



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 02:42 PM UTC

PDB ID : 2HPY / pdb_00002hpy
Title : Crystallographic model of lumirhodopsin
Authors : Nakamichi, H.; Okada, T.
Deposited on : 2006-07-18
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

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

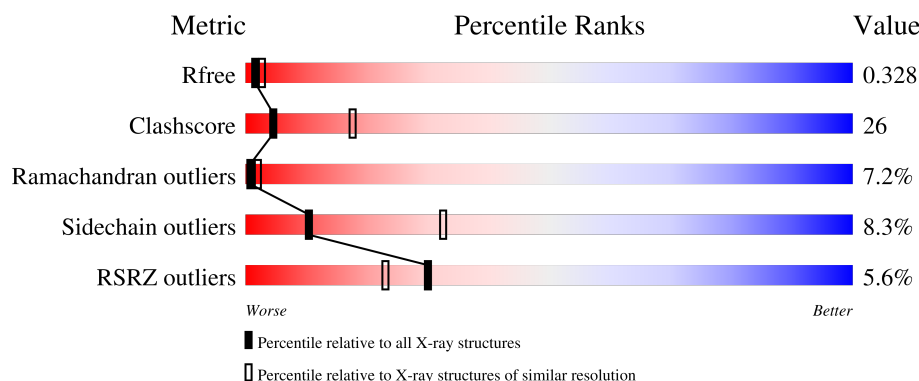
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 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}	180053	3866 (2.80-2.80)
Clashscore	190562	4276 (2.80-2.80)
Ramachandran outliers	187476	4196 (2.80-2.80)
Sidechain outliers	187428	4198 (2.80-2.80)
RSRZ outliers	180081	3869 (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 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	349	4% 66% 27% 7%
1	B	349	7% 62% 30% 7%
2	C	3	33% 33% 33%
3	D	2	100%
3	F	2	50% 50%

Continued on next page...

Continued from previous page...

Mol	Chain	Length	Quality of chain
4	E	4	

2 Entry composition [i](#)

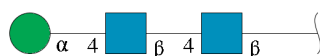
There are 11 unique types of molecules in this entry. The entry contains 5948 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Rhodopsin.

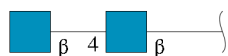
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	349	Total	C	N	O	S	0	0	0
			2749	1818	424	481	26			
1	B	349	Total	C	N	O	S	0	0	0
			2749	1818	424	481	26			

- Molecule 2 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.



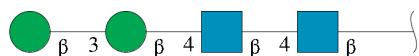
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	C	3	Total	C	N	O	0	0	0
			39	22	2	15			

- 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	F	2	Total	C	N	O	0	0	0
			28	16	2	10			

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



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
4	E	4	Total	C	N	O	0	0	0
			50	28	2	20			

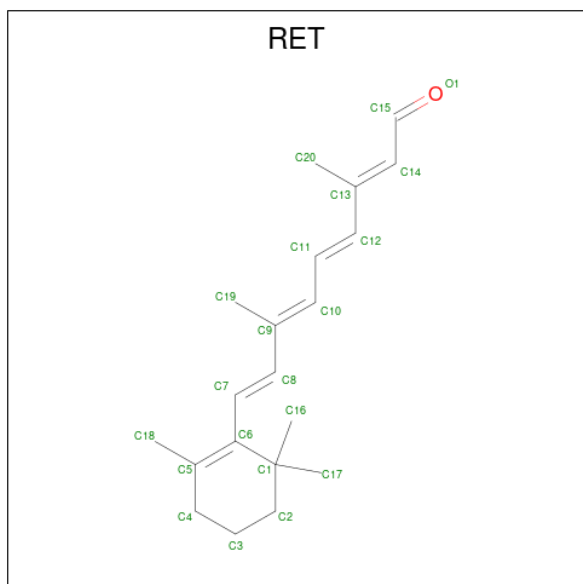
- Molecule 5 is MERCURY (II) ION (CCD ID: HG) (formula: Hg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	3	Total	Hg	0	0
			3	3		
5	B	3	Total	Hg	0	0
			3	3		

- Molecule 6 is ZINC ION (CCD ID: ZN) (formula: Zn).

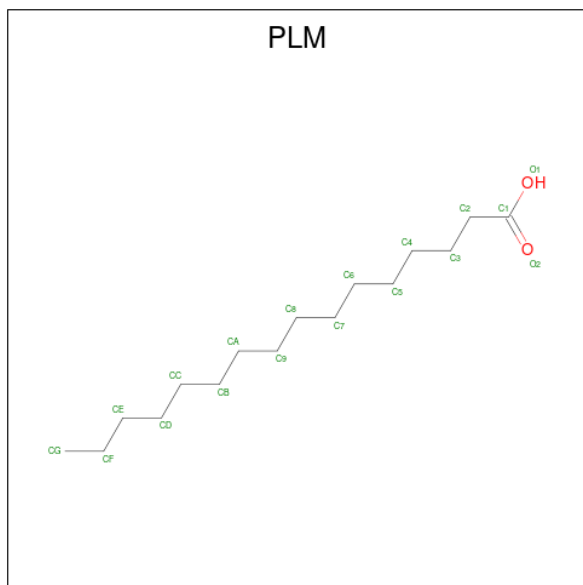
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	4	Total	Zn	0	0
			4	4		
6	B	3	Total	Zn	0	0
			3	3		

- Molecule 7 is RETINAL (CCD ID: RET) (formula: C₂₀H₂₈O).



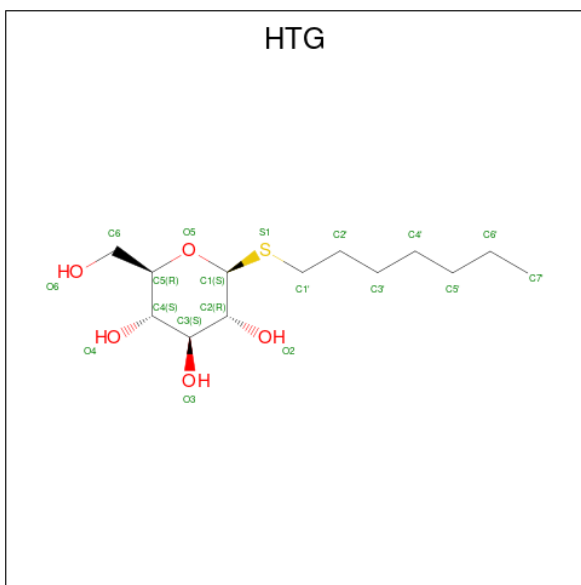
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
7	A	1	Total C 20 20	0	0
7	B	1	Total C 20 20	0	0

- Molecule 8 is PALMITIC ACID (CCD ID: PLM) (formula: $C_{16}H_{32}O_2$).



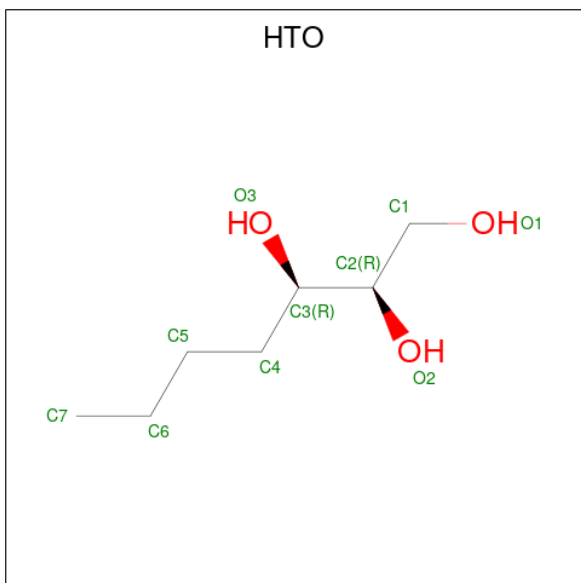
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
8	A	1	Total C O 17 16 1	0	0
8	A	1	Total C O 17 16 1	0	0
8	A	1	Total C 16 16	0	0
8	B	1	Total C O 17 16 1	0	0
8	B	1	Total C O 17 16 1	0	0
8	B	1	Total C 16 16	0	0

- Molecule 9 is heptyl 1-thio-beta-D-glucopyranoside (CCD ID: HTG) (formula: $C_{13}H_{26}O_5S$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
9	A	1	Total	C	O	S	0	0
			19	13	5	1		
9	A	1	Total	C	O	S	0	0
			19	13	5	1		
9	B	1	Total	C	O	S	0	0
			19	13	5	1		
9	B	1	Total	C	O	S	0	0
			19	13	5	1		

- Molecule 10 is HEPTANE-1,2,3-TRIOL (CCD ID: HTO) (formula: $C_7H_{16}O_3$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
10	B	1	Total	C	O	0	0
			10	7	3		

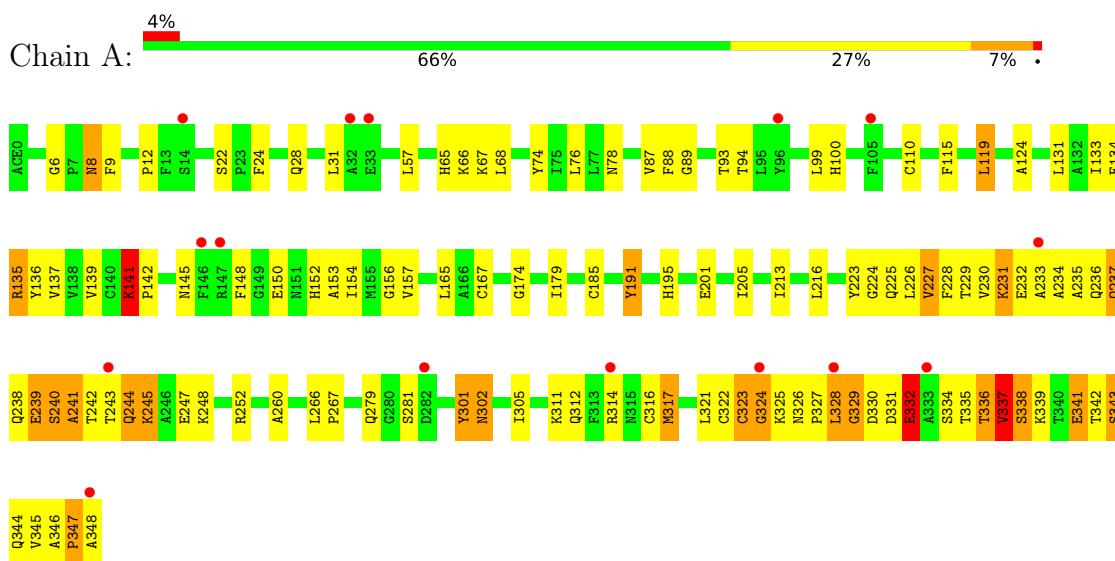
- Molecule 11 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
11	A	37	Total	O	0	0
			37	37		
11	B	29	Total	O	0	0
			29	29		

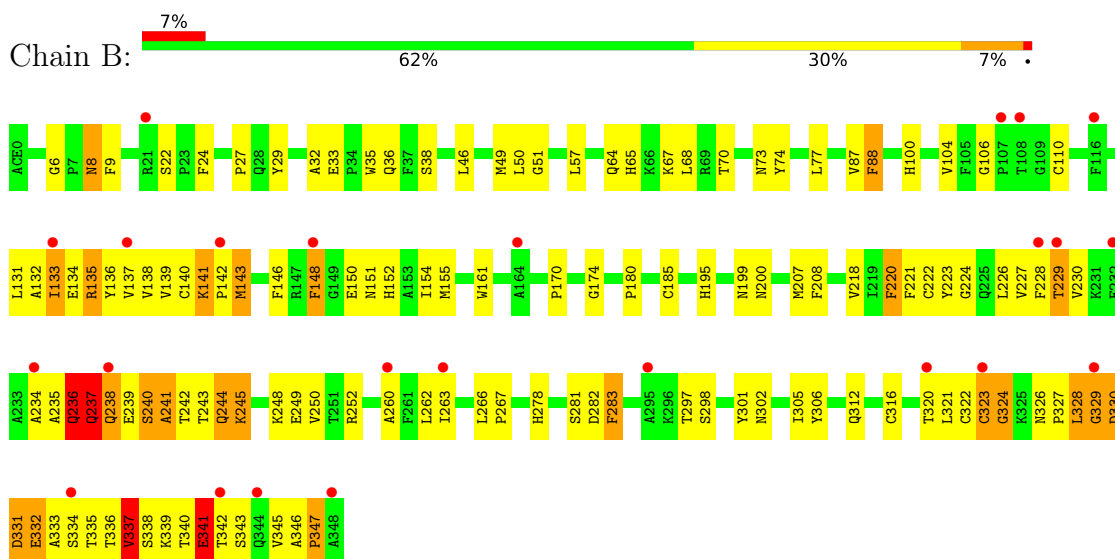
3 Residue-property plots

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.

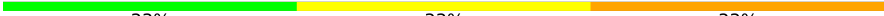
• Molecule 1: Rhodopsin



• Molecule 1: Rhodopsin

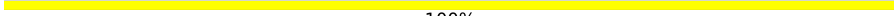


• Molecule 2: alpha-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain C:  33% 33% 33%



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

Chain D:  100%



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

Chain F:  50% 50%



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

Chain E:  50% 50%



4 Data and refinement statistics

Property	Value	Source
Space group	P 41	Depositor
Cell constants a, b, c, α , β , γ	96.72Å 96.72Å 150.40Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 2.80 50.00 – 2.80	Depositor EDS
% Data completeness (in resolution range)	(Not available) (50.00-2.80) 87.6 (50.00-2.80)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.88 (at 2.81Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.218 , 0.238 0.343 , 0.328	Depositor DCC
R_{free} test set	1531 reflections (4.47%)	wwPDB-VP
Wilson B-factor (Å ²)	64.8	Xtriage
Anisotropy	0.474	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 242.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	0.138 for h,-k,-l	Xtriage
F_o, F_c correlation	0.85	EDS
Total number of atoms	5948	wwPDB-VP
Average B, all atoms (Å ²)	75.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.70% 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: ACE, RET, HTG, MAN, HTO, PLM, NAG, ZN, HG, BMA

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	0.88	2/2831 (0.1%)	1.08	10/3859 (0.3%)
1	B	0.85	1/2831 (0.0%)	1.06	9/3859 (0.2%)
All	All	0.87	3/5662 (0.1%)	1.07	19/7718 (0.2%)

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	B	0	1

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	88	PHE	CA-C	-5.21	1.45	1.52
1	A	339	LYS	CA-C	5.15	1.56	1.52
1	A	317	MET	SD-CE	5.12	1.92	1.79

All (19) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	174	GLY	N-CA-C	8.27	124.92	115.08
1	A	191	TYR	N-CA-C	-7.40	104.50	113.97
1	B	170	PRO	N-CA-C	7.17	119.45	110.70
1	B	106	GLY	CA-C-N	6.89	126.36	118.85
1	B	106	GLY	C-N-CA	6.89	126.36	118.85
1	A	239	GLU	N-CA-C	-6.67	106.09	114.56
1	A	343	SER	N-CA-C	-6.54	105.60	112.93
1	A	141	LYS	CA-C-N	6.51	126.44	119.28

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	141	LYS	C-N-CA	6.51	126.44	119.28
1	B	77	LEU	N-CA-C	-6.42	104.21	111.14
1	A	156	GLY	N-CA-C	-5.82	105.75	112.50
1	A	227	VAL	N-CA-C	-5.72	106.24	111.67
1	B	174	GLY	CA-C-N	-5.67	115.47	122.84
1	B	174	GLY	C-N-CA	-5.67	115.47	122.84
1	B	250	VAL	N-CA-C	-5.59	104.10	111.09
1	B	297	THR	N-CA-C	-5.54	106.36	113.01
1	B	38	SER	N-CA-C	-5.28	105.70	111.82
1	A	332	GLU	N-CA-C	-5.24	106.28	113.30
1	A	185	CYS	N-CA-C	-5.19	107.62	114.31

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	306	TYR	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	2749	0	2709	134	1
1	B	2749	0	2709	170	0
2	C	39	0	34	1	0
3	D	28	0	25	0	0
3	F	28	0	25	1	0
4	E	50	0	43	0	0
5	A	3	0	0	0	0
5	B	3	0	0	0	0
6	A	4	0	0	0	0
6	B	3	0	0	0	0
7	A	20	0	27	0	0
7	B	20	0	27	1	0
8	A	50	0	89	7	0
8	B	50	0	89	4	0
9	A	38	0	52	3	0

Continued on next page...

