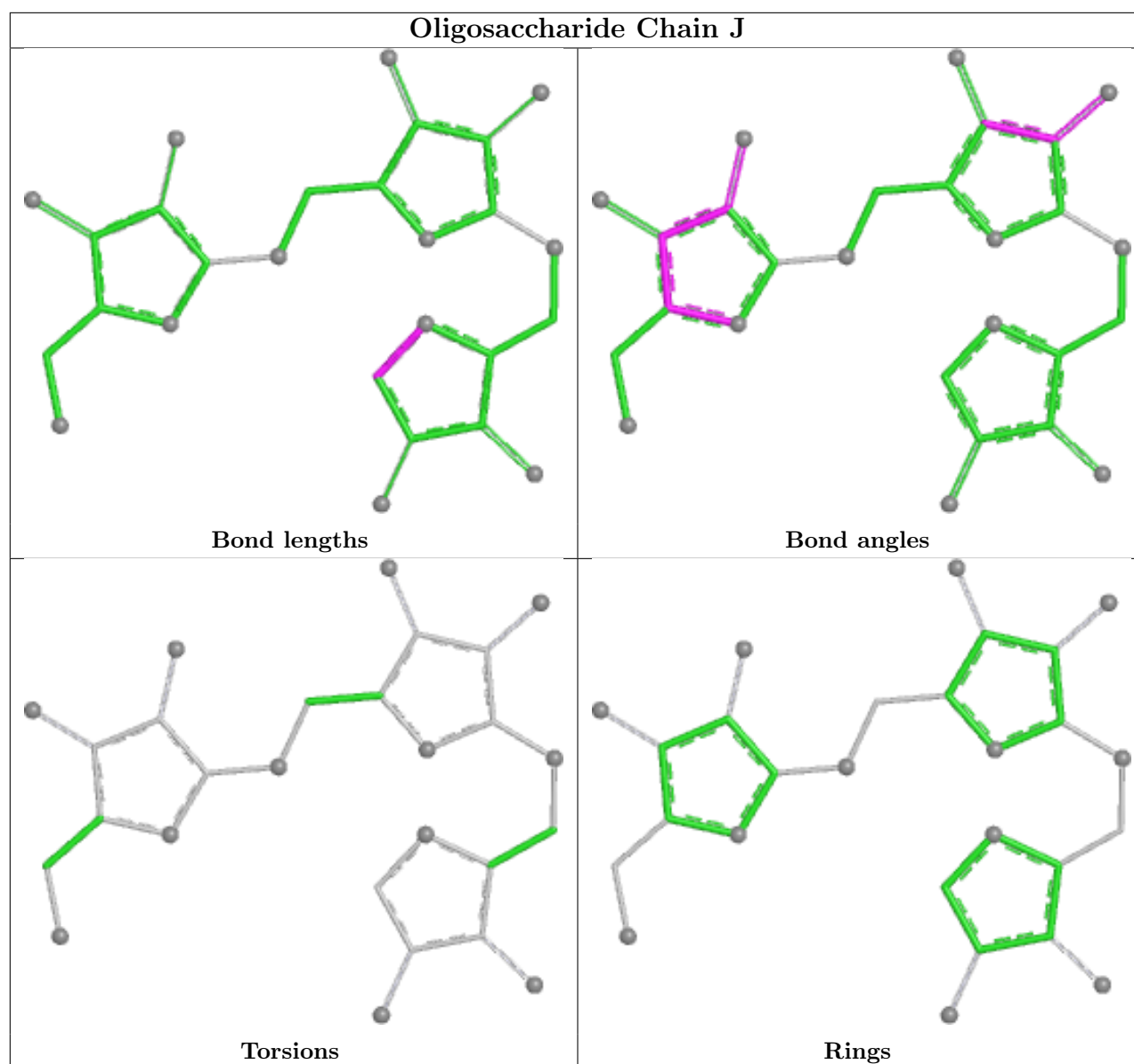
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.











5.6 Ligand geometry [i](#)

Of 37 ligands modelled in this entry, 8 are monoatomic - leaving 29 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	$\# Z > 2$
10	MES	B	505	-	12,12,12	0.70	0	15,16,16	0.68	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	GOL	B	503	-	5,5,5	0.27	0	5,5,5	0.58	0
6	PEG	D	504	-	6,6,6	0.18	0	5,5,5	0.25	0
4	GOL	C	505	-	5,5,5	0.27	0	5,5,5	0.46	0
6	PEG	B	506	-	6,6,6	0.33	0	5,5,5	0.19	0
9	OYO	D	509	3	14,14,14	0.54	0	22,22,22	1.61	6 (27%)
6	PEG	C	504	-	6,6,6	0.29	0	5,5,5	0.19	0
5	PGE	B	502	-	9,9,9	0.15	0	8,8,8	0.18	0
4	GOL	B	501	-	5,5,5	0.16	0	5,5,5	0.42	0
5	PGE	D	502	-	9,9,9	0.16	0	8,8,8	0.22	0
5	PGE	C	502	-	9,9,9	0.16	0	8,8,8	0.30	0
4	GOL	B	509	-	5,5,5	0.19	0	5,5,5	0.37	0
4	GOL	B	504	-	5,5,5	0.28	0	5,5,5	1.16	0
4	GOL	B	511	-	5,5,5	0.14	0	5,5,5	0.44	0
6	PEG	D	505	-	6,6,6	0.21	0	5,5,5	0.33	0
4	GOL	C	508	-	5,5,5	0.14	0	5,5,5	0.34	0
4	GOL	D	501	-	5,5,5	0.13	0	5,5,5	0.39	0
9	OYO	C	509	3	14,14,14	0.66	0	22,22,22	1.23	2 (9%)
9	OYO	B	510	3	14,14,14	0.39	0	22,22,22	1.49	4 (18%)
9	OYO	A	508	3	14,14,14	0.44	0	22,22,22	1.14	3 (13%)
5	PGE	A	502	-	9,9,9	0.26	0	8,8,8	0.29	0
4	GOL	A	501	-	5,5,5	0.15	0	5,5,5	0.44	0
4	GOL	D	503	-	5,5,5	0.18	0	5,5,5	0.50	0
4	GOL	A	507	-	5,5,5	0.28	0	5,5,5	0.32	0
6	PEG	A	504	-	6,6,6	0.16	0	5,5,5	0.22	0
4	GOL	D	508	-	5,5,5	0.12	0	5,5,5	0.54	0
4	GOL	C	501	-	5,5,5	0.15	0	5,5,5	0.25	0
5	PGE	A	503	-	9,9,9	0.24	0	8,8,8	0.45	0
6	PEG	C	503	-	6,6,6	0.16	0	5,5,5	0.39	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
10	MES	B	505	-	-	0/6/14/14	0/1/1/1
4	GOL	B	503	-	-	2/4/4/4	-
6	PEG	D	504	-	-	3/4/4/4	-
4	GOL	C	505	-	-	2/4/4/4	-
6	PEG	B	506	-	-	2/4/4/4	-
9	OYO	D	509	3	-	2/2/28/28	0/2/2/2

