



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 12:17 PM UTC

PDB ID : 7LAN / pdb_00007lan
Title : CRYSTAL STRUCTURE OF MYELOPEROXIDASE SUBFORM C (MPO)
COMPLEX WITH COMPOUND-30 AKA 7-[(3 {S},4 {R},6 {R})-4-benzyl-2
-oxa-7,13,14-triazatetracyclo[14.3.1.1[^]{3,6}.1[^]{11,14}]docosa-1(19),11(21),12
,16(20),17-pentaen-10-yl]-3 {H}-triazolo[4,5-b]pyridin-5-amine
Authors : Khan, J.A.
Deposited on : 2021-01-06
Resolution : 2.28 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](#) ①) were used in the production of this report:

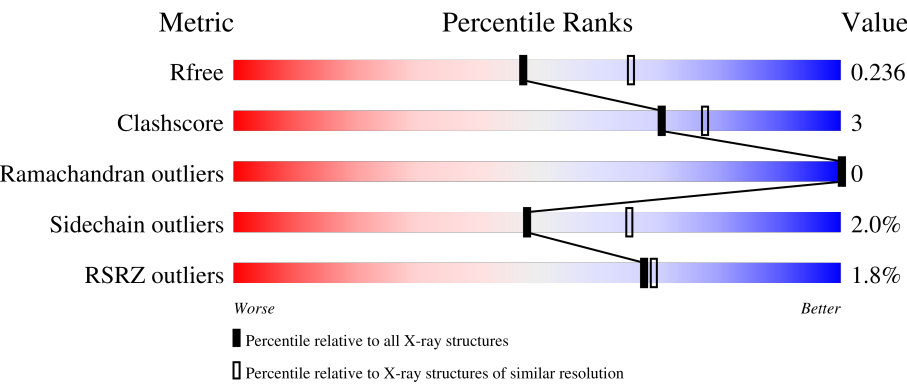
MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal 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 i

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

The reported resolution of this entry is 2.28 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	9078 (2.30-2.26)
Clashscore	190562	9802 (2.30-2.26)
Ramachandran outliers	187476	9690 (2.30-2.26)
Sidechain outliers	187428	9691 (2.30-2.26)
RSRZ outliers	180081	9085 (2.30-2.26)

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	105	2% 90% 8% ..
1	D	105	2% 92% 7% .
1	F	105	% 89% 9% ..
1	H	105	92% 6% .
2	B	466	4% 93% 7%

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Mol	Chain	Length	Quality of chain
2	E	466	% 93% 6%
2	G	466	% 91% 8%
2	I	466	2% 94% 6%
3	C	2	100%
3	K	2	100%
3	M	2	50% 50%
4	J	6	33% 67%
4	L	6	33% 67%
4	N	6	50% 50%
4	O	6	67% 33%

2 Entry composition [i](#)

There are 10 unique types of molecules in this entry. The entry contains 19728 atoms, of which 123 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 Myeloperoxidase light chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	103	Total	C	N	O	S	3	0	0
			821	520	145	151	5			
1	D	104	Total	C	N	O	S	5	0	0
			835	528	147	155	5			
1	F	103	Total	C	N	O	S	3	0	0
			821	520	145	151	5			
1	H	103	Total	C	N	O	S	0	0	0
			809	515	144	145	5			

- Molecule 2 is a protein called Isoform H14 of Myeloperoxidase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	464	Total	C	N	O	S	32	0	0
			3636	2298	654	658	26			
2	E	465	Total	C	N	O	S	36	0	0
			3651	2311	659	654	27			
2	G	464	Total	C	N	O	S	32	0	0
			3662	2314	664	657	27			
2	I	464	Total	C	N	O	S	36	0	0
			3636	2302	655	652	27			

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



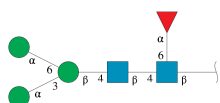
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	C	2	Total	C	N	O	0	0	0
			28	16	2	10			

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

- Molecule 4 is an oligosaccharide called alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-6)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-[alpha-L-fucopyranose-(1-6)]2-acetamido-2-deoxy-beta-D-glucopyranose.



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
4	J	6	Total	C	N	O	0	0	0
			71	40	2	29			
4	L	6	Total	C	N	O	0	0	0
			71	40	2	29			
4	N	6	Total	C	N	O	0	0	0
			71	40	2	29			
4	O	6	Total	C	N	O	0	0	0
			71	40	2	29			

- Molecule 5 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	1	Total	Cl	0	0
			1	1		
5	B	1	Total	Cl	0	0
			1	1		
5	F	1	Total	Cl	0	0
			1	1		
5	G	1	Total	Cl	0	0
			1	1		