Continued from previous page...

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
9	B	38	0	52	4	1
10	B	10	0	16	1	0
11	A	37	0	0	6	0
11	B	29	0	0	11	0
All	All	5948	0	5897	307	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 26.

All (307) 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:67:LYS:HB2	1:A:337:VAL:HB	1.38	1.02
1:B:64:GLN:HG3	1:B:339:LYS:HE2	1.39	1.02
1:A:345:VAL:HG12	1:A:347:PRO:HD3	1.45	0.98
1:B:337:VAL:HB	1:B:343:SER:HA	1.45	0.97
1:B:67:LYS:HZ3	1:B:336:THR:HA	1.30	0.96
1:B:239:GLU:HB2	1:B:245:LYS:HD2	1.44	0.95
1:A:65:HIS:ND1	1:A:338:SER:HA	1.81	0.94
1:A:88:PHE:HB3	8:B:1323:PLM:HB1	1.49	0.94
1:B:245:LYS:HE3	1:B:245:LYS:HA	1.51	0.92
1:A:239:GLU:HG2	1:A:248:LYS:NZ	1.85	0.91
1:A:332:GLU:C	1:A:334:SER:H	1.77	0.88
1:B:298:SER:HB2	11:B:2036:HOH:O	1.73	0.88
1:B:245:LYS:NZ	1:B:248:LYS:HD2	1.91	0.86
1:A:316:CYS:SG	1:A:337:VAL:HG13	2.15	0.85
1:A:65:HIS:HB3	1:A:337:VAL:HG22	1.60	0.84
1:B:143:MET:HA	1:B:143:MET:HE3	1.59	0.83
1:B:346:ALA:N	1:B:347:PRO:HD3	1.93	0.83
1:B:67:LYS:HB3	1:B:337:VAL:CG1	2.10	0.82
1:B:267:PRO:HG2	11:B:2019:HOH:O	1.80	0.81
1:A:341:GLU:CG	1:A:342:THR:H	1.94	0.81
1:A:239:GLU:HG2	1:A:248:LYS:HZ1	1.46	0.79
1:A:326:ASN:O	1:A:328:LEU:HD22	1.83	0.79
1:A:341:GLU:HG3	1:A:342:THR:N	1.98	0.77
1:B:312:GLN:HA	1:B:332:GLU:HG2	1.65	0.77
1:A:267:PRO:HG2	11:A:964:HOH:O	1.84	0.77
1:B:67:LYS:HE3	1:B:312:GLN:HG3	1.67	0.77
1:A:245:LYS:NZ	1:A:245:LYS:HA	2.00	0.77
1:B:248:LYS:O	1:B:252:ARG:HG3	1.83	0.76
1:B:332:GLU:HB3	1:B:335:THR:O	1.85	0.76

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:338:SER:HB2	1:B:341:GLU:HG3	1.66	0.76
1:A:345:VAL:HG12	1:A:347:PRO:CD	2.16	0.75
1:A:65:HIS:HD1	1:A:338:SER:HA	1.47	0.75
1:B:67:LYS:H	1:B:337:VAL:CG2	2.00	0.75
1:A:67:LYS:HB2	1:A:337:VAL:CB	2.15	0.73
1:B:230:VAL:HG23	1:B:248:LYS:HD3	1.70	0.73
1:B:337:VAL:HA	1:B:342:THR:C	2.13	0.73
1:B:337:VAL:HA	1:B:342:THR:O	1.89	0.72
1:B:139:VAL:HG11	1:B:230:VAL:HG12	1.71	0.72
1:B:337:VAL:CB	1:B:343:SER:HA	2.19	0.72
1:B:278:HIS:NE2	11:B:2038:HOH:O	2.23	0.72
1:A:150:GLU:O	1:A:154:ILE:HG13	1.90	0.72
1:A:341:GLU:OE1	1:A:342:THR:HG22	1.88	0.72
1:A:325:LYS:HE3	1:A:341:GLU:OE1	1.90	0.71
1:A:100:HIS:CE1	9:B:1509:HTG:O6	2.44	0.71
1:B:234:ALA:HA	1:B:245:LYS:NZ	2.05	0.71
1:A:245:LYS:HA	1:A:245:LYS:HZ2	1.55	0.71
1:B:237:GLN:H	1:B:237:GLN:NE2	1.89	0.71
1:A:302:ASN:HB2	11:A:2017:HOH:O	1.89	0.70
1:A:341:GLU:HG3	1:A:342:THR:H	1.55	0.70
1:A:326:ASN:HD21	1:A:328:LEU:HD11	1.55	0.70
1:B:67:LYS:N	1:B:337:VAL:HG21	2.06	0.70
1:B:74:TYR:HE2	1:B:150:GLU:HG2	1.57	0.70
1:A:239:GLU:HB3	1:A:244:GLN:CD	2.17	0.70
1:B:64:GLN:HG3	1:B:339:LYS:HB2	1.74	0.69
1:B:301:TYR:HE1	11:B:2008:HOH:O	1.75	0.69
1:A:332:GLU:C	1:A:334:SER:N	2.49	0.69
1:B:64:GLN:CG	1:B:339:LYS:HB2	2.22	0.69
1:A:67:LYS:H	1:A:337:VAL:HG23	1.57	0.69
1:B:338:SER:HB2	1:B:341:GLU:CG	2.23	0.69
1:B:239:GLU:HB3	1:B:244:GLN:HE22	1.59	0.68
1:A:65:HIS:HD1	1:A:338:SER:CA	2.06	0.68
1:B:337:VAL:HG23	1:B:338:SER:N	2.09	0.68
1:B:143:MET:HG2	1:B:146:PHE:HB3	1.76	0.68
1:A:65:HIS:ND1	1:A:337:VAL:O	2.27	0.68
1:B:238:GLN:HG3	1:B:241:ALA:H	1.59	0.68
1:B:143:MET:HG2	1:B:146:PHE:CB	2.24	0.67
1:A:341:GLU:CG	1:A:342:THR:N	2.57	0.67
1:B:332:GLU:OE2	1:B:334:SER:HB2	1.94	0.67
1:A:65:HIS:HD1	1:A:337:VAL:C	2.04	0.66
1:B:65:HIS:HB2	1:B:68:LEU:HD12	1.77	0.66

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:239:GLU:HG2	1:A:248:LYS:HZ2	1.61	0.66
1:B:67:LYS:HB3	1:B:337:VAL:HG11	1.78	0.65
1:B:67:LYS:HD3	1:B:337:VAL:HG13	1.78	0.65
1:A:238:GLN:HA	1:A:242:THR:HA	1.79	0.65
1:B:67:LYS:HZ3	1:B:336:THR:CA	2.06	0.65
1:A:137:VAL:HA	1:A:142:PRO:CD	2.27	0.65
1:A:325:LYS:HE2	1:A:327:PRO:HD3	1.79	0.65
1:A:99:LEU:HD11	9:B:1509:HTG:H3'1	1.77	0.65
1:B:237:GLN:O	1:B:238:GLN:HB3	1.97	0.65
1:A:338:SER:HB2	1:A:341:GLU:HG2	1.79	0.64
1:B:238:GLN:HA	1:B:242:THR:HA	1.79	0.63
1:A:235:ALA:HB3	1:A:239:GLU:OE1	1.98	0.63
1:A:337:VAL:O	1:A:338:SER:OG	2.17	0.63
1:A:110:CYS:HB3	11:A:2045:HOH:O	1.99	0.63
1:A:326:ASN:OD1	1:A:328:LEU:HD13	1.99	0.63
1:A:227:VAL:HG13	1:A:228:PHE:N	2.13	0.63
1:B:199:ASN:ND2	11:B:2005:HOH:O	2.31	0.62
1:A:322:CYS:O	1:A:324:GLY:N	2.32	0.62
1:A:57:LEU:HD23	8:A:1322:PLM:HD2	1.81	0.62
1:B:6:GLY:HA3	1:B:9:PHE:CZ	2.34	0.61
1:B:150:GLU:O	1:B:154:ILE:HD13	2.00	0.61
1:A:342:THR:O	1:A:342:THR:HG23	2.01	0.61
1:A:234:ALA:HB2	1:A:248:LYS:HG2	1.82	0.61
1:B:236:GLN:HB2	1:B:237:GLN:NE2	2.15	0.61
1:B:302:ASN:HB2	11:B:2016:HOH:O	2.00	0.61
1:B:337:VAL:HB	1:B:343:SER:CA	2.26	0.60
1:A:67:LYS:H	1:A:337:VAL:CG2	2.13	0.60
1:A:224:GLY:O	1:A:227:VAL:HG12	2.01	0.60
1:A:65:HIS:HB3	1:A:337:VAL:CG2	2.31	0.60
1:A:312:GLN:NE2	1:A:337:VAL:HG12	2.16	0.60
1:B:345:VAL:C	1:B:347:PRO:HD3	2.26	0.60
1:B:245:LYS:HZ1	1:B:248:LYS:HD2	1.67	0.60
1:B:327:PRO:HB2	1:B:331:ASP:OD2	2.02	0.60
1:B:143:MET:HA	1:B:143:MET:CE	2.32	0.60
1:B:64:GLN:HG3	1:B:339:LYS:CE	2.24	0.59
1:A:137:VAL:HA	1:A:142:PRO:HD2	1.85	0.59
1:B:67:LYS:NZ	1:B:332:GLU:HG3	2.17	0.59
1:A:66:LYS:HD2	1:A:66:LYS:N	2.18	0.59
1:B:46:LEU:HD12	8:B:1407:PLM:HD2	1.85	0.59
1:B:67:LYS:HB3	1:B:337:VAL:CG2	2.32	0.59
1:B:220:PHE:O	1:B:223:TYR:HB3	2.04	0.58

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:6:GLY:HA3	1:A:9:PHE:CZ	2.38	0.58
1:B:234:ALA:HB1	1:B:245:LYS:HE2	1.86	0.58
1:B:241:ALA:HB1	1:B:243:THR:HG22	1.86	0.58
1:B:326:ASN:C	1:B:328:LEU:HD22	2.29	0.58
1:A:239:GLU:HB3	1:A:244:GLN:NE2	2.18	0.58
1:B:180:PRO:HG2	11:B:2023:HOH:O	2.03	0.57
1:B:332:GLU:O	1:B:333:ALA:HB3	2.04	0.57
1:A:326:ASN:ND2	1:A:328:LEU:HD21	2.19	0.56
1:B:234:ALA:HA	1:B:245:LYS:HZ3	1.70	0.56
1:B:312:GLN:HA	1:B:332:GLU:CG	2.34	0.56
1:B:239:GLU:HB3	1:B:244:GLN:NE2	2.19	0.56
1:B:326:ASN:O	1:B:328:LEU:HD22	2.04	0.56
1:A:237:GLN:CD	1:A:237:GLN:H	2.14	0.56
1:B:239:GLU:HB2	1:B:245:LYS:CD	2.28	0.56
1:A:65:HIS:HB2	1:A:68:LEU:HD12	1.86	0.56
1:B:226:LEU:N	1:B:226:LEU:HD12	2.20	0.56
1:A:230:VAL:HG23	1:A:248:LYS:HE2	1.88	0.56
8:A:1322:PLM:H61	1:B:88:PHE:HE1	1.69	0.55
1:B:244:GLN:O	1:B:248:LYS:HG3	2.05	0.55
1:A:65:HIS:ND1	1:A:338:SER:CA	2.61	0.55
1:A:346:ALA:N	1:A:347:PRO:HD3	2.21	0.55
1:B:245:LYS:CE	1:B:248:LYS:HD2	2.36	0.55
1:A:338:SER:CB	1:A:341:GLU:HG2	2.36	0.55
1:B:57:LEU:HD21	1:B:321:LEU:HD21	1.89	0.55
1:B:326:ASN:HB3	1:B:328:LEU:CD2	2.37	0.55
1:B:33:GLU:HB2	1:B:36:GLN:HG3	1.89	0.55
1:A:311:LYS:HG2	1:A:331:ASP:O	2.06	0.54
1:B:135:ARG:NE	1:B:135:ARG:HA	2.22	0.54
1:B:337:VAL:HA	1:B:342:THR:HG23	1.90	0.54
1:B:143:MET:CG	1:B:146:PHE:HB3	2.37	0.54
1:B:136:TYR:HA	1:B:226:LEU:HD11	1.88	0.54
1:B:238:GLN:HG3	1:B:241:ALA:N	2.21	0.54
1:A:267:PRO:HA	9:A:1507:HTG:H6'1	1.90	0.54
1:A:139:VAL:HG11	1:A:226:LEU:HG	1.90	0.54
1:B:67:LYS:CE	1:B:312:GLN:HG3	2.35	0.54
1:B:237:GLN:H	1:B:237:GLN:HE21	1.55	0.53
1:A:232:GLU:HA	1:A:252:ARG:HD3	1.91	0.53
1:A:237:GLN:O	1:A:238:GLN:HB3	2.07	0.53
1:A:332:GLU:O	1:A:334:SER:N	2.42	0.53
1:A:227:VAL:CG1	1:A:228:PHE:N	2.71	0.53
1:A:311:LYS:CE	1:A:330:ASP:HB2	2.38	0.53

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:321:LEU:C	1:A:323:CYS:H	2.16	0.53
1:B:64:GLN:HG2	1:B:339:LYS:HB2	1.91	0.53
1:B:322:CYS:HA	8:B:1322:PLM:O1	2.08	0.53
1:B:139:VAL:HG21	1:B:226:LEU:HD23	1.92	0.52
1:B:238:GLN:HG3	1:B:241:ALA:CA	2.39	0.52
1:A:321:LEU:HD13	8:A:1322:PLM:H72	1.91	0.52
1:B:235:ALA:HB3	1:B:239:GLU:OE1	2.09	0.52
1:B:332:GLU:CB	1:B:335:THR:O	2.57	0.52
1:B:221:PHE:O	1:B:224:GLY:N	2.37	0.52
1:B:139:VAL:CG1	1:B:230:VAL:HG12	2.39	0.52
1:B:234:ALA:HA	1:B:245:LYS:HZ2	1.75	0.52
1:B:32:ALA:HB1	1:B:36:GLN:OE1	2.10	0.52
1:B:67:LYS:HD3	1:B:336:THR:C	2.35	0.52
1:B:321:LEU:C	1:B:323:CYS:H	2.16	0.52
1:B:337:VAL:CA	1:B:342:THR:HG23	2.40	0.52
1:B:245:LYS:HZ3	1:B:248:LYS:HD2	1.73	0.52
1:B:238:GLN:HG3	1:B:241:ALA:C	2.34	0.52
1:B:328:LEU:O	1:B:329:GLY:C	2.53	0.51
1:A:133:ILE:O	1:A:136:TYR:N	2.44	0.51
1:A:139:VAL:HG11	1:A:230:VAL:HG12	1.91	0.51
1:B:332:GLU:O	1:B:332:GLU:OE1	2.29	0.51
1:B:146:PHE:HE1	1:B:152:HIS:NE2	2.09	0.51
2:C:1:NAG:H61	2:C:2:NAG:C7	2.40	0.51
1:B:239:GLU:CB	1:B:244:GLN:HE22	2.23	0.51
1:A:239:GLU:HB3	1:A:244:GLN:OE1	2.10	0.50
8:A:1322:PLM:H91	1:B:49:MET:SD	2.51	0.50
1:B:230:VAL:HG23	1:B:248:LYS:CD	2.39	0.50
1:B:67:LYS:HZ1	1:B:332:GLU:HG3	1.75	0.50
1:A:67:LYS:HB2	1:A:337:VAL:CG2	2.42	0.50
1:B:330:ASP:O	1:B:331:ASP:O	2.29	0.50
1:B:64:GLN:HA	1:B:339:LYS:HD3	1.93	0.49
1:A:247:GLU:HA	1:A:247:GLU:OE1	2.11	0.49
1:B:340:THR:HG23	1:B:341:GLU:N	2.27	0.49
1:B:346:ALA:N	1:B:347:PRO:CD	2.71	0.49
1:B:227:VAL:HG13	1:B:228:PHE:N	2.26	0.49
1:A:139:VAL:HG21	1:A:230:VAL:HG11	1.95	0.49
1:B:227:VAL:CG1	1:B:228:PHE:N	2.75	0.49
1:B:239:GLU:HB3	1:B:244:GLN:OE1	2.13	0.49
1:B:67:LYS:N	1:B:337:VAL:CG2	2.66	0.48
1:A:12:PRO:HD3	11:A:2064:HOH:O	2.11	0.48
1:A:240:SER:O	1:A:241:ALA:HB2	2.12	0.48