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
6	PEG	C	504	-	-	0/4/4/4	-
5	PGE	B	502	-	-	1/7/7/7	-
4	GOL	B	501	-	-	2/4/4/4	-
5	PGE	D	502	-	-	2/7/7/7	-
5	PGE	C	502	-	-	2/7/7/7	-
4	GOL	B	509	-	-	0/4/4/4	-
4	GOL	B	504	-	-	2/4/4/4	-
4	GOL	B	511	-	-	0/4/4/4	-
6	PEG	D	505	-	-	3/4/4/4	-
4	GOL	C	508	-	-	0/4/4/4	-
4	GOL	D	501	-	-	2/4/4/4	-
9	OYO	C	509	3	-	2/2/28/28	0/2/2/2
9	OYO	B	510	3	-	2/2/28/28	0/2/2/2
9	OYO	A	508	3	-	1/2/28/28	0/2/2/2
5	PGE	A	502	-	-	4/7/7/7	-
4	GOL	A	501	-	-	2/4/4/4	-
4	GOL	D	503	-	-	2/4/4/4	-
4	GOL	A	507	-	-	0/4/4/4	-
6	PEG	A	504	-	-	0/4/4/4	-
4	GOL	D	508	-	-	0/4/4/4	-
4	GOL	C	501	-	-	2/4/4/4	-
5	PGE	A	503	-	-	4/7/7/7	-
6	PEG	C	503	-	-	1/4/4/4	-

There are no bond length outliers.

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
9	B	510	OYO	O2-C5-C6	4.28	118.27	109.22
9	D	509	OYO	O2-C5-C7	-3.58	98.05	105.15
9	D	509	OYO	C4-C8-C7	-3.26	100.61	104.10
9	C	509	OYO	O4-C7-C5	-3.11	102.16	111.08
9	D	509	OYO	O4-C7-C5	-3.01	102.43	111.08
9	A	508	OYO	O4-C7-C5	-2.83	102.94	111.08
9	B	510	OYO	O5-C2-O1	-2.80	100.99	105.04
9	B	510	OYO	O2-C5-C7	-2.67	99.85	105.15
9	A	508	OYO	O2-C5-C6	2.57	114.64	109.22
9	C	509	OYO	O2-C5-C6	2.53	114.57	109.22
9	D	509	OYO	O5-C2-O1	-2.40	101.57	105.04

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
9	D	509	OYO	O5-C2-C3	2.09	113.66	109.52
9	D	509	OYO	O2-C5-C6	2.07	113.60	109.22
9	B	510	OYO	O4-C7-C5	-2.06	105.16	111.08
9	A	508	OYO	O5-C2-O1	-2.04	102.09	105.04

There are no chirality outliers.

All (45) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	B	503	GOL	C1-C2-C3-O3
4	D	501	GOL	O1-C1-C2-C3
4	D	503	GOL	C1-C2-C3-O3
4	D	503	GOL	O2-C2-C3-O3
5	A	502	PGE	O2-C3-C4-O3
4	D	501	GOL	O1-C1-C2-O2
5	A	502	PGE	O1-C1-C2-O2
5	A	502	PGE	O3-C5-C6-O4
5	A	503	PGE	O1-C1-C2-O2
5	A	503	PGE	O2-C3-C4-O3
5	C	502	PGE	O3-C5-C6-O4
6	B	506	PEG	O1-C1-C2-O2
9	D	509	OYO	C7-C5-C6-O3
4	A	501	GOL	C1-C2-C3-O3
4	B	501	GOL	C1-C2-C3-O3
4	B	504	GOL	C1-C2-C3-O3
4	C	501	GOL	O1-C1-C2-C3
4	C	505	GOL	O1-C1-C2-C3
5	B	502	PGE	O3-C5-C6-O4
5	D	502	PGE	O3-C5-C6-O4
4	A	501	GOL	O2-C2-C3-O3
4	B	503	GOL	O2-C2-C3-O3
4	B	504	GOL	O2-C2-C3-O3
6	D	505	PEG	O2-C3-C4-O4
9	B	510	OYO	C7-C5-C6-O3
9	D	509	OYO	O2-C5-C6-O3
4	C	505	GOL	O1-C1-C2-O2
9	C	509	OYO	O2-C5-C6-O3
9	B	510	OYO	O2-C5-C6-O3
6	D	504	PEG	O2-C3-C4-O4
4	B	501	GOL	O2-C2-C3-O3
5	A	502	PGE	C6-C5-O3-C4
6	B	506	PEG	O2-C3-C4-O4

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Mol	Chain	Res	Type	Atoms
5	D	502	PGE	C3-C4-O3-C5
9	A	508	OYO	O2-C5-C6-O3
5	A	503	PGE	C4-C3-O2-C2
6	D	505	PEG	O1-C1-C2-O2
9	C	509	OYO	C7-C5-C6-O3
4	C	501	GOL	O1-C1-C2-O2
6	D	504	PEG	O1-C1-C2-O2
5	A	503	PGE	C3-C4-O3-C5
5	C	502	PGE	O2-C3-C4-O3
6	D	504	PEG	C4-C3-O2-C2
6	C	503	PEG	C4-C3-O2-C2
6	D	505	PEG	C4-C3-O2-C2