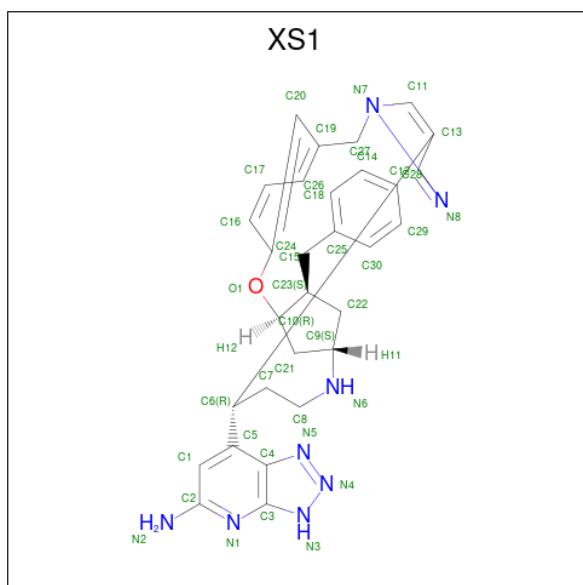
- Molecule 6 is PROTOPORPHYRIN IX CONTAINING FE (CCD ID: HEM) (formula: C₃₄H₃₂FeN₄O₄).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
7	B	1	Total C N O 14 8 1 5	0	0
7	E	1	Total C N O 14 8 1 5	0	0
7	G	1	Total C N O 14 8 1 5	0	0
7	I	1	Total C N O 14 8 1 5	0	0
7	I	1	Total C N O 14 8 1 5	0	0

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

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
8	B	1	Total Ca 1 1	0	0
8	E	1	Total Ca 1 1	0	0
8	G	1	Total Ca 1 1	0	0
8	I	1	Total Ca 1 1	0	0

- Molecule 9 is 7-[(3R,4S,6S,10R)-4-benzyl-2-oxa-7,13,14-triazatetracyclo[14.3.1.1 3,6 .1 11,14]docosa-1(20),11(21),12,16,18-pentaen-10-yl]-3H-[1,2,3]triazolo[4,5-b]pyridin-5-amine (CCD ID: XS1) (formula: C₃₀H₃₂N₈O) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
9	B	1	Total	C	H	N	O	27	0
			60	24	27	8	1		
9	E	1	Total	C	H	N	O	32	0
			71	30	32	8	1		
9	G	1	Total	C	H	N	O	32	0
			71	30	32	8	1		
9	I	1	Total	C	H	N	O	32	0
			71	30	32	8	1		

- Molecule 10 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
10	A	59	Total	O	0	0
			59	59		
10	B	162	Total	O	0	0
			162	162		
10	D	61	Total	O	0	0
			61	61		
10	E	213	Total	O	0	0
			213	213		
10	F	56	Total	O	0	0
			56	56		
10	G	188	Total	O	0	0
			188	188		
10	H	47	Total	O	0	0
			47	47		
10	I	180	Total	O	0	0
			180	180		

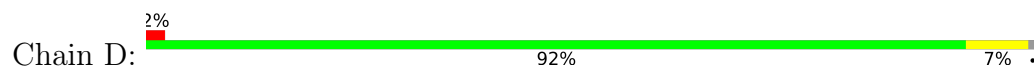
3 Residue-property plots [i](#)

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.

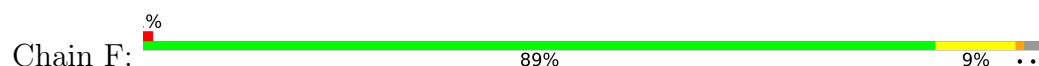
- Molecule 1: Myeloperoxidase light chain



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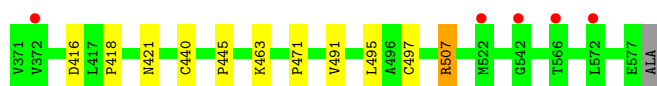


- Molecule 1: Myeloperoxidase light chain



- Molecule 2: Isoform H14 of Myeloperoxidase





- Molecule 2: Isoform H14 of Myeloperoxidase



- Molecule 2: Isoform H14 of Myeloperoxidase



- Molecule 2: Isoform H14 of Myeloperoxidase



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



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



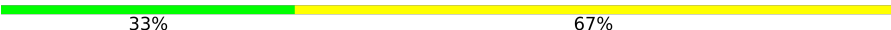
MAG1
MAG2

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

Chain M:  50% 50%

MAG1
MAG2

- Molecule 4: alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-6)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-[alpha-L-fucopyranose-(1-6)]2-acetamido-2-deoxy-beta-D-glucopyranose

Chain J:  33% 67%

MAG1
MAG2
BMA3
MAN4
MAN5
FUC6

- Molecule 4: alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-6)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-[alpha-L-fucopyranose-(1-6)]2-acetamido-2-deoxy-beta-D-glucopyranose

Chain L:  33% 67%

MAG1
MAG2
BMA3
MAN4
MAN5
FUC6

- Molecule 4: alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-6)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-[alpha-L-fucopyranose-(1-6)]2-acetamido-2-deoxy-beta-D-glucopyranose

Chain N:  50% 50%

MAG1
MAG2
BMA3
MAN4
MAN5
FUC6

- Molecule 4: alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-6)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-[alpha-L-fucopyranose-(1-6)]2-acetamido-2-deoxy-beta-D-glucopyranose

Chain O:  67% 33%

MAG1
MAG2
BMA3
MAN4
MAN5
FUC6

4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	144.21Å 150.39Å 231.01Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	47.45 – 2.28 47.45 – 2.28	Depositor EDS
% Data completeness (in resolution range)	100.0 (47.45-2.28) 99.9 (47.45-2.28)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.78 (at 2.27Å)	Xtriage
Refinement program	BUSTER 2.11.5	Depositor
R, R_{free}	0.200 , 0.239 0.195 , 0.236	Depositor DCC
R_{free} test set	5724 reflections (4.99%)	wwPDB-VP
Wilson B-factor (Å ²)	39.1	Xtriage
Anisotropy	0.281	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 53.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.016 for -k,-h,-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	19728	wwPDB-VP