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
10:B:1401:HTO:H72	9:B:1509:HTG:H3'2	1.95	0.48
1:A:327:PRO:O	1:A:328:LEU:C	2.56	0.48
1:A:245:LYS:NZ	1:A:245:LYS:CA	2.76	0.48
1:B:282:ASP:HB2	3:F:1:NAG:H62	1.96	0.48
1:A:153:ALA:O	1:A:157:VAL:HG23	2.13	0.48
1:B:67:LYS:HB3	1:B:337:VAL:HG13	1.91	0.48
1:A:135:ARG:HA	1:A:135:ARG:NE	2.29	0.48
1:B:136:TYR:CE2	1:B:142:PRO:HG2	2.49	0.48
1:B:221:PHE:CD1	1:B:221:PHE:C	2.92	0.48
1:A:232:GLU:HB3	1:A:252:ARG:HD3	1.96	0.47
1:B:151:ASN:HD21	1:B:152:HIS:CE1	2.32	0.47
1:A:311:LYS:HE3	1:A:331:ASP:O	2.15	0.47
1:A:336:THR:C	1:A:337:VAL:HG12	2.38	0.47
1:B:238:GLN:HB2	1:B:242:THR:OG1	2.14	0.47
1:A:311:LYS:NZ	1:A:330:ASP:HB2	2.29	0.47
1:A:342:THR:C	1:A:344:GLN:H	2.21	0.47
1:B:67:LYS:H	1:B:337:VAL:HG22	1.77	0.47
1:A:225:GLN:C	1:A:227:VAL:H	2.21	0.47
1:B:240:SER:O	1:B:241:ALA:CB	2.62	0.47
1:A:232:GLU:CA	1:A:252:ARG:HD3	2.45	0.47
1:B:134:GLU:CD	1:B:135:ARG:HH21	2.22	0.47
1:B:200:ASN:HB3	11:B:2044:HOH:O	2.14	0.47
1:B:8:ASN:HD22	1:B:8:ASN:HA	1.50	0.47
1:A:148:PHE:HA	1:A:152:HIS:ND1	2.30	0.47
1:B:22:SER:C	1:B:24:PHE:H	2.22	0.47
1:B:64:GLN:HA	1:B:339:LYS:NZ	2.29	0.47
1:B:154:ILE:N	1:B:154:ILE:HD12	2.30	0.46
8:B:1322:PLM:H52	8:B:1323:PLM:H61	1.97	0.46
1:B:68:LEU:HB3	1:B:73:ASN:HD22	1.80	0.46
1:A:231:LYS:HB2	1:A:231:LYS:NZ	2.30	0.46
1:B:150:GLU:O	1:B:154:ILE:CD1	2.63	0.46
1:B:341:GLU:OE1	1:B:342:THR:HG22	2.15	0.46
1:A:22:SER:C	1:A:24:PHE:H	2.22	0.46
1:B:132:ALA:O	1:B:222:CYS:SG	2.70	0.46
1:A:231:LYS:HZ1	1:A:233:ALA:HB3	1.81	0.46
1:A:267:PRO:HA	9:A:1507:HTG:C6'	2.45	0.46
1:B:33:GLU:O	1:B:36:GLN:HB2	2.16	0.46
1:B:336:THR:O	1:B:337:VAL:O	2.33	0.46
1:B:338:SER:CB	1:B:341:GLU:CD	2.88	0.46
1:B:140:CYS:O	1:B:141:LYS:C	2.58	0.45
1:A:223:TYR:C	1:A:225:GLN:N	2.74	0.45

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:321:LEU:HD13	8:A:1322:PLM:H51	1.97	0.45
1:B:64:GLN:HA	1:B:339:LYS:HZ2	1.81	0.45
1:B:237:GLN:HE21	1:B:237:GLN:N	2.14	0.45
1:B:68:LEU:HA	1:B:73:ASN:ND2	2.32	0.45
1:B:244:GLN:OE1	1:B:248:LYS:HE2	2.16	0.45
1:A:28:GLN:HB3	1:A:31:LEU:HD12	1.98	0.45
1:A:74:TYR:CE2	1:A:150:GLU:HG2	2.52	0.45
1:A:137:VAL:O	1:A:141:LYS:HA	2.15	0.45
1:A:134:GLU:HG2	1:A:148:PHE:CD2	2.51	0.45
1:A:337:VAL:C	1:A:338:SER:OG	2.60	0.45
1:B:33:GLU:HB3	1:B:35:TRP:CD1	2.52	0.45
1:B:65:HIS:NE2	1:B:320:THR:OG1	2.49	0.45
1:A:231:LYS:HB2	1:A:231:LYS:HZ1	1.81	0.44
1:B:237:GLN:O	1:B:238:GLN:CB	2.65	0.44
1:A:134:GLU:HG2	1:A:148:PHE:CE2	2.53	0.44
1:A:305:ILE:HG12	9:A:1508:HTG:H6'2	1.99	0.44
1:A:312:GLN:HB2	1:A:332:GLU:HG3	1.99	0.44
1:A:245:LYS:HA	1:A:245:LYS:HZ1	1.81	0.44
1:B:67:LYS:HE2	1:B:316:CYS:SG	2.57	0.44
1:B:240:SER:O	1:B:241:ALA:HB2	2.18	0.44
1:A:322:CYS:HA	8:A:1322:PLM:O1	2.17	0.44
1:A:341:GLU:CD	1:A:342:THR:HG22	2.43	0.44
1:B:244:GLN:HE21	1:B:245:LYS:N	2.16	0.44
1:A:326:ASN:ND2	1:A:328:LEU:CD2	2.81	0.44
1:B:148:PHE:CD1	1:B:152:HIS:HB2	2.52	0.44
1:B:143:MET:CB	1:B:146:PHE:HB3	2.48	0.44
1:B:238:GLN:CG	1:B:241:ALA:H	2.29	0.44
1:A:338:SER:HB2	1:A:341:GLU:CG	2.47	0.43
1:A:301:TYR:HE1	11:A:2020:HOH:O	2.01	0.43
1:B:244:GLN:CD	1:B:248:LYS:HE2	2.43	0.43
1:B:110:CYS:HB3	11:B:2060:HOH:O	2.18	0.43
1:A:137:VAL:O	1:A:137:VAL:HG12	2.17	0.43
1:A:224:GLY:O	1:A:227:VAL:CG1	2.66	0.43
1:A:314:ARG:O	1:A:317:MET:HB3	2.19	0.43
1:B:332:GLU:C	1:B:334:SER:H	2.26	0.43
1:A:338:SER:OG	1:A:341:GLU:OE2	2.34	0.43
1:A:238:GLN:HA	1:A:242:THR:CA	2.48	0.43
1:B:152:HIS:O	1:B:155:MET:HB2	2.20	0.42
1:A:8:ASN:HD22	1:A:8:ASN:HA	1.50	0.42
1:A:328:LEU:O	1:A:329:GLY:C	2.62	0.42
1:A:330:ASP:OD2	1:B:100:HIS:CE1	2.72	0.42

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:227:VAL:CG1	1:A:228:PHE:H	2.32	0.42
1:A:238:GLN:HG2	1:A:239:GLU:N	2.34	0.42
1:B:224:GLY:O	1:B:227:VAL:HG12	2.19	0.42
1:A:241:ALA:C	1:A:243:THR:N	2.78	0.42
1:B:199:ASN:HB3	11:B:2005:HOH:O	2.19	0.42
1:A:124:ALA:HB2	11:A:2030:HOH:O	2.20	0.42
1:A:239:GLU:CG	1:A:248:LYS:HZ2	2.31	0.41
1:B:146:PHE:CE1	1:B:152:HIS:NE2	2.86	0.41
1:A:345:VAL:CG1	1:A:347:PRO:HD3	2.33	0.41
7:B:1296:RET:H201	11:B:2018:HOH:O	2.20	0.41
1:B:135:ARG:NH2	1:B:138:VAL:HG21	2.35	0.41
1:B:330:ASP:C	1:B:331:ASP:O	2.63	0.41
1:B:239:GLU:HB3	1:B:244:GLN:CD	2.45	0.41
1:A:347:PRO:HB2	1:A:348:ALA:H	1.58	0.41
1:A:260:ALA:HB2	8:A:1410:PLM:HE2	2.02	0.41
1:A:337:VAL:CG2	1:A:338:SER:N	2.83	0.41
1:B:68:LEU:CA	1:B:73:ASN:HD22	2.34	0.41
1:B:224:GLY:O	1:B:227:VAL:HB	2.21	0.41
1:B:322:CYS:O	1:B:324:GLY:N	2.53	0.41
1:B:64:GLN:CG	1:B:339:LYS:HE2	2.29	0.41
1:B:283:PHE:HA	9:B:1506:HTG:H61	2.03	0.41
1:A:223:TYR:C	1:A:225:GLN:H	2.28	0.41
1:A:227:VAL:O	1:A:231:LYS:HG2	2.21	0.41
1:B:27:PRO:HB3	1:B:29:TYR:CE2	2.56	0.41
1:B:239:GLU:OE1	1:B:239:GLU:HA	2.20	0.41
1:B:226:LEU:N	1:B:226:LEU:CD1	2.83	0.41
1:B:244:GLN:NE2	1:B:245:LYS:N	2.68	0.40
1:B:338:SER:HB2	1:B:341:GLU:CD	2.46	0.40
1:A:232:GLU:CB	1:A:252:ARG:HD3	2.51	0.40
1:A:236:GLN:O	1:A:245:LYS:HD2	2.21	0.40
1:B:332:GLU:O	1:B:333:ALA:CB	2.68	0.40
1:B:137:VAL:HG12	1:B:137:VAL:O	2.21	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:344:GLN:NE2	9:B:1509:HTG:O6[4_564]	2.19	0.01

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	347/349 (99%)	264 (76%)	57 (16%)	26 (8%)	1	2
1	B	347/349 (99%)	254 (73%)	69 (20%)	24 (7%)	1	2
All	All	694/698 (99%)	518 (75%)	126 (18%)	50 (7%)	1	2

All (50) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	240	SER
1	A	241	ALA
1	A	323	CYS
1	A	328	LEU
1	A	341	GLU
1	B	229	THR
1	B	240	SER
1	B	241	ALA
1	B	281	SER
1	B	331	ASP
1	B	337	VAL
1	B	341	GLU
1	B	347	PRO
1	A	195	HIS
1	A	229	THR
1	A	279	GLN
1	A	281	SER
1	A	301	TYR
1	A	324	GLY
1	A	329	GLY
1	A	335	THR
1	A	347	PRO
1	B	195	HIS
1	B	237	GLN
1	B	323	CYS

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	B	324	GLY
1	A	87	VAL
1	A	93	THR
1	A	94	THR
1	A	115	PHE
1	A	119	LEU
1	A	167	CYS
1	A	338	SER
1	B	236	GLN
1	B	238	GLN
1	A	89	GLY
1	A	201	GLU
1	B	51	GLY
1	B	161	TRP
1	B	260	ALA
1	A	141	LYS
1	A	145	ASN
1	B	262	LEU
1	B	87	VAL
1	B	218	VAL
1	B	329	GLY
1	B	133	ILE
1	A	337	VAL
1	B	141	LYS
1	B	305	ILE

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	296/296 (100%)	274 (93%)	22 (7%)	13	37
1	B	296/296 (100%)	269 (91%)	27 (9%)	9	28
All	All	592/592 (100%)	543 (92%)	49 (8%)	10	32

All (49) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	8	ASN
1	A	76	LEU
1	A	78	ASN
1	A	119	LEU
1	A	131	LEU
1	A	135	ARG
1	A	165	LEU
1	A	179	ILE
1	A	191	TYR
1	A	205	ILE
1	A	213	ILE
1	A	216	LEU
1	A	231	LYS
1	A	237	GLN
1	A	244	GLN
1	A	245	LYS
1	A	266	LEU
1	A	302	ASN
1	A	332	GLU
1	A	336	THR
1	A	337	VAL
1	A	343	SER
1	B	8	ASN
1	B	50	LEU
1	B	70	THR
1	B	104	VAL
1	B	131	LEU
1	B	133	ILE
1	B	135	ARG
1	B	143	MET
1	B	148	PHE
1	B	185	CYS
1	B	207	MET
1	B	208	PHE
1	B	220	PHE
1	B	229	THR
1	B	236	GLN
1	B	237	GLN
1	B	244	GLN
1	B	245	LYS
1	B	249	GLU
1	B	263	ILE
1	B	266	LEU

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	B	283	PHE
1	B	328	LEU
1	B	330	ASP
1	B	332	GLU
1	B	337	VAL
1	B	341	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (18) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	8	ASN
1	A	64	GLN
1	A	73	ASN
1	A	100	HIS
1	A	237	GLN
1	A	302	ASN
1	A	326	ASN
1	B	8	ASN
1	B	64	GLN
1	B	73	ASN
1	B	111	ASN
1	B	151	ASN
1	B	184	GLN
1	B	236	GLN
1	B	237	GLN
1	B	244	GLN
1	B	302	ASN
1	B	326	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 ⓘ

11 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	0.63	0	17,19,21	0.85	0
2	NAG	C	2	2	14,14,15	0.77	0	17,19,21	0.93	1 (5%)
2	MAN	C	3	2	11,11,12	0.79	0	15,15,17	0.72	0
3	NAG	D	1	1,3	14,14,15	0.53	0	17,19,21	0.84	1 (5%)
3	NAG	D	2	3	14,14,15	0.63	0	17,19,21	1.00	1 (5%)
4	NAG	E	1	1,4	14,14,15	0.60	0	17,19,21	0.67	0
4	NAG	E	2	4	14,14,15	0.65	0	17,19,21	0.96	1 (5%)
4	BMA	E	3	4	11,11,12	0.81	0	15,15,17	0.60	0
4	BMA	E	4	4	11,11,12	0.83	0	15,15,17	0.81	1 (6%)
3	NAG	F	1	1,3	14,14,15	0.64	0	17,19,21	0.84	1 (5%)
3	NAG	F	2	3	14,14,15	0.66	0	17,19,21	0.68	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	-	2/6/23/26	0/1/1/1
2	NAG	C	2	2	-	4/6/23/26	0/1/1/1
2	MAN	C	3	2	-	2/2/19/22	0/1/1/1
3	NAG	D	1	1,3	-	4/6/23/26	0/1/1/1
3	NAG	D	2	3	-	0/6/23/26	0/1/1/1
4	NAG	E	1	1,4	-	2/6/23/26	0/1/1/1
4	NAG	E	2	4	-	2/6/23/26	0/1/1/1
4	BMA	E	3	4	-	2/2/19/22	0/1/1/1
4	BMA	E	4	4	-	0/2/19/22	0/1/1/1
3	NAG	F	1	1,3	-	2/6/23/26	0/1/1/1
3	NAG	F	2	3	-	4/6/23/26	0/1/1/1

There are no bond length outliers.

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	2	NAG	C2-N2-C7	-3.20	118.61	122.90
4	E	2	NAG	C2-N2-C7	-2.81	119.13	122.90
3	D	1	NAG	C2-N2-C7	-2.77	119.18	122.90
4	E	4	BMA	C1-O5-C5	2.32	115.30	112.19
2	C	2	NAG	C8-C7-N2	2.24	119.83	116.12
3	F	1	NAG	C2-N2-C7	-2.20	119.95	122.90

There are no chirality outliers.

All (24) torsion outliers are listed below:

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

There are no ring outliers.