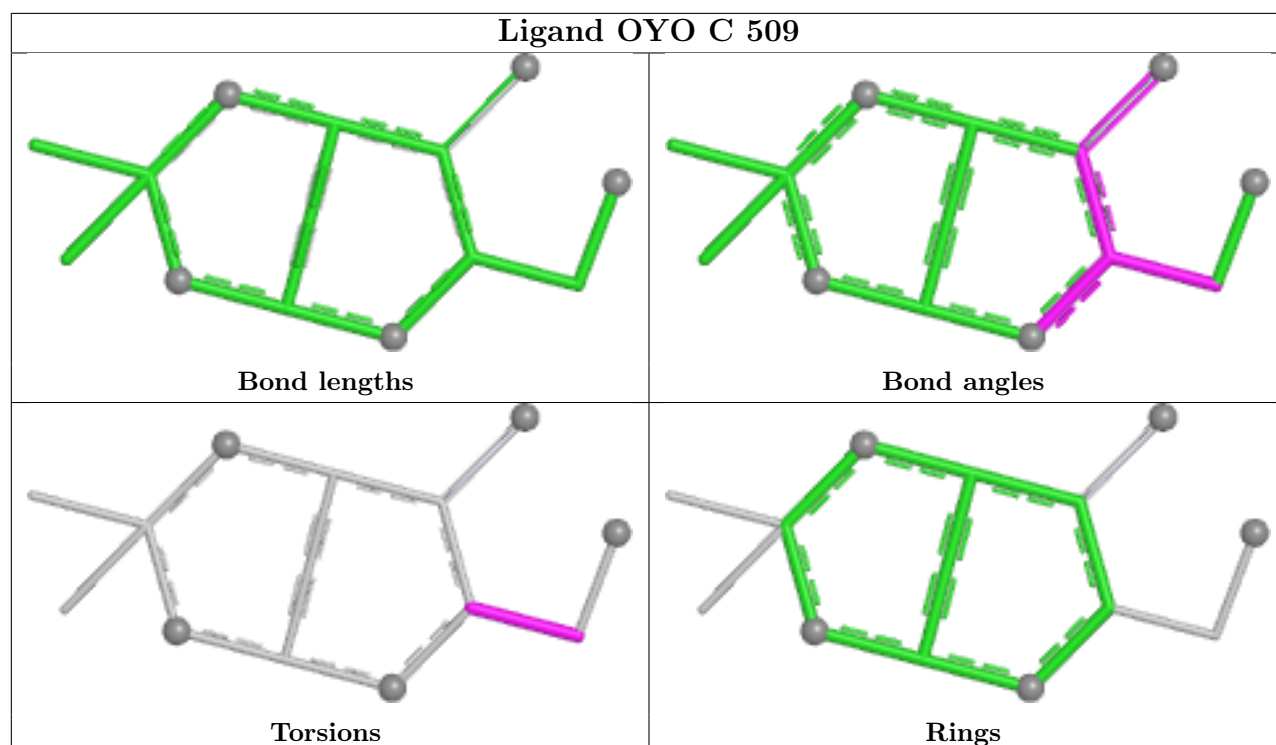
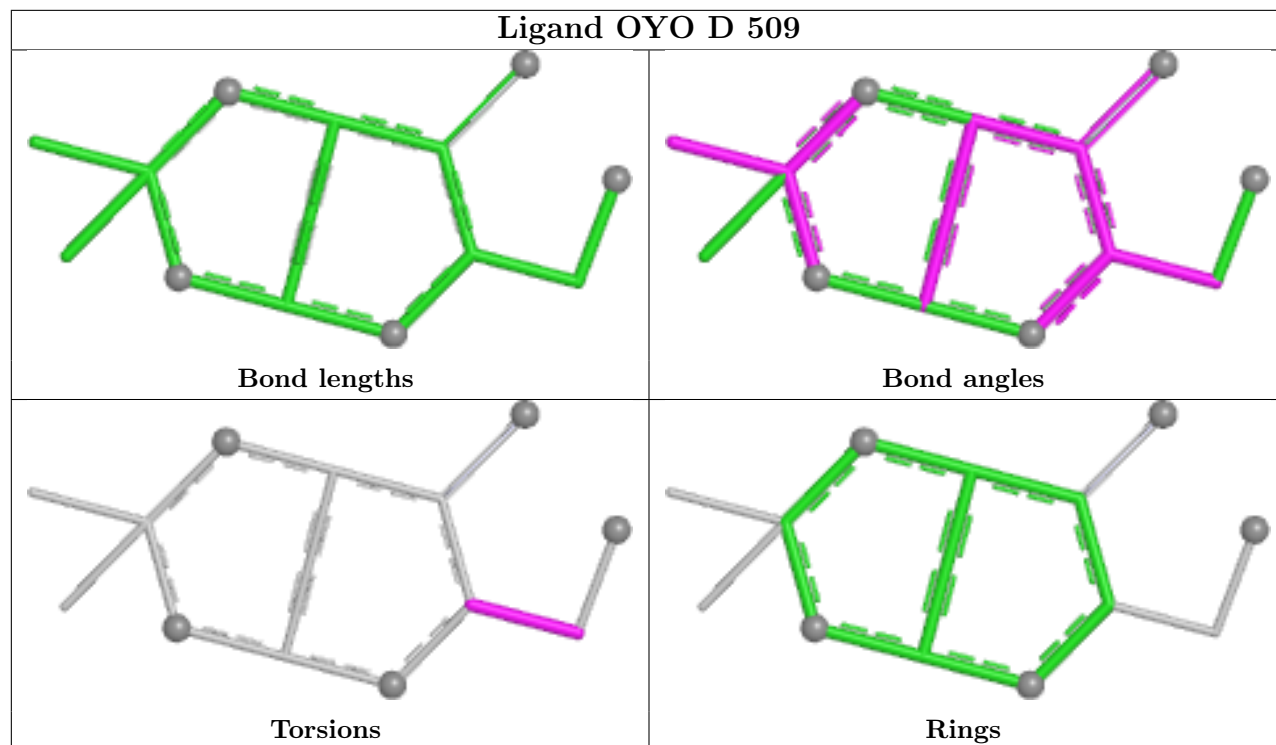
There are no ring outliers.