3 monomers are involved in 2 short contacts:

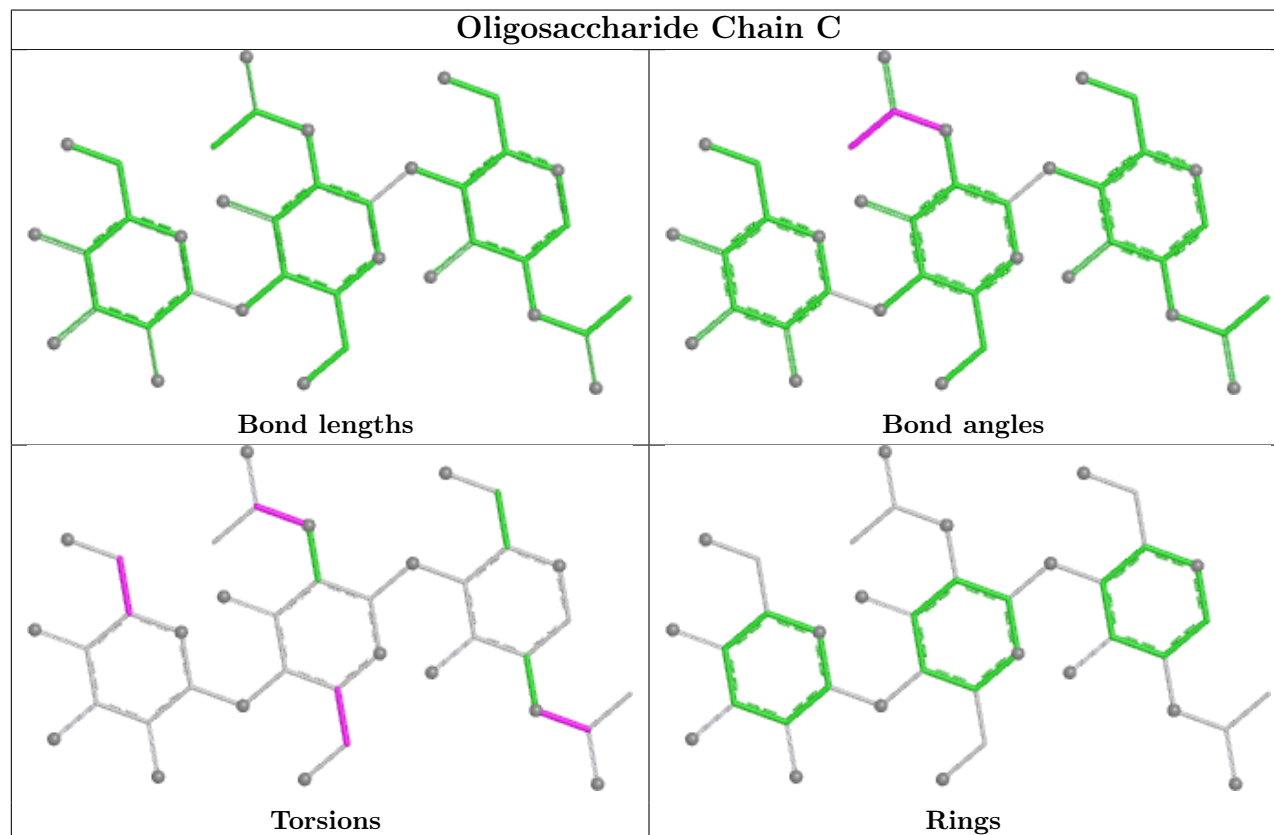
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	2	NAG	1	0

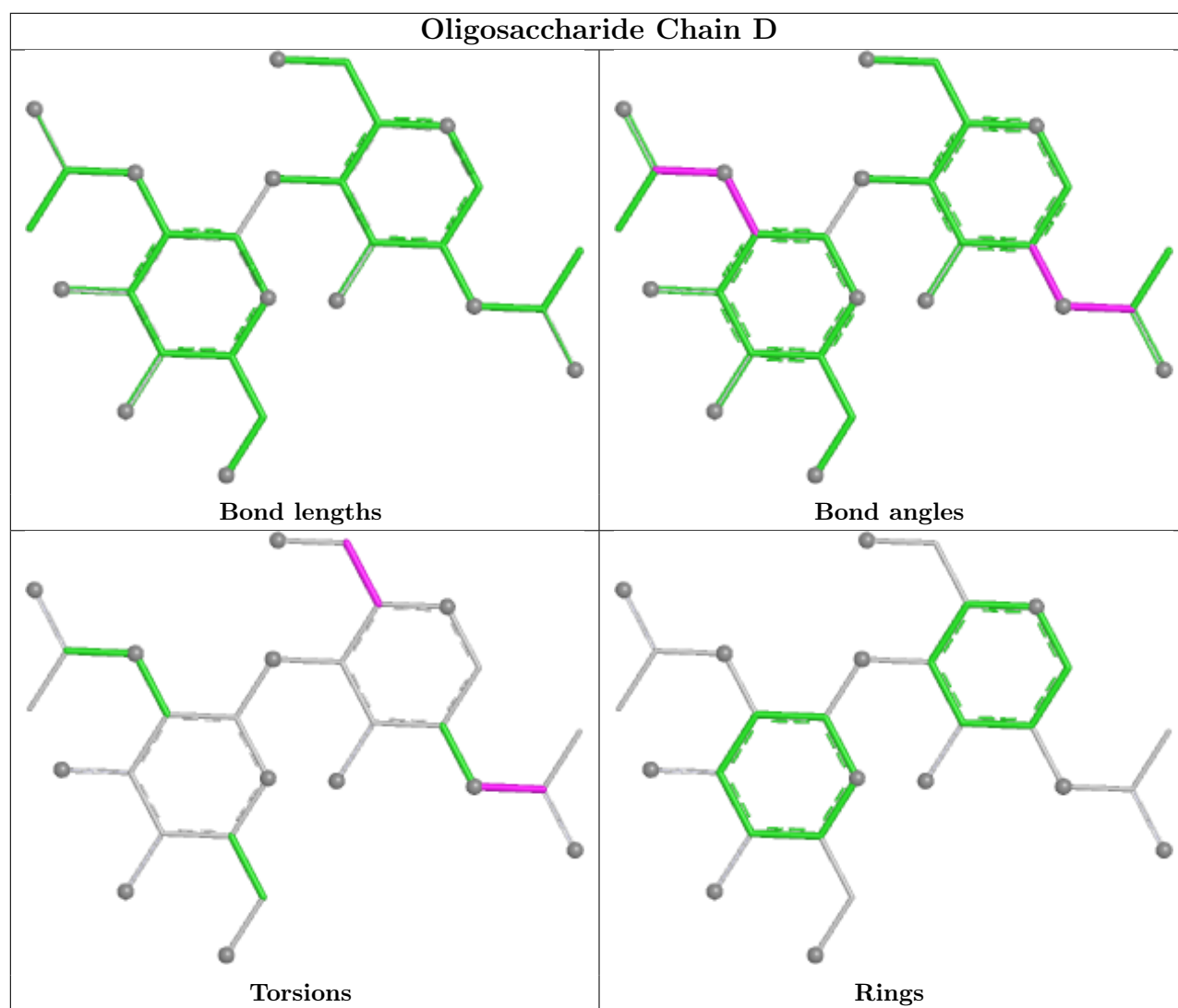
Continued on next page...

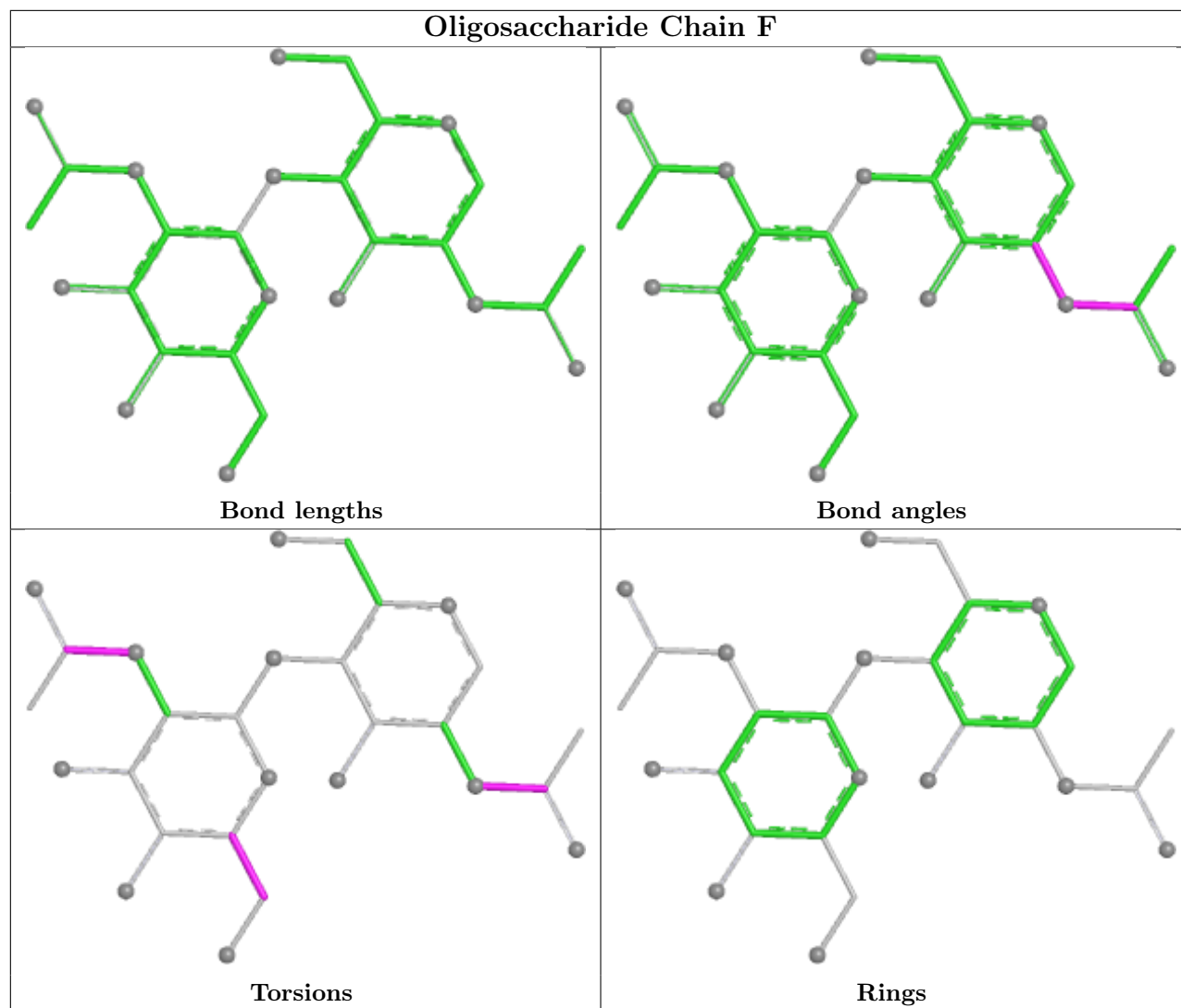
Continued from previous page...

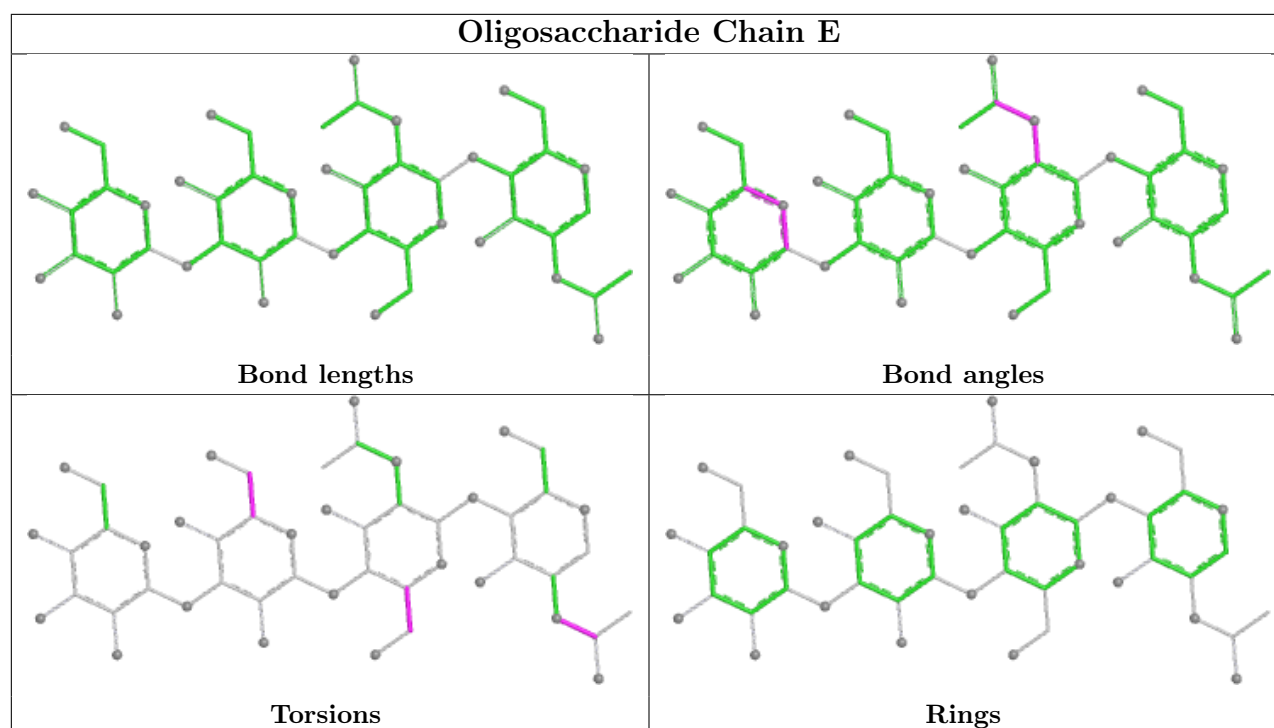
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	F	1	NAG	1	0
2	C	1	NAG	1	0

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 [i](#)

Of 26 ligands modelled in this entry, 13 are monoatomic - leaving 13 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	$\# Z > 2$
8	PLM	A	1323	1	16,16,17	1.00	1 (6%)	15,15,17	0.52	0
8	PLM	B	1323	1	16,16,17	1.01	1 (6%)	15,15,17	0.45	0
9	HTG	B	1506	-	19,19,19	3.06	8 (42%)	23,24,24	1.76	2 (8%)
8	PLM	A	1322	1	16,16,17	0.99	1 (6%)	15,15,17	0.61	0
7	RET	A	1296	1	20,20,21	2.56	3 (15%)	27,27,28	2.03	7 (25%)
8	PLM	B	1322	1	16,16,17	1.04	1 (6%)	15,15,17	0.43	0
9	HTG	B	1509	-	19,19,19	2.80	8 (42%)	23,24,24	1.98	1 (4%)
8	PLM	A	1410	-	15,15,17	2.07	3 (20%)	14,14,17	3.42	7 (50%)
9	HTG	A	1507	-	19,19,19	3.25	8 (42%)	23,24,24	1.99	2 (8%)
8	PLM	B	1407	-	15,15,17	2.01	3 (20%)	14,14,17	3.41	8 (57%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
9	HTG	A	1508	-	19,19,19	2.48	8 (42%)	23,24,24	2.29	2 (8%)
7	RET	B	1296	1	20,20,21	2.30	4 (20%)	27,27,28	1.90	6 (22%)
10	HTO	B	1401	-	9,9,9	1.81	2 (22%)	10,10,10	1.10	1 (10%)

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
8	PLM	A	1323	1	-	10/14/14/15	-
8	PLM	B	1323	1	-	9/14/14/15	-
9	HTG	B	1506	-	-	2/10/30/30	0/1/1/1
8	PLM	A	1322	1	-	11/14/14/15	-
7	RET	A	1296	1	-	4/13/30/31	0/1/1/1
8	PLM	B	1322	1	-	11/14/14/15	-
9	HTG	B	1509	-	-	5/10/30/30	0/1/1/1
8	PLM	A	1410	-	-	5/13/13/15	-
9	HTG	A	1507	-	-	5/10/30/30	0/1/1/1
8	PLM	B	1407	-	-	5/13/13/15	-
9	HTG	A	1508	-	-	4/10/30/30	0/1/1/1
7	RET	B	1296	1	-	5/13/30/31	0/1/1/1
10	HTO	B	1401	-	-	2/10/10/10	-

All (51) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
9	A	1507	HTG	O5-C1	9.41	1.57	1.42
9	B	1506	HTG	O5-C1	8.98	1.56	1.42
9	B	1509	HTG	O5-C1	7.97	1.54	1.42
7	A	1296	RET	C1-C6	7.82	1.63	1.53
9	A	1508	HTG	O5-C1	7.05	1.53	1.42
7	A	1296	RET	C5-C6	6.62	1.45	1.34
7	B	1296	RET	C1-C6	6.17	1.61	1.53
7	B	1296	RET	C5-C6	5.75	1.44	1.34
8	A	1410	PLM	CC-CB	-5.50	1.24	1.51
8	B	1407	PLM	CC-CB	-5.27	1.25	1.51
9	A	1507	HTG	C1-S1	5.01	1.89	1.80
9	A	1507	HTG	C1-C2	4.77	1.61	1.53

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
9	B	1509	HTG	C4-C5	4.61	1.62	1.53
9	B	1509	HTG	C1-C2	4.35	1.60	1.53
10	B	1401	HTO	C3-C2	4.25	1.63	1.53
9	B	1506	HTG	C1-S1	4.09	1.87	1.80
9	B	1506	HTG	C3-C2	4.04	1.62	1.52
8	A	1410	PLM	CB-CA	-3.84	1.32	1.51
8	A	1322	PLM	O1-C1	-3.81	1.22	1.42
9	B	1506	HTG	C4-C5	3.80	1.61	1.53
9	A	1507	HTG	C3-C2	3.78	1.62	1.52
9	A	1508	HTG	C4-C5	3.74	1.61	1.53
8	B	1322	PLM	O1-C1	-3.74	1.23	1.42
8	A	1323	PLM	O1-C1	-3.70	1.23	1.42
9	A	1507	HTG	C4-C5	3.62	1.60	1.53
8	B	1407	PLM	CB-CA	-3.60	1.33	1.51
9	B	1506	HTG	C1-C2	3.57	1.59	1.53
9	A	1507	HTG	O5-C5	3.48	1.52	1.44
8	B	1323	PLM	O1-C1	-3.45	1.24	1.42
8	A	1410	PLM	CD-CC	-3.42	1.34	1.51
9	B	1506	HTG	O5-C5	3.39	1.52	1.44
9	A	1507	HTG	C4-C3	3.31	1.60	1.52
9	B	1506	HTG	C4-C3	3.31	1.60	1.52
9	B	1509	HTG	O5-C5	3.30	1.52	1.44
7	B	1296	RET	C14-C13	3.29	1.36	1.33
9	A	1508	HTG	C1-C2	3.29	1.59	1.53
9	B	1509	HTG	C3-C2	3.21	1.60	1.52
9	A	1508	HTG	O5-C5	3.20	1.52	1.44
8	B	1407	PLM	CD-CC	-3.20	1.35	1.51
9	B	1509	HTG	C4-C3	3.12	1.60	1.52
9	A	1508	HTG	C3-C2	3.09	1.60	1.52
9	A	1508	HTG	C4-C3	3.07	1.60	1.52
9	B	1506	HTG	C1'-S1	2.89	1.85	1.81
9	A	1507	HTG	C1'-S1	2.83	1.85	1.81
7	B	1296	RET	C17-C1	2.66	1.58	1.53
9	B	1509	HTG	C1-S1	2.59	1.85	1.80
7	A	1296	RET	C15-C14	-2.33	1.40	1.49
9	B	1509	HTG	C6-C5	2.27	1.59	1.51
9	A	1508	HTG	C1'-S1	2.19	1.84	1.81
9	A	1508	HTG	C1-S1	2.13	1.84	1.80
10	B	1401	HTO	C5-C4	2.07	1.60	1.52

All (36) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
9	A	1508	HTG	C1'-S1-C1	10.12	122.30	100.45
9	B	1509	HTG	C1'-S1-C1	8.41	118.62	100.45
9	A	1507	HTG	C1'-S1-C1	8.35	118.47	100.45
8	A	1410	PLM	CD-CC-CB	8.34	156.51	114.37
8	B	1407	PLM	CD-CC-CB	8.21	155.85	114.37
8	A	1410	PLM	CA-C9-C8	7.57	152.62	114.37
8	B	1407	PLM	CA-C9-C8	7.41	151.83	114.37
9	B	1506	HTG	C1'-S1-C1	7.15	115.88	100.45
7	A	1296	RET	C18-C5-C6	5.23	130.19	124.48
7	A	1296	RET	C11-C10-C9	-4.93	120.36	127.28
7	B	1296	RET	C18-C5-C6	4.69	129.60	124.48
7	B	1296	RET	C11-C10-C9	-3.81	121.94	127.28
7	B	1296	RET	C8-C9-C10	-3.67	113.23	119.01
8	B	1407	PLM	CC-CB-CA	3.21	130.59	114.37
8	A	1410	PLM	CC-CB-CA	3.18	130.46	114.37
7	B	1296	RET	C8-C7-C6	3.13	135.37	127.00
7	A	1296	RET	C8-C9-C10	-2.96	114.36	119.01
7	B	1296	RET	C2-C1-C6	2.70	114.36	110.44
7	B	1296	RET	C19-C9-C10	2.66	127.12	122.82
7	A	1296	RET	C8-C7-C6	2.65	134.09	127.00
8	B	1407	PLM	CE-CD-CC	2.58	127.40	114.37
7	A	1296	RET	C2-C1-C6	2.54	114.13	110.44
8	B	1407	PLM	C5-C4-C3	2.50	127.00	114.37
8	A	1410	PLM	C5-C4-C3	2.50	126.99	114.37
9	A	1507	HTG	O5-C1-C2	-2.38	107.03	110.32
7	A	1296	RET	C19-C9-C10	2.33	126.59	122.82
8	A	1410	PLM	CE-CD-CC	2.27	125.86	114.37
7	A	1296	RET	C18-C5-C4	-2.26	108.77	113.60
8	B	1407	PLM	C9-C8-C7	2.20	125.50	114.37
8	B	1407	PLM	CG-CF-CE	2.20	128.21	113.36
10	B	1401	HTO	O3-C3-C4	-2.19	104.17	109.14
8	A	1410	PLM	CG-CF-CE	2.16	127.98	113.36
8	A	1410	PLM	CF-CE-CD	-2.14	96.51	115.25
8	B	1407	PLM	CF-CE-CD	-2.12	96.67	115.25
9	B	1506	HTG	O5-C5-C6	2.12	111.68	106.44
9	A	1508	HTG	O5-C5-C6	2.05	111.52	106.44

There are no chirality outliers.

All (78) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
7	A	1296	RET	C1-C6-C7-C8
7	A	1296	RET	C5-C6-C7-C8

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms
7	B	1296	RET	C5-C6-C7-C8
9	B	1509	HTG	O5-C5-C6-O6
8	B	1407	PLM	CA-CB-CC-CD
9	B	1509	HTG	C4-C5-C6-O6
9	A	1507	HTG	O5-C5-C6-O6
9	A	1508	HTG	C1'-C2'-C3'-C4'
7	B	1296	RET	C11-C10-C9-C8
9	B	1506	HTG	C1'-C2'-C3'-C4'
8	A	1322	PLM	CC-CD-CE-CF
8	A	1410	PLM	C5-C6-C7-C8
8	A	1323	PLM	C4-C5-C6-C7
8	B	1323	PLM	CA-CB-CC-CD
8	A	1323	PLM	C1-C2-C3-C4
8	A	1323	PLM	CB-CC-CD-CE
8	A	1410	PLM	C7-C8-C9-CA
8	B	1323	PLM	C6-C7-C8-C9
8	B	1323	PLM	C9-CA-CB-CC
8	B	1322	PLM	CB-CC-CD-CE
7	B	1296	RET	C1-C6-C7-C8
8	A	1323	PLM	C7-C8-C9-CA
8	B	1323	PLM	C3-C4-C5-C6
7	A	1296	RET	C10-C11-C12-C13
8	B	1322	PLM	C9-CA-CB-CC
9	A	1507	HTG	C2'-C3'-C4'-C5'
9	B	1509	HTG	C2'-C3'-C4'-C5'
9	B	1509	HTG	C3'-C4'-C5'-C6'
8	A	1323	PLM	C3-C4-C5-C6
8	A	1322	PLM	C9-CA-CB-CC
8	B	1323	PLM	CB-CC-CD-CE
8	B	1323	PLM	C8-C9-CA-CB
8	B	1322	PLM	C8-C9-CA-CB
9	A	1508	HTG	C2'-C3'-C4'-C5'
8	B	1322	PLM	C5-C6-C7-C8
9	B	1506	HTG	C2'-C3'-C4'-C5'
8	A	1322	PLM	C5-C6-C7-C8
8	A	1410	PLM	C9-CA-CB-CC
8	B	1407	PLM	C2-C3-C4-C5
8	A	1322	PLM	C7-C8-C9-CA
8	A	1322	PLM	CB-CC-CD-CE
8	B	1322	PLM	C7-C8-C9-CA
9	A	1508	HTG	C3'-C4'-C5'-C6'
7	B	1296	RET	C10-C11-C12-C13

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms
8	A	1322	PLM	C1-C2-C3-C4
8	A	1322	PLM	C6-C7-C8-C9
8	B	1323	PLM	C4-C5-C6-C7
8	A	1322	PLM	C3-C4-C5-C6
8	B	1322	PLM	CC-CD-CE-CF
8	A	1323	PLM	CA-CB-CC-CD
8	A	1410	PLM	C2-C3-C4-C5
10	B	1401	HTO	C4-C5-C6-C7
8	A	1410	PLM	CA-CB-CC-CD
8	B	1407	PLM	C5-C6-C7-C8
8	B	1322	PLM	C4-C5-C6-C7
9	A	1507	HTG	C4'-C5'-C6'-C7'
8	A	1322	PLM	C4-C5-C6-C7
9	A	1508	HTG	C4'-C5'-C6'-C7'
8	A	1323	PLM	CC-CD-CE-CF
8	A	1322	PLM	C8-C9-CA-CB
8	A	1323	PLM	C2-C3-C4-C5
8	B	1407	PLM	C3-C4-C5-C6
10	B	1401	HTO	C1-C2-C3-O3
8	B	1323	PLM	C5-C6-C7-C8
8	B	1322	PLM	C3-C4-C5-C6
8	A	1323	PLM	C8-C9-CA-CB
8	B	1322	PLM	C1-C2-C3-C4
8	A	1322	PLM	O1-C1-C2-C3
7	A	1296	RET	C11-C10-C9-C8
8	B	1407	PLM	C7-C8-C9-CA
9	A	1507	HTG	C4-C5-C6-O6
8	A	1323	PLM	C6-C7-C8-C9
8	B	1323	PLM	O1-C1-C2-C3
8	B	1322	PLM	C2-C3-C4-C5
7	B	1296	RET	C6-C7-C8-C9
9	A	1507	HTG	C3'-C4'-C5'-C6'
8	B	1322	PLM	CA-CB-CC-CD
9	B	1509	HTG	C4'-C5'-C6'-C7'

There are no ring outliers.

11 monomers are involved in 20 short contacts:

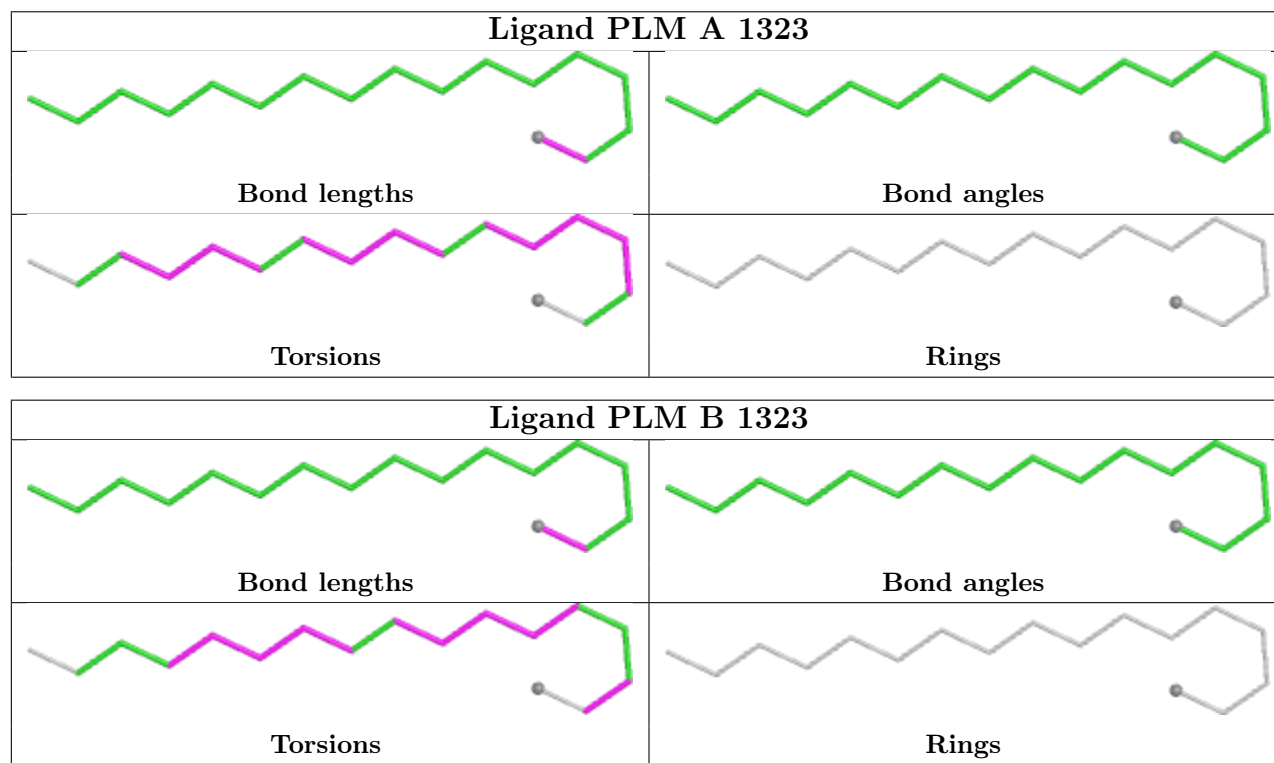
Mol	Chain	Res	Type	Clashes	Symm-Clashes
8	B	1323	PLM	2	0
9	B	1506	HTG	1	0
8	A	1322	PLM	6	0