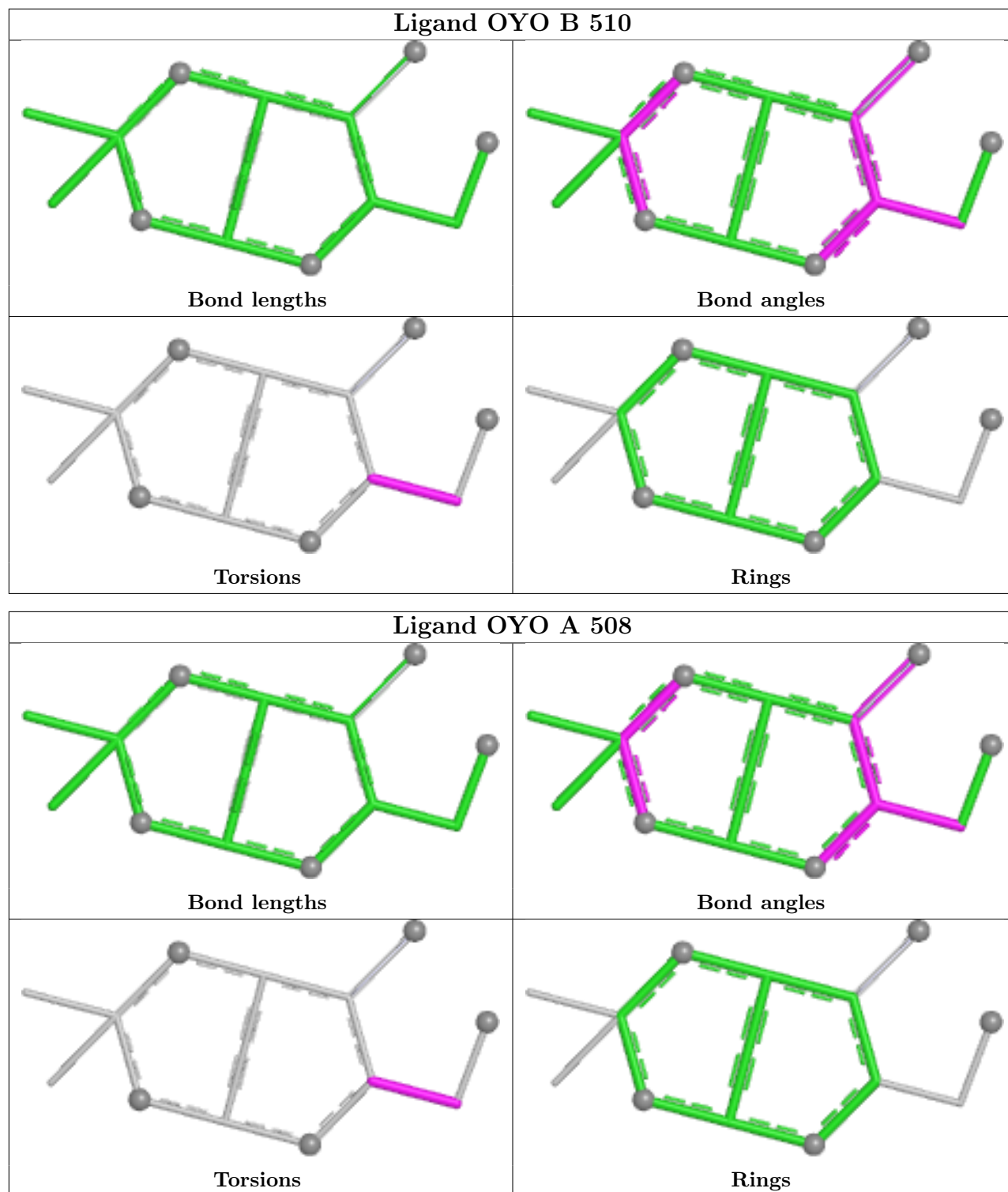
13 monomers are involved in 17 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	503	GOL	1	0
6	D	504	PEG	3	0
6	B	506	PEG	1	0
5	B	502	PGE	1	0
4	B	501	GOL	1	0
5	D	502	PGE	1	0
5	C	502	PGE	1	0
6	D	505	PEG	4	0
4	D	501	GOL	1	0
5	A	502	PGE	1	0
4	A	501	GOL	1	0
4	D	503	GOL	2	0
5	A	503	PGE	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 ⓘ

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	479/488 (98%)	-0.44	2 (0%) 88 89	10, 19, 31, 43	1 (0%)
1	B	479/488 (98%)	-0.27	0 100 100	11, 22, 35, 50	1 (0%)
1	C	479/488 (98%)	-0.22	3 (0%) 85 86	11, 23, 38, 53	3 (0%)
1	D	479/488 (98%)	-0.09	0 100 100	12, 23, 39, 57	2 (0%)
All	All	1916/1952 (98%)	-0.25	5 (0%) 90 90	10, 22, 36, 57	7 (0%)

All (5) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	171	ALA	2.8
1	C	138	TRP	2.7
1	A	138	TRP	2.4
1	C	231	SER	2.3
1	A	171	ALA	2.1

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

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

6.3 Carbohydrates [i](#)

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

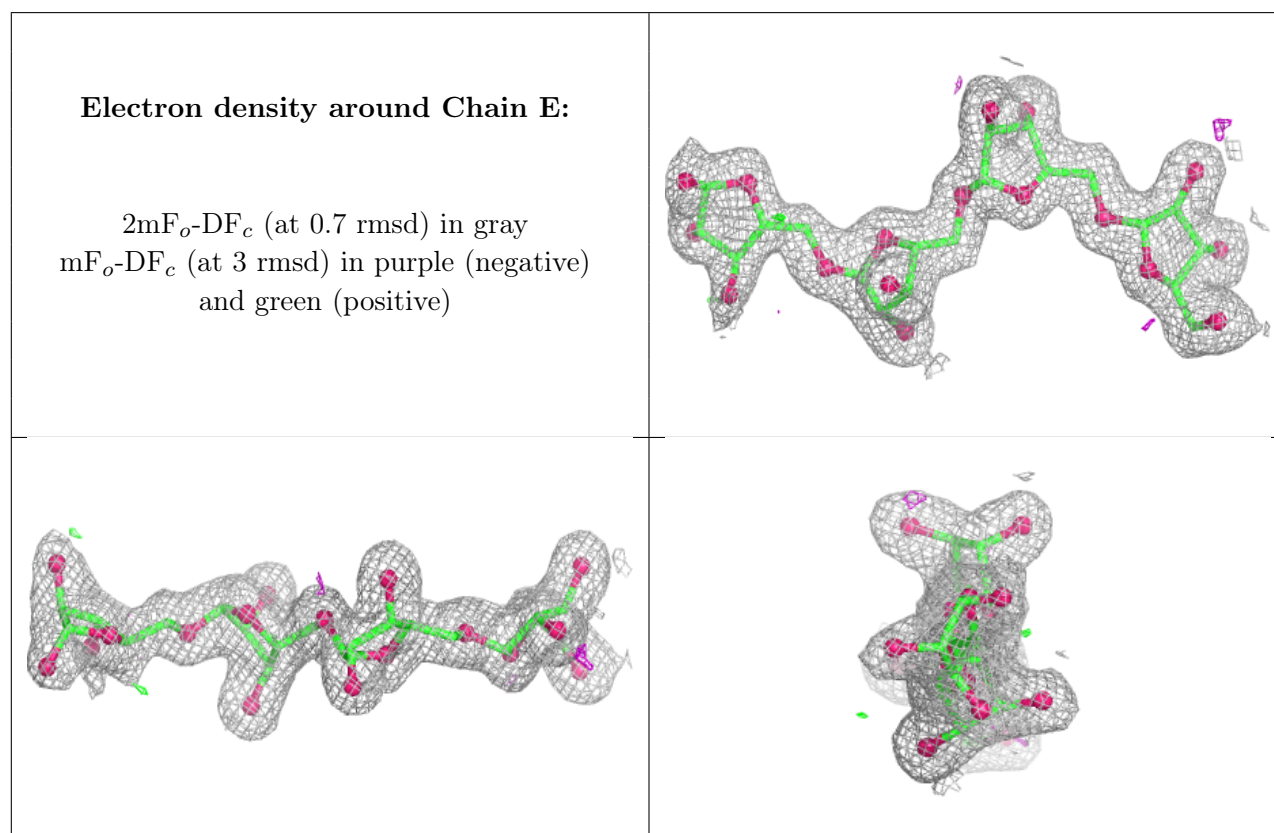
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	BXY	E	1	10/10	0.90	0.11	29,43,49,50	0
2	BXY	I	1	10/10	0.90	0.12	24,41,48,53	0