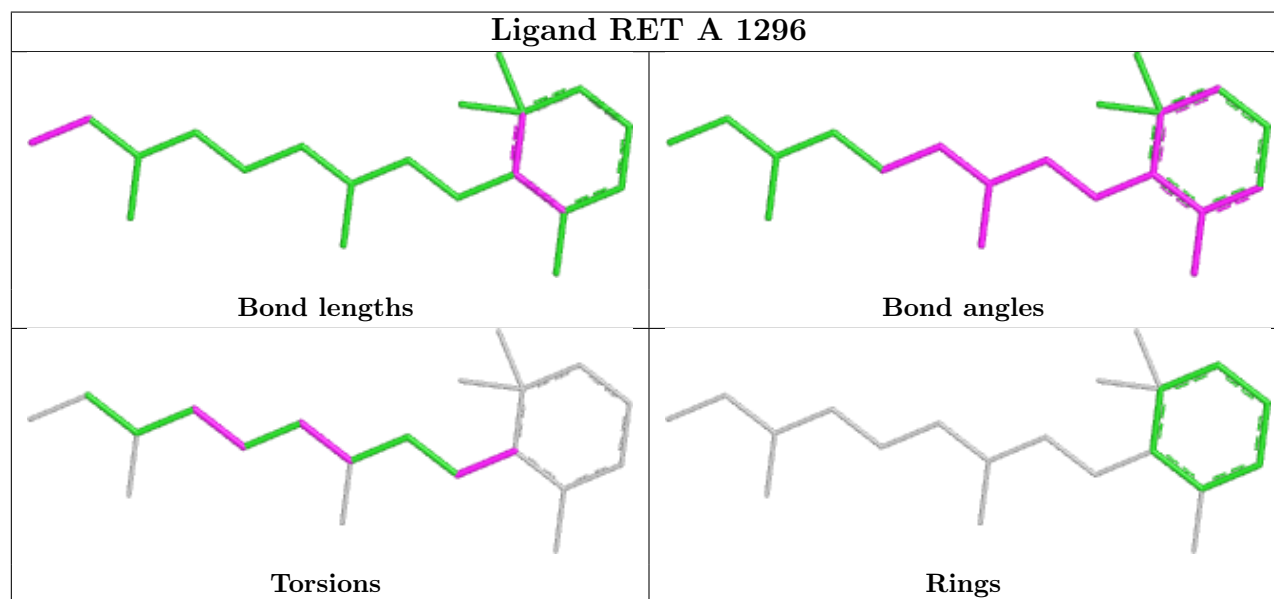
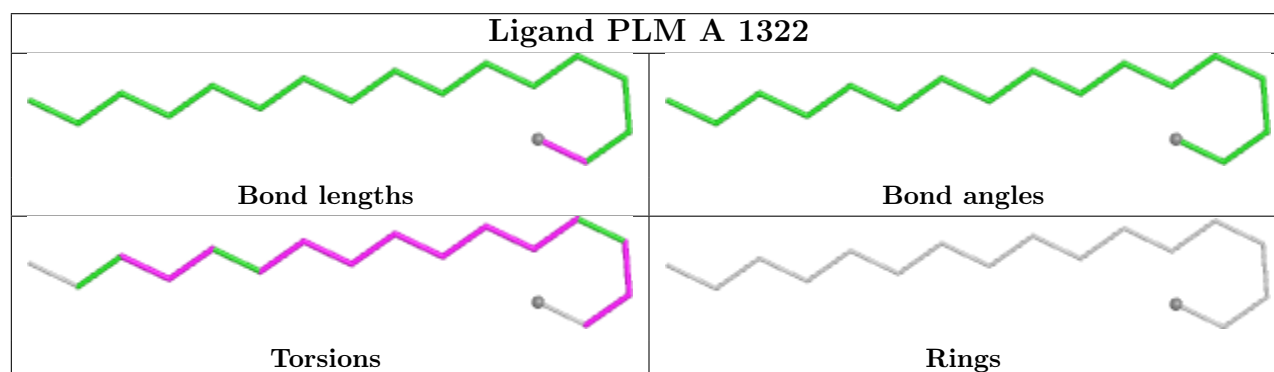
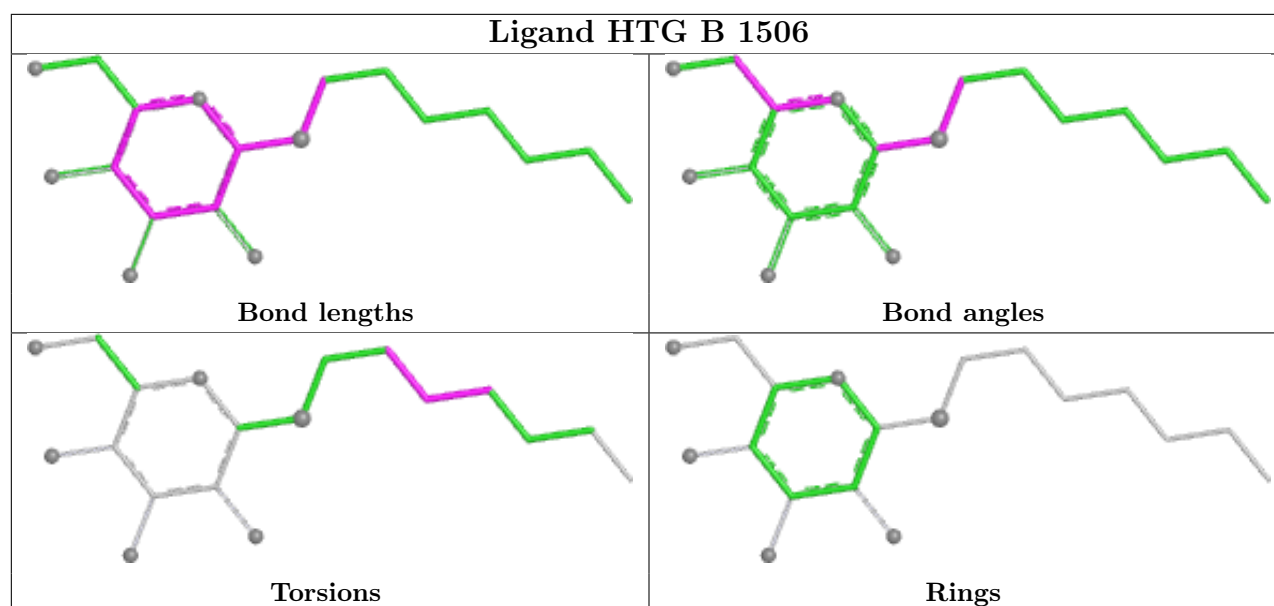
Continued on next page...

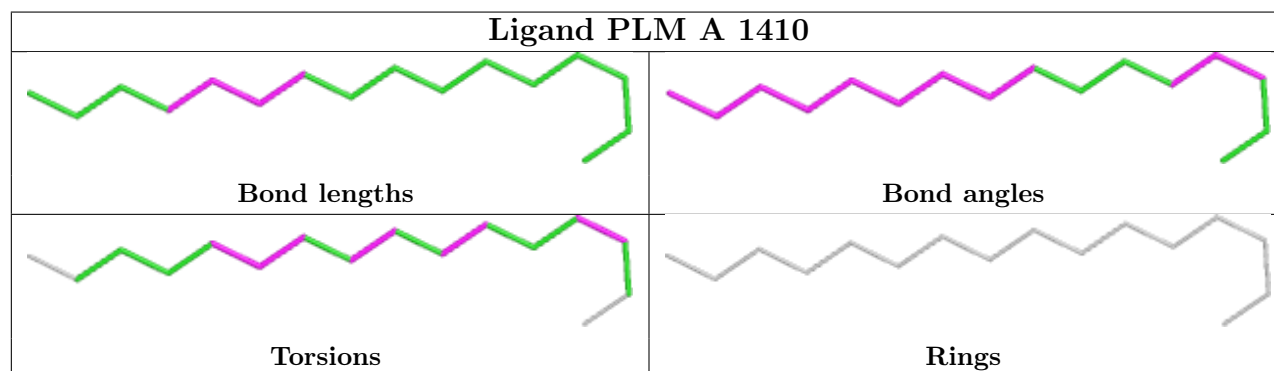
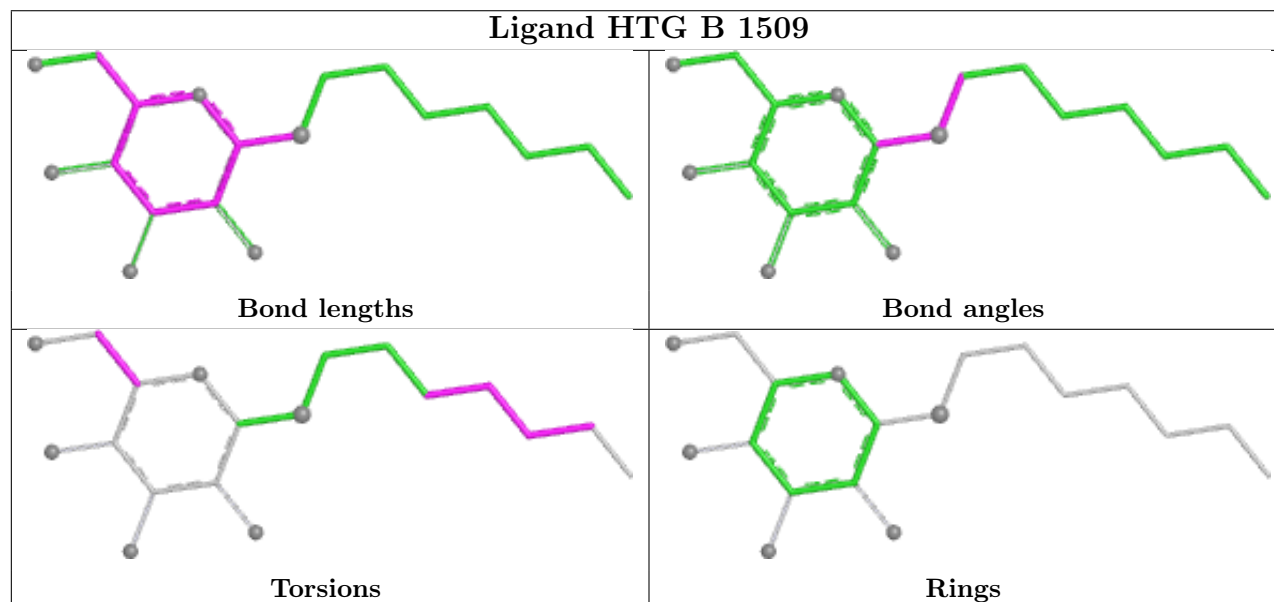
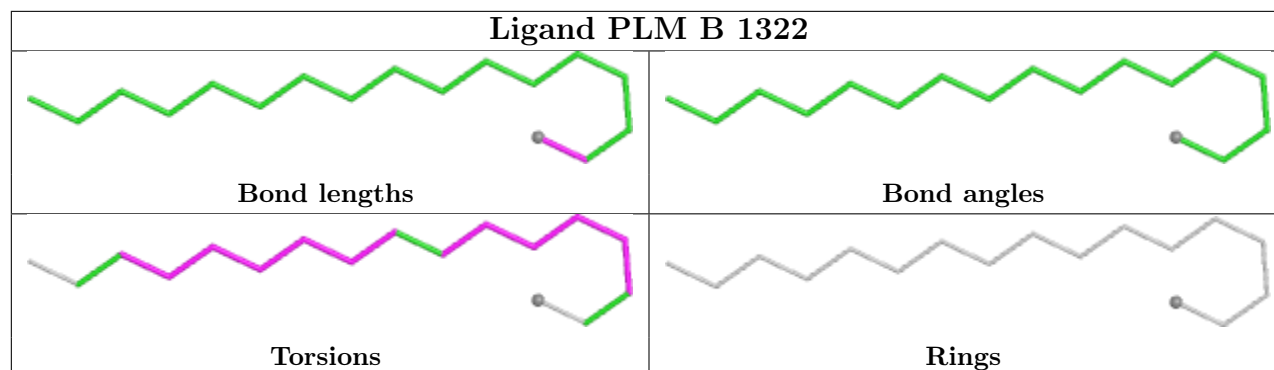
Continued from previous page...

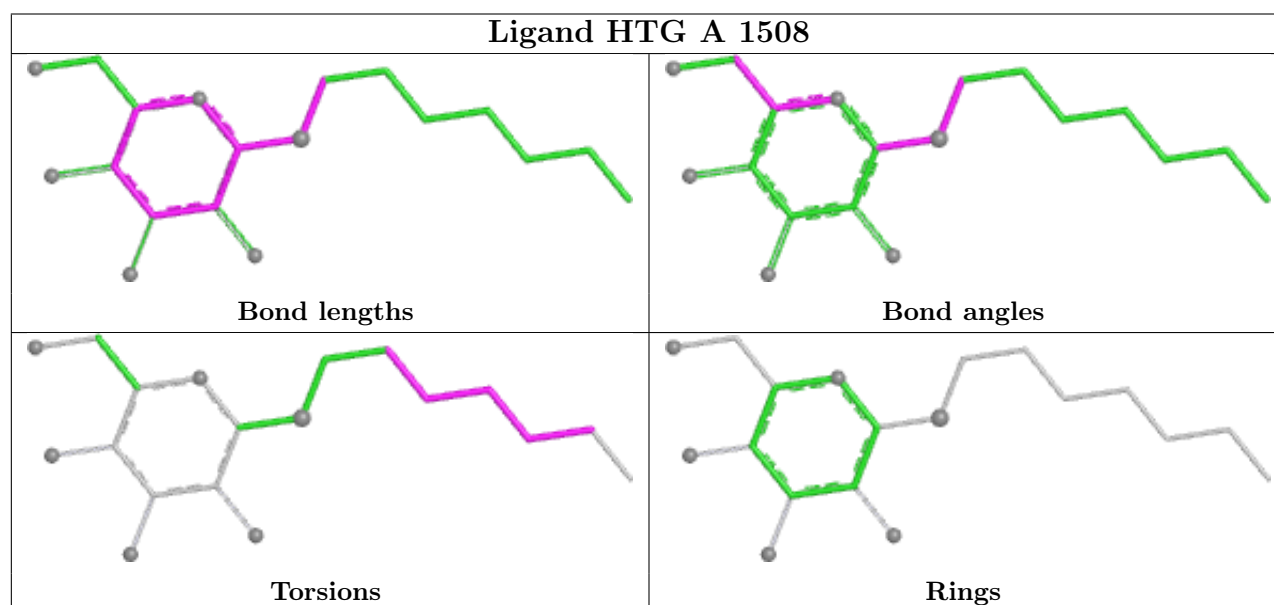
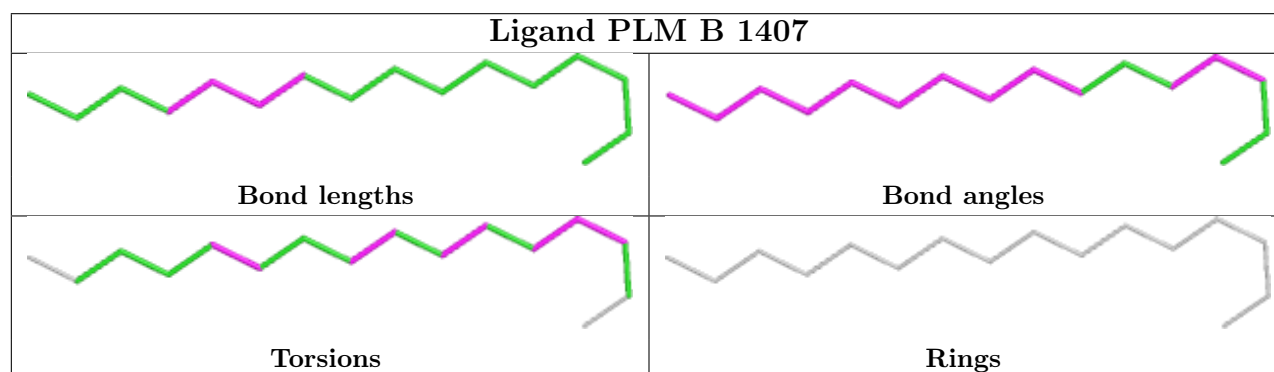
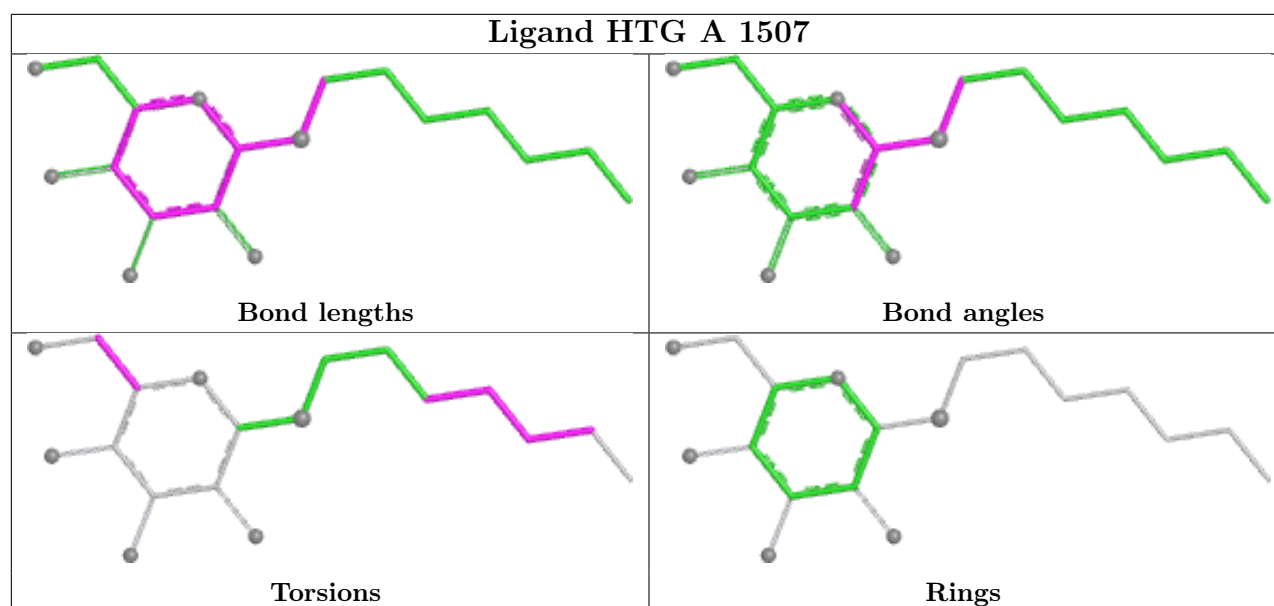
Mol	Chain	Res	Type	Clashes	Symm-Clashes
8	B	1322	PLM	2	0
9	B	1509	HTG	3	1
8	A	1410	PLM	1	0
9	A	1507	HTG	2	0
8	B	1407	PLM	1	0
9	A	1508	HTG	1	0
7	B	1296	RET	1	0
10	B	1401	HTO	1	0

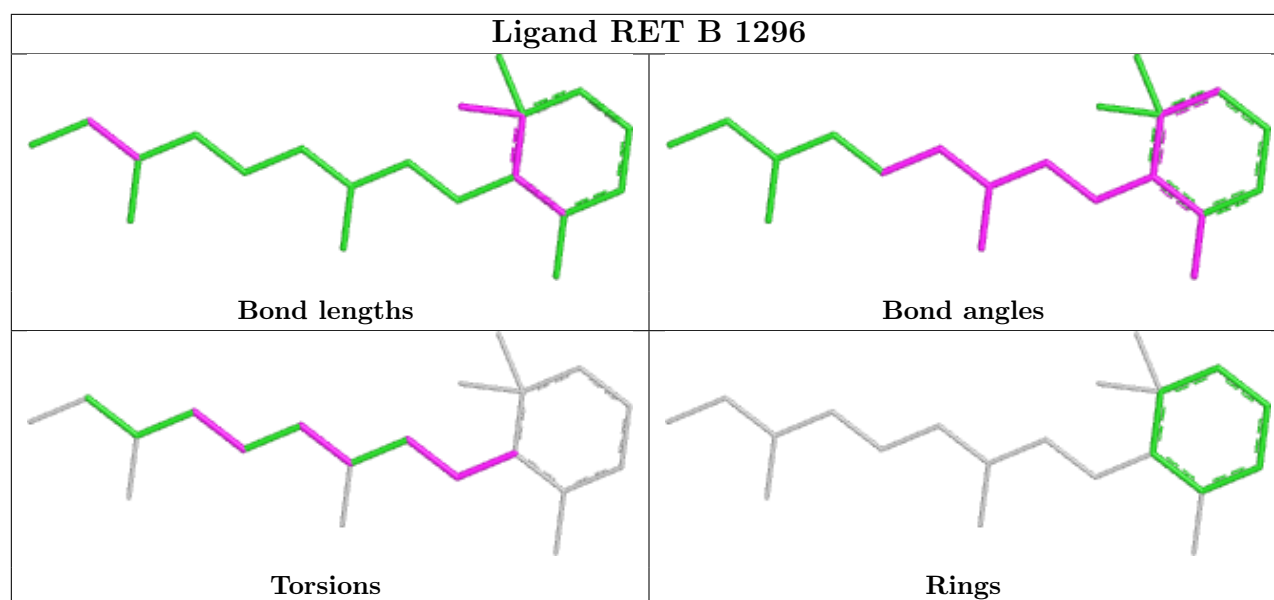
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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











5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2		OWAB(Å ²)	Q<0.9
1	A	348/349 (99%)	0.21	15 (4%)	40 31	19, 26, 135, 149	228 (65%)
1	B	348/349 (99%)	0.38	24 (6%)	23 16	19, 27, 154, 163	228 (65%)
All	All	696/698 (99%)	0.30	39 (5%)	30 23	19, 26, 147, 163	456 (65%)

All (39) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	146	PHE	7.0
1	B	229	THR	5.4
1	A	328	LEU	5.1
1	B	142	PRO	5.0
1	B	344	GLN	4.9
1	B	238	GLN	4.2
1	B	133	ILE	4.2
1	B	228	PHE	3.8
1	B	323	CYS	3.7
1	B	232	GLU	3.7
1	B	263	ILE	3.7
1	B	148	PHE	3.5
1	B	107	PRO	3.4
1	B	320	THR	3.1
1	B	21	ARG	3.1
1	A	233	ALA	3.0
1	B	348	ALA	3.0
1	A	14	SER	3.0
1	A	147	ARG	2.8
1	A	333	ALA	2.8
1	B	334	SER	2.7
1	B	234	ALA	2.6
1	A	32	ALA	2.6
1	B	116	PHE	2.5

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	A	324	GLY	2.5
1	B	137	VAL	2.4
1	B	260	ALA	2.4
1	A	282	ASP	2.4
1	A	105	PHE	2.3
1	A	314	ARG	2.3
1	A	348	ALA	2.3
1	B	164	ALA	2.2
1	B	329	GLY	2.2
1	A	96	TYR	2.2
1	A	33	GLU	2.2
1	B	342	THR	2.1
1	B	295	ALA	2.1
1	A	243	THR	2.1
1	B	108	THR	2.0

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

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

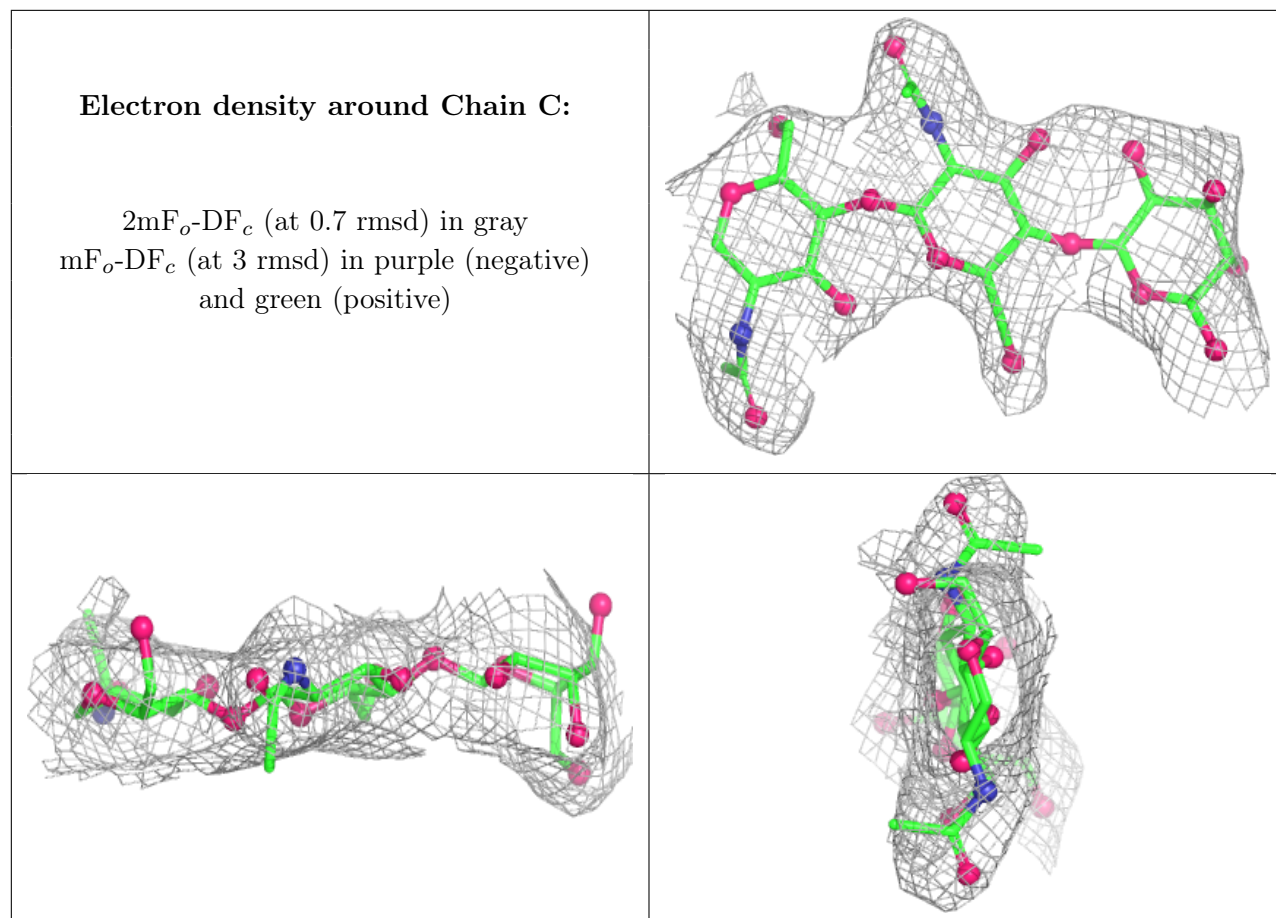
6.3 Carbohydrates [i](#)

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	NAG	F	2	14/15	0.95	0.08	84,89,91,92	0
4	BMA	E	3	11/12	0.95	0.10	104,108,112,116	0
3	NAG	D	1	14/15	0.96	0.07	70,75,80,84	0
3	NAG	F	1	14/15	0.96	0.08	65,69,71,77	0
3	NAG	D	2	14/15	0.97	0.07	91,94,99,101	0
4	NAG	E	1	14/15	0.97	0.07	63,64,69,75	0
4	NAG	E	2	14/15	0.97	0.09	77,83,90,97	0
2	MAN	C	3	11/12	0.97	0.05	94,97,97,99	0
2	NAG	C	1	14/15	0.98	0.06	66,68,74,75	0
2	NAG	C	2	14/15	0.98	0.06	79,82,87,92	0
4	BMA	E	4	11/12	0.98	0.10	120,122,122,124	0

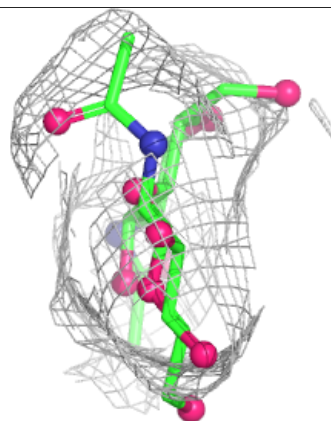
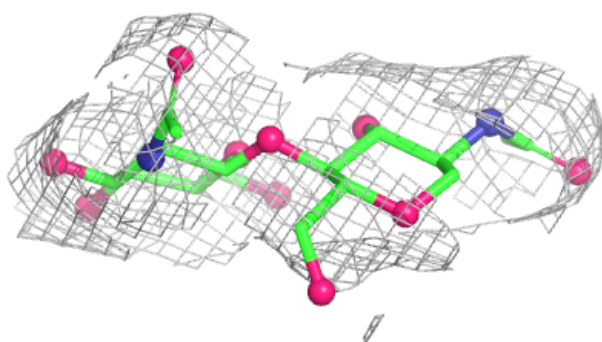
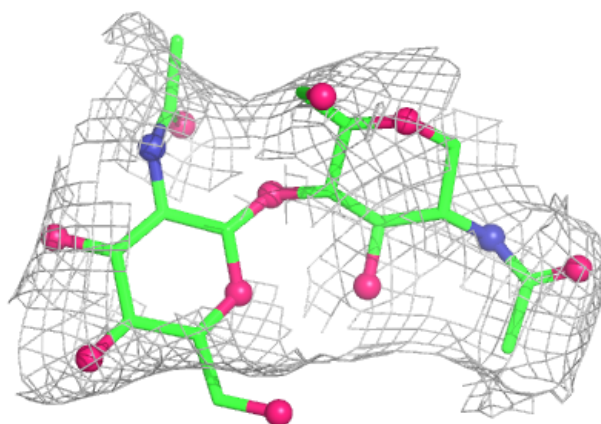
The following is a graphical depiction of the model fit to experimental electron density for oligosac-

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

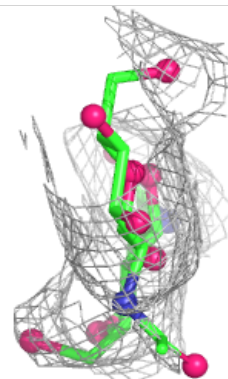
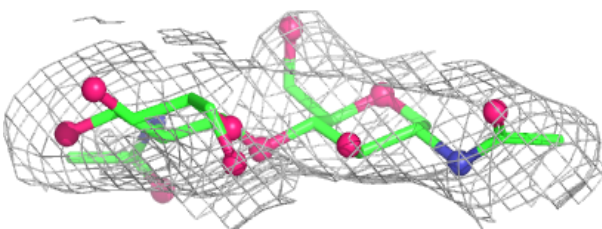
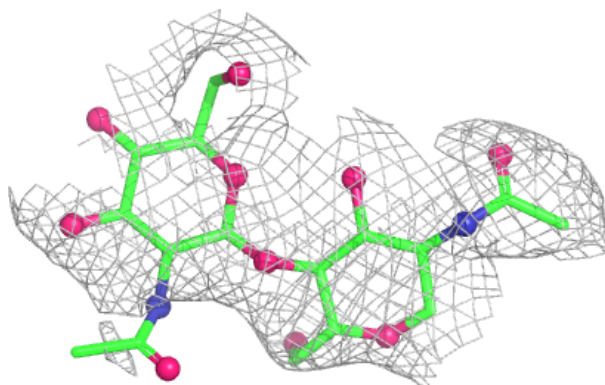


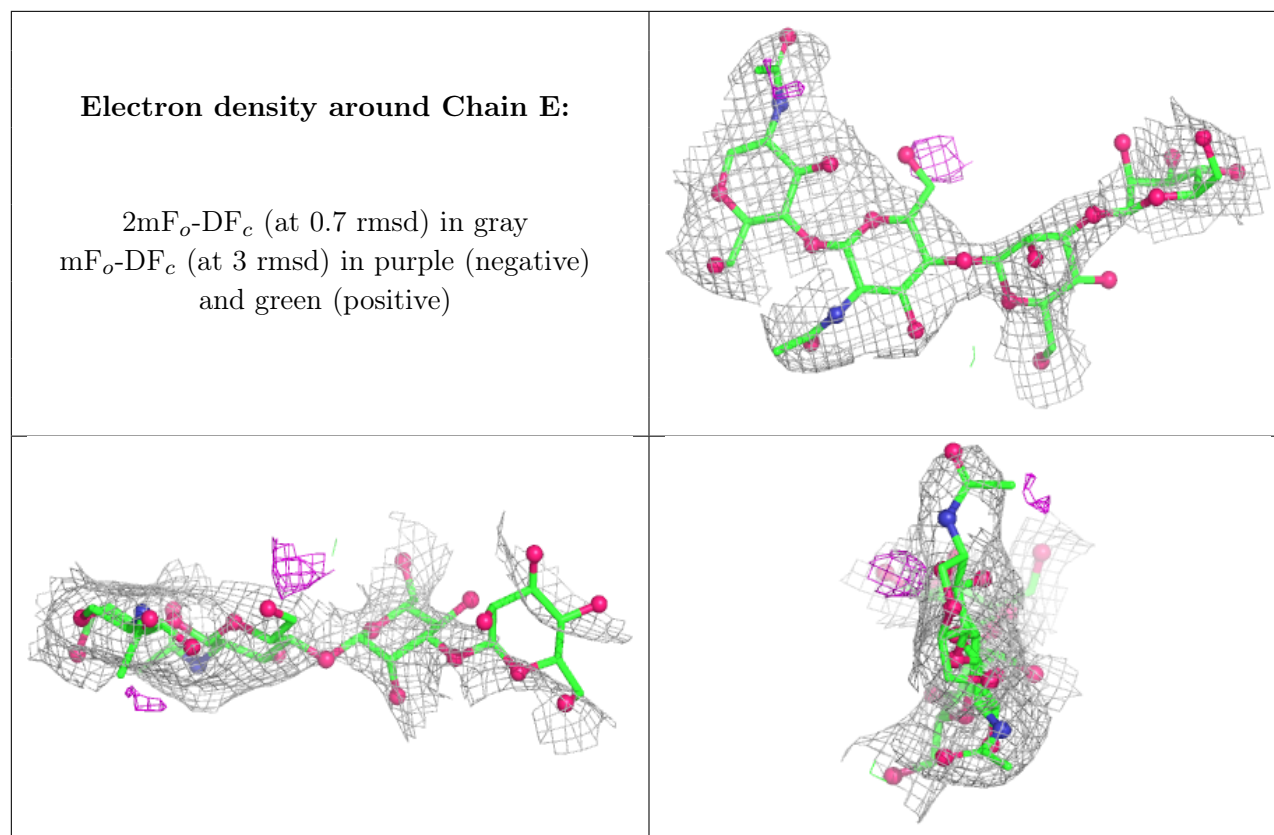
Electron density around Chain D:

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

**Electron density around Chain 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, 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
8	PLM	A	1323	17/18	0.94	0.08	98,105,113,114	0
6	ZN	A	962	1/1	0.95	0.12	80,80,80,80	1
8	PLM	B	1407	16/18	0.95	0.12	75,82,87,88	0
9	HTG	B	1506	19/19	0.95	0.09	88,93,97,99	0
10	HTO	B	1401	10/10	0.95	0.10	56,60,62,63	0
8	PLM	A	1322	17/18	0.96	0.15	94,104,110,110	0
8	PLM	A	1410	16/18	0.96	0.08	74,77,79,79	0
8	PLM	B	1323	17/18	0.96	0.11	117,122,124,124	0
6	ZN	B	963	1/1	0.97	0.14	93,93,93,93	1
7	RET	A	1296	20/21	0.97	0.14	50,52,54,55	20
9	HTG	A	1508	19/19	0.97	0.07	72,78,82,85	0
7	RET	B	1296	20/21	0.97	0.12	51,52,54,54	20
9	HTG	B	1509	19/19	0.97	0.10	93,95,100,100	0
8	PLM	B	1322	17/18	0.97	0.09	105,108,109,110	0

Continued on next page...