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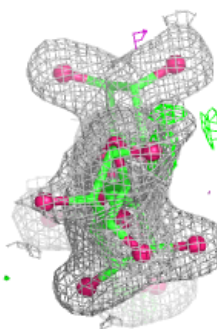
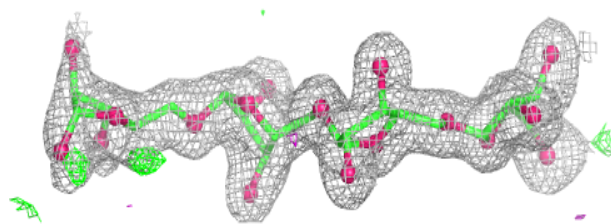
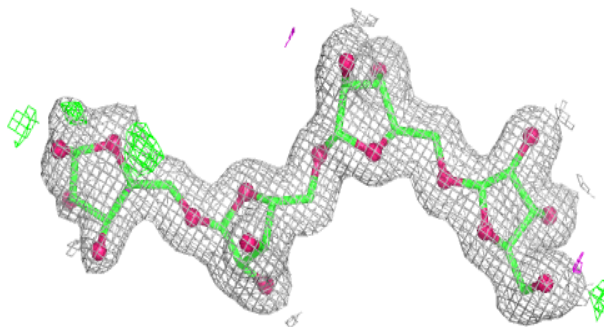
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	BXY	G	3	9/10	0.91	0.10	22,31,40,45	0
3	BXY	G	2	9/10	0.93	0.09	24,28,32,32	0
3	BXY	J	3	9/10	0.93	0.10	22,27,37,46	0
3	BXY	H	3	9/10	0.95	0.07	18,24,29,35	0
3	BXY	F	3	9/10	0.96	0.07	17,19,30,39	0
3	BXY	H	1	9/10	0.96	0.06	20,22,24,24	0
3	BXY	H	2	9/10	0.96	0.06	20,22,23,25	0
3	BXY	G	1	9/10	0.96	0.06	20,24,27,28	0
3	BXY	J	1	9/10	0.96	0.07	20,26,27,28	0
3	BXY	J	2	9/10	0.96	0.07	21,24,27,27	0
2	BXY	I	2	9/10	0.96	0.06	17,19,21,24	0
2	BXY	I	4	9/10	0.97	0.05	16,17,18,19	0
3	BXY	F	2	9/10	0.97	0.05	16,17,20,20	0
2	BXY	E	4	9/10	0.97	0.04	17,18,19,21	0
3	BXY	F	1	9/10	0.98	0.04	15,18,19,19	0
2	BXY	E	2	9/10	0.98	0.05	20,21,22,24	0
2	BXY	I	3	9/10	0.98	0.04	15,17,18,20	0
2	BXY	E	3	9/10	0.98	0.04	15,17,19,19	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.

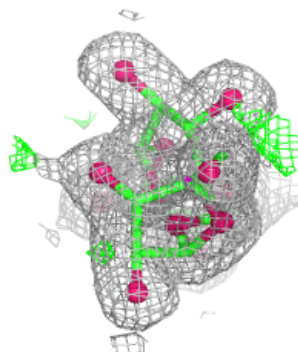
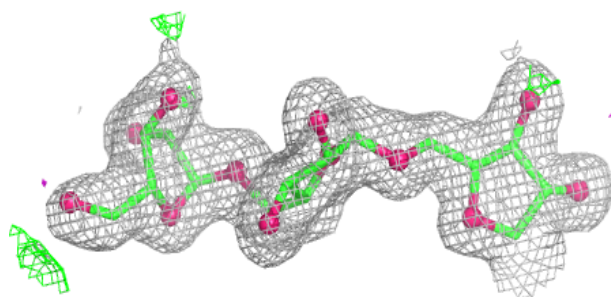
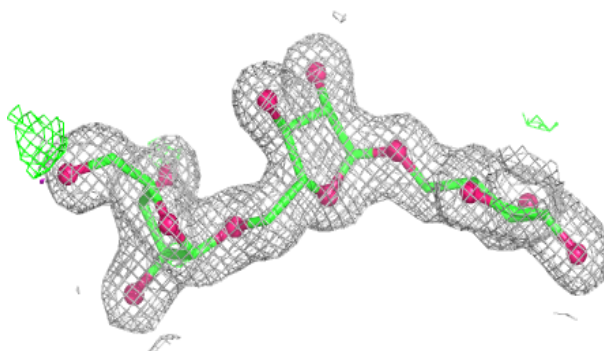


Electron density around Chain I:

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

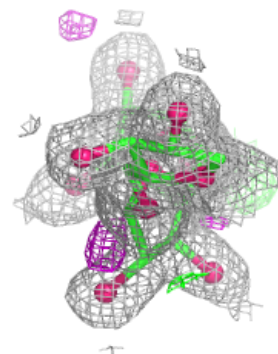
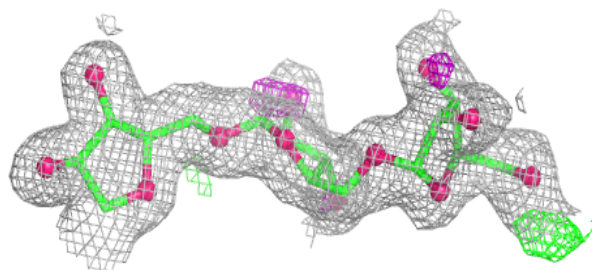
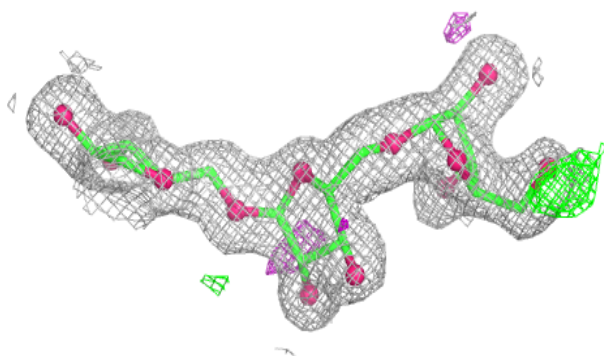
**Electron density around Chain F:**

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

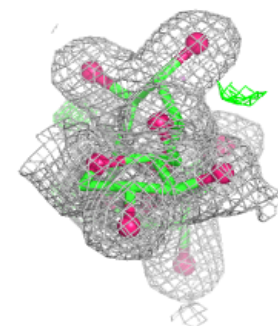
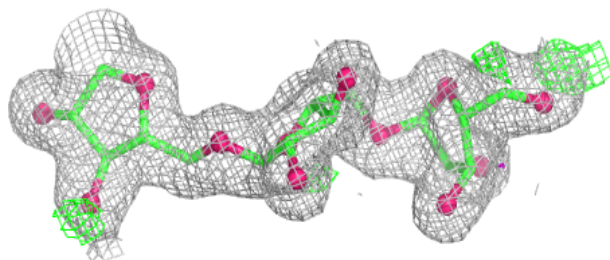
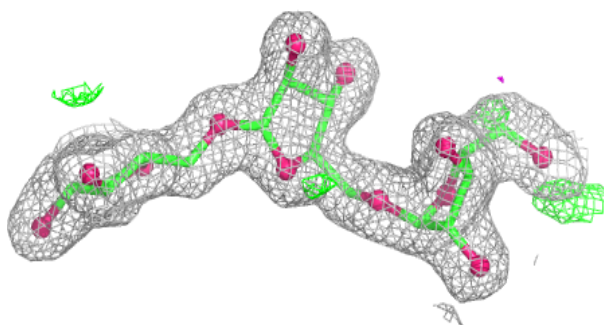


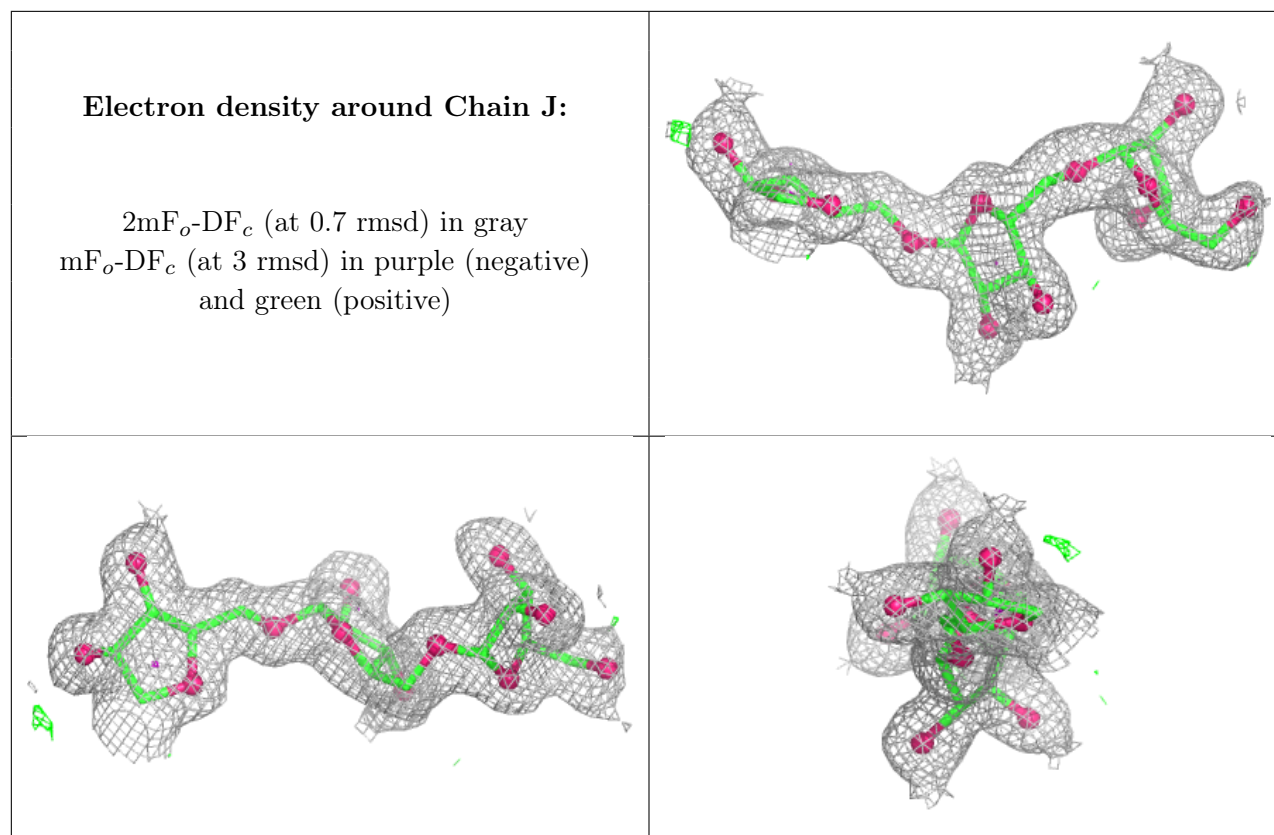
Electron density around Chain G:

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

**Electron density around Chain H:**

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





6.4 Ligands [i](#)

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
9	OYO	B	510	13/13	0.75	0.24	29,66,73,82	0
9	OYO	D	509	13/13	0.77	0.21	32,53,62,63	0
6	PEG	D	505	7/7	0.78	0.20	34,44,49,51	0
9	OYO	C	509	13/13	0.81	0.17	23,50,55,59	0
4	GOL	B	504	6/6	0.82	0.17	28,38,49,50	0
9	OYO	A	508	13/13	0.83	0.16	18,40,46,50	0
6	PEG	B	506	7/7	0.84	0.17	41,47,51,54	0
4	GOL	B	503	6/6	0.84	0.17	31,42,45,49	0
6	PEG	D	504	7/7	0.86	0.15	41,43,46,47	0
5	PGE	C	502	10/10	0.88	0.14	43,50,54,60	0
5	PGE	A	503	10/10	0.88	0.14	31,43,46,48	0
4	GOL	B	511	6/6	0.89	0.12	35,39,40,41	0
4	GOL	A	501	6/6	0.89	0.14	33,40,44,48	0
10	MES	B	505	12/12	0.89	0.13	35,40,45,48	0