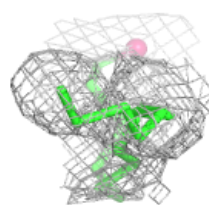
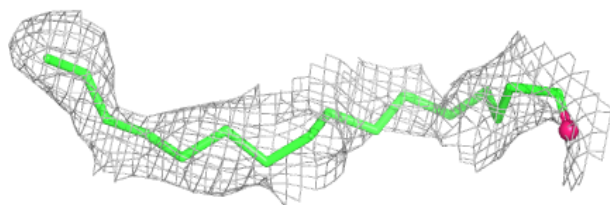
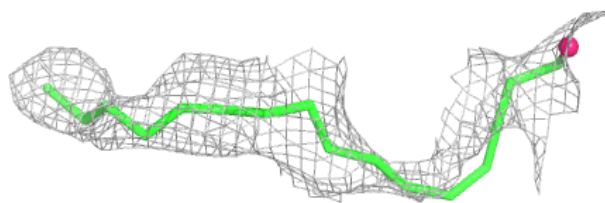
Continued from previous page...

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
5	HG	A	905	1/1	0.98	0.12	99,99,99,99	1
9	HTG	A	1507	19/19	0.98	0.10	99,113,115,116	0
6	ZN	A	2011	1/1	0.98	0.04	53,53,53,53	1
6	ZN	B	956	1/1	0.99	0.03	60,60,60,60	1
5	HG	B	902	1/1	0.99	0.04	80,80,80,80	1
5	HG	B	906	1/1	0.99	0.04	121,121,121,121	1
5	HG	A	903	1/1	0.99	0.05	96,96,96,96	1
5	HG	A	901	1/1	0.99	0.05	82,82,82,82	1
6	ZN	A	959	1/1	1.00	0.02	92,92,92,92	1
6	ZN	B	958	1/1	1.00	0.02	43,43,43,43	1
5	HG	B	904	1/1	1.00	0.03	109,109,109,109	1
6	ZN	A	957	1/1	1.00	0.02	53,53,53,53	1

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.

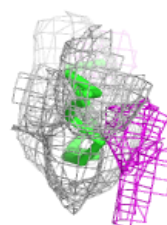
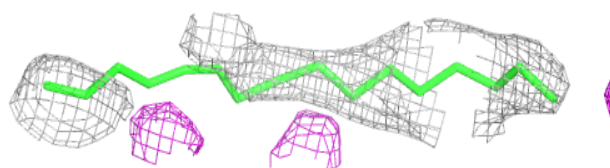
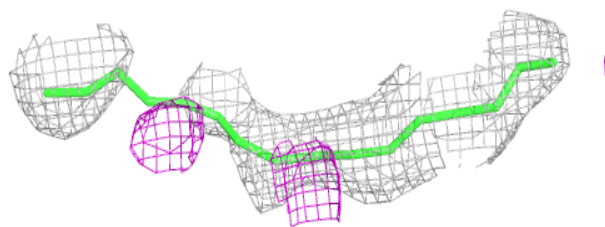
Electron density around PLM A 1323:

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

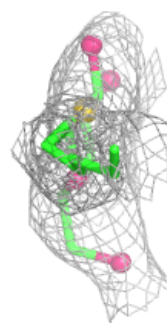
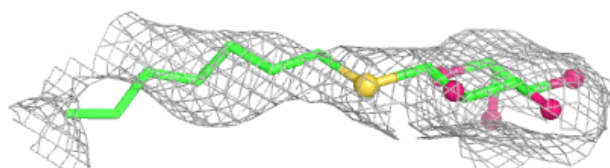
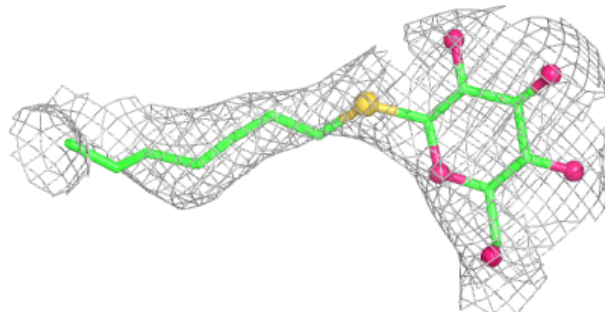


Electron density around PLM B 1407:

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

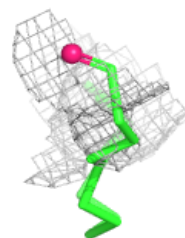
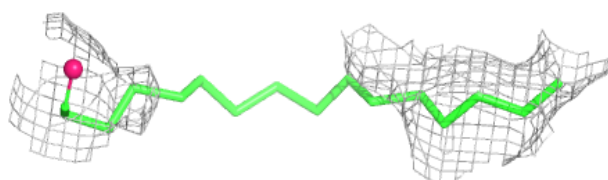
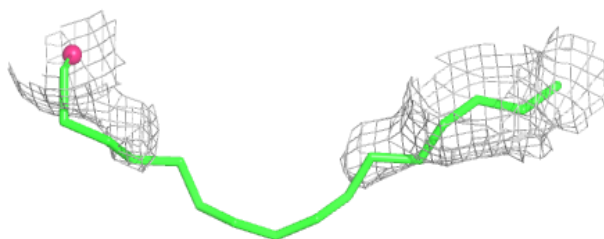
**Electron density around HTG B 1506:**

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

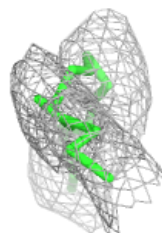
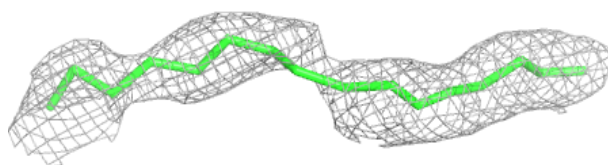
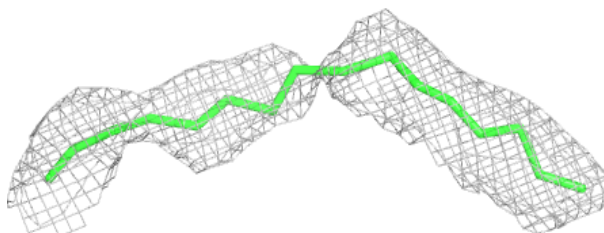


Electron density around PLM A 1322:

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

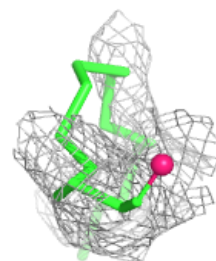
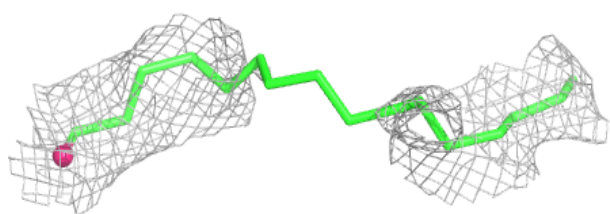
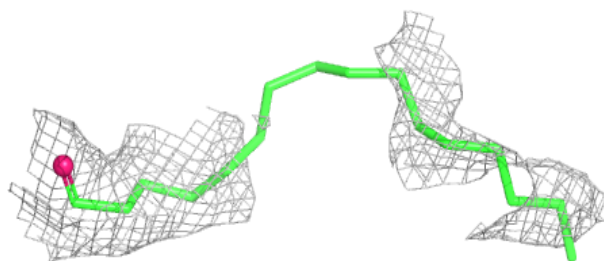
**Electron density around PLM A 1410:**

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

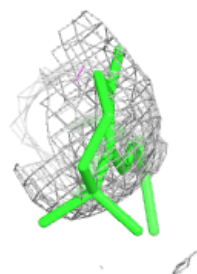
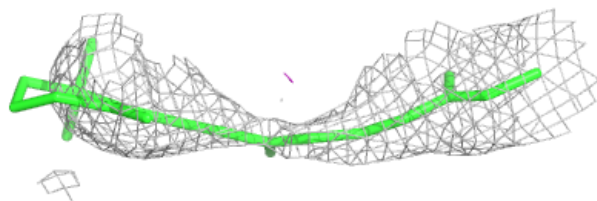
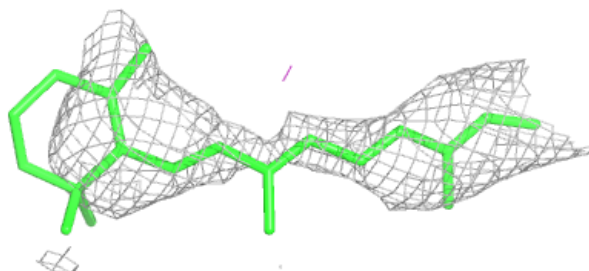


Electron density around PLM B 1323:

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

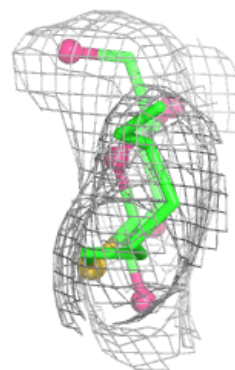
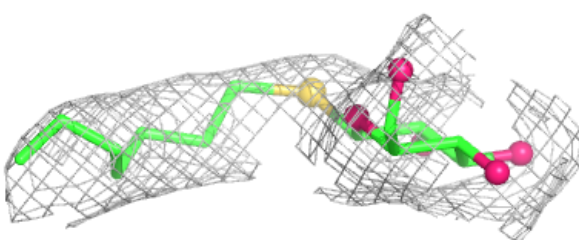
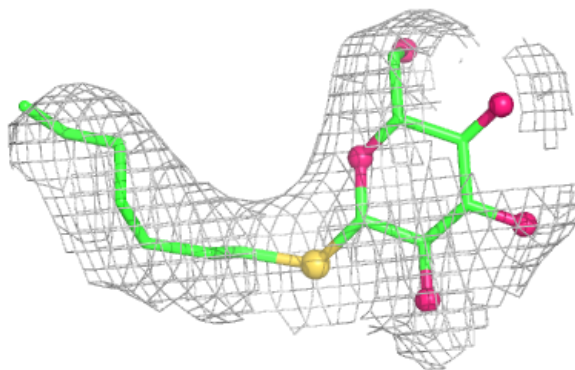
**Electron density around RET A 1296:**

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

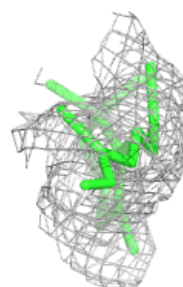
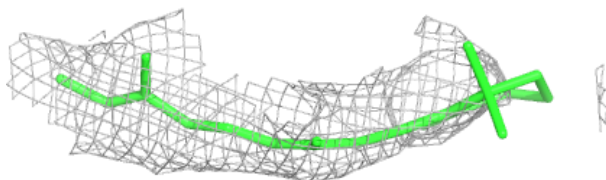
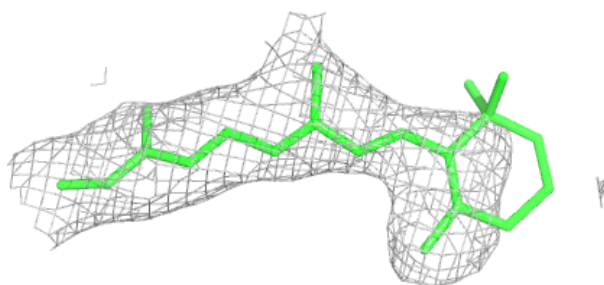


Electron density around HTG A 1508:

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

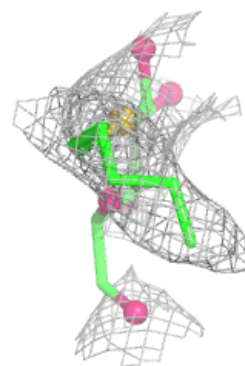
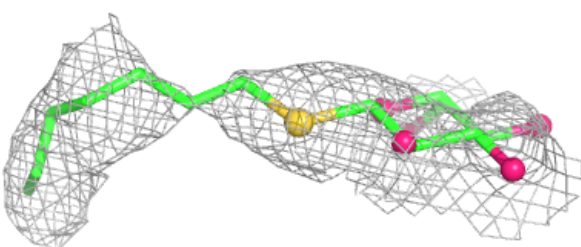
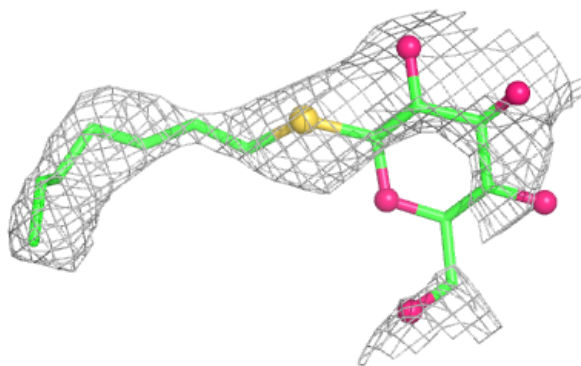
**Electron density around RET B 1296:**

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

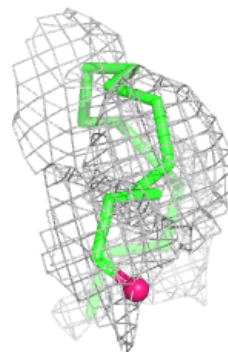
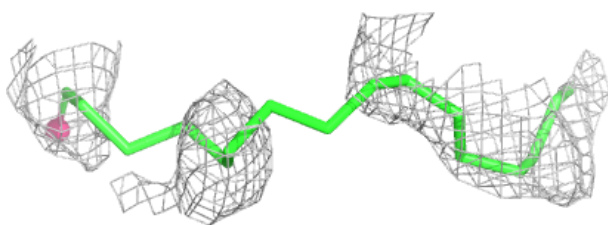
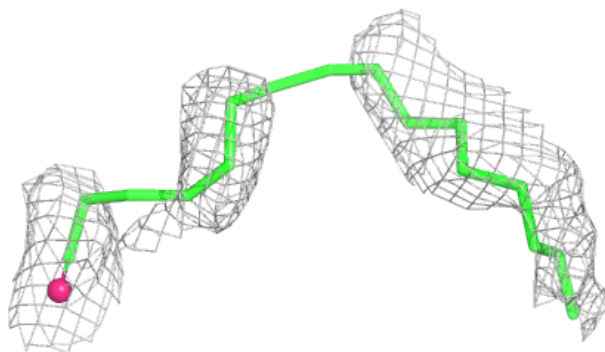


Electron density around HTG B 1509:

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

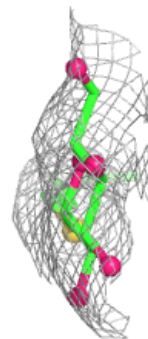
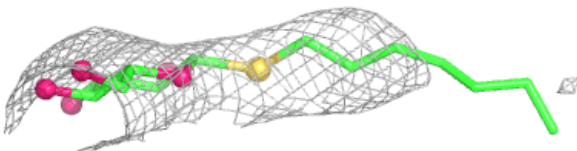
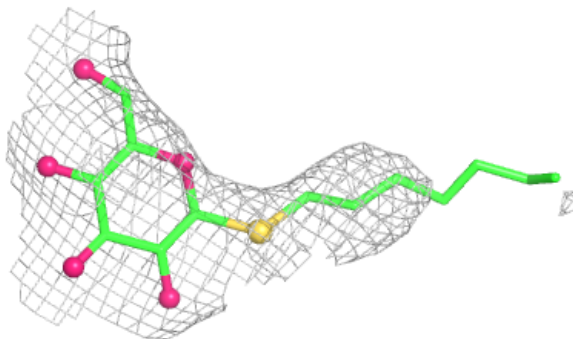
**Electron density around PLM B 1322:**

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



Electron density around HTG A 1507:

$2mF_o-DF_c$ (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.