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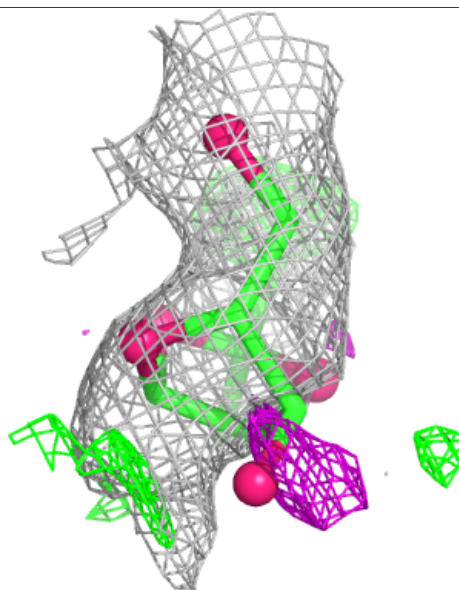
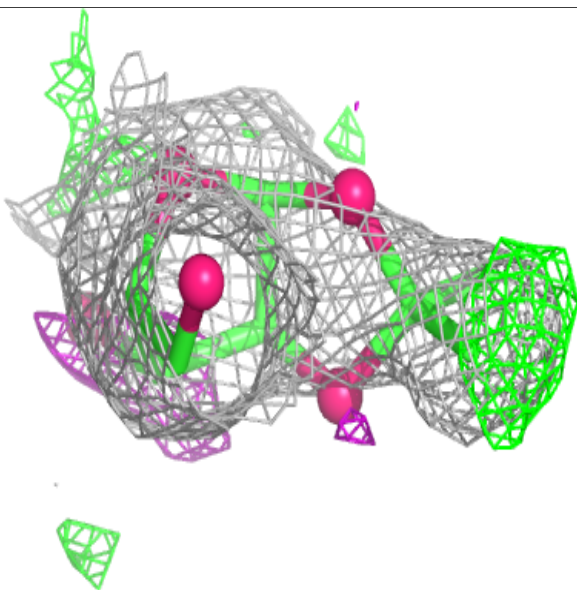
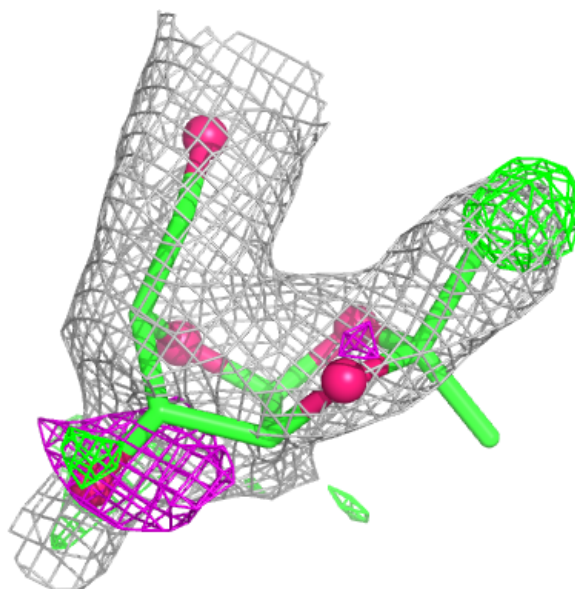
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
4	GOL	D	503	6/6	0.90	0.13	30,36,45,45	0
5	PGE	D	502	10/10	0.91	0.12	40,47,49,52	0
4	GOL	D	501	6/6	0.91	0.14	43,53,54,57	0
4	GOL	C	505	6/6	0.91	0.12	32,39,42,44	0
6	PEG	C	504	7/7	0.92	0.12	39,40,41,42	0
5	PGE	B	502	10/10	0.92	0.12	43,48,56,59	0
6	PEG	C	503	7/7	0.93	0.11	32,35,39,40	0
4	GOL	B	501	6/6	0.93	0.13	36,40,46,50	0
4	GOL	C	501	6/6	0.93	0.12	44,46,49,50	0
4	GOL	A	507	6/6	0.94	0.07	23,24,26,26	0
5	PGE	A	502	10/10	0.95	0.11	42,46,54,57	0
6	PEG	A	504	7/7	0.95	0.10	33,35,46,56	0
4	GOL	C	508	6/6	0.96	0.07	25,28,29,30	0
4	GOL	D	508	6/6	0.96	0.06	26,26,28,31	0
7	NA	D	506	1/1	0.97	0.07	28,28,28,28	0
7	NA	A	505	1/1	0.97	0.04	23,23,23,23	0
4	GOL	B	509	6/6	0.98	0.05	26,29,29,30	0
7	NA	B	507	1/1	0.98	0.05	25,25,25,25	0
7	NA	C	506	1/1	0.98	0.04	29,29,29,29	0
8	CA	B	508	1/1	1.00	0.02	18,18,18,18	0
8	CA	C	507	1/1	1.00	0.01	15,15,15,15	0
8	CA	D	507	1/1	1.00	0.02	19,19,19,19	0
8	CA	A	506	1/1	1.00	0.03	16,16,16,16	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.

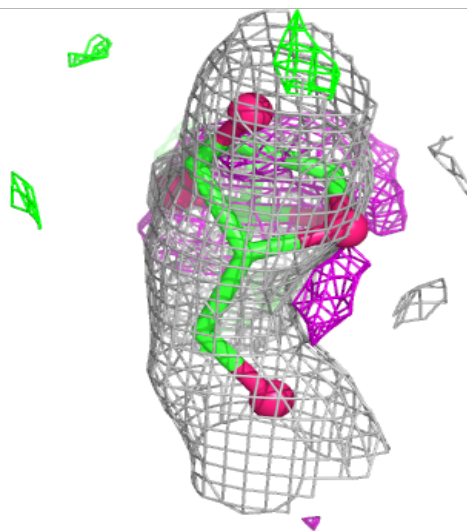
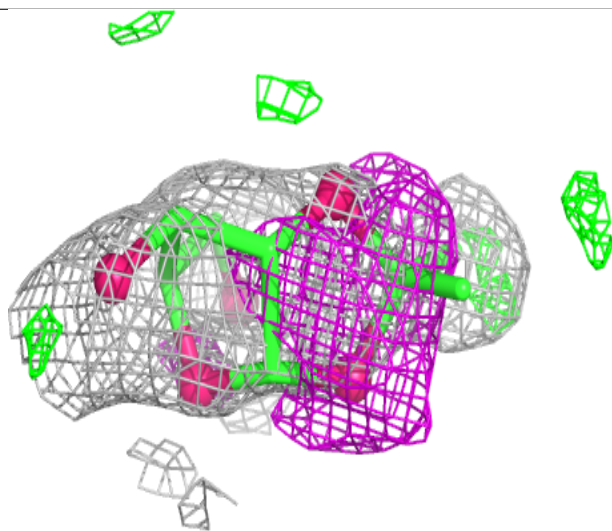
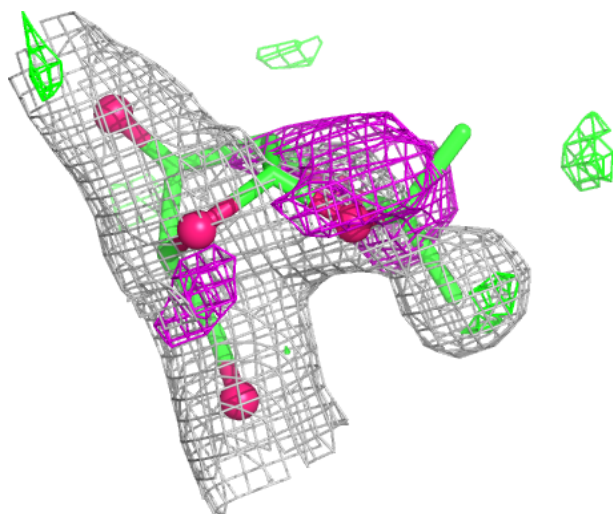
Electron density around OYO B 510:

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



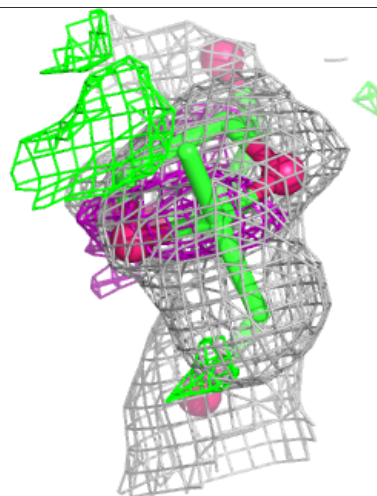
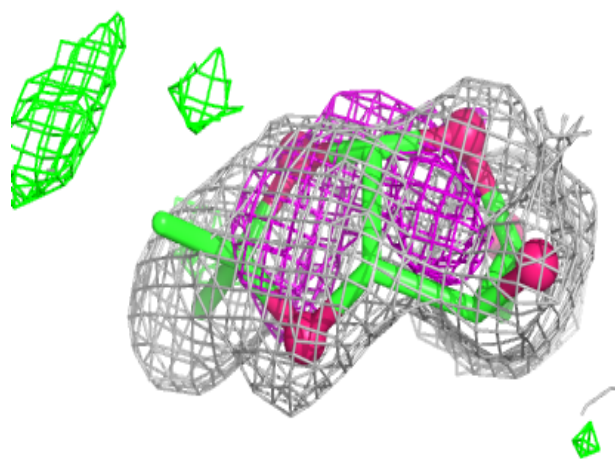
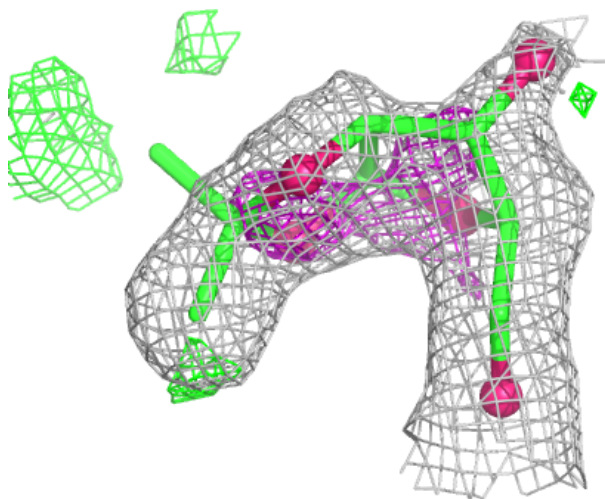
Electron density around OYO D 509:

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



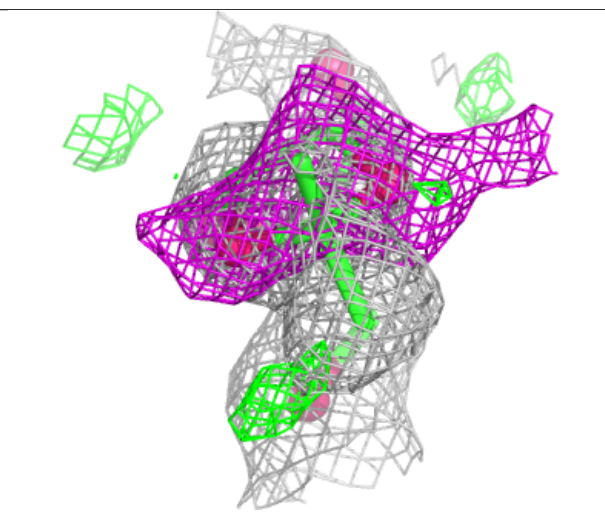
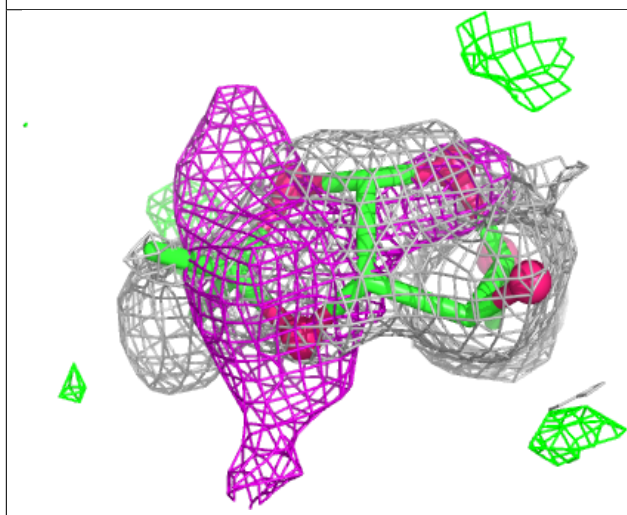
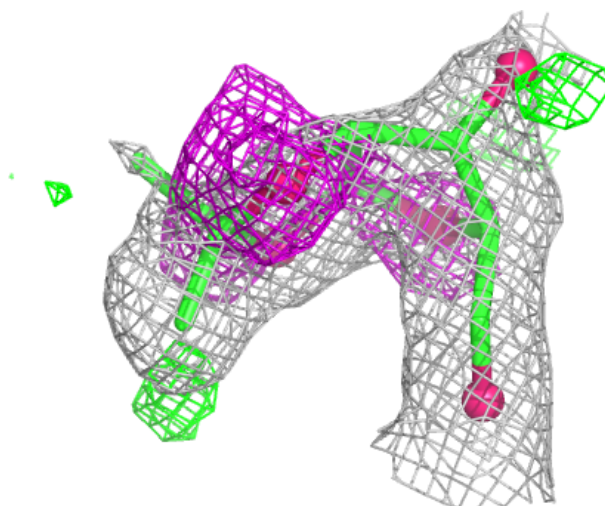
Electron density around OYO C 509:

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



Electron density around OYO A 508:

$2mF_o-DF_c$ (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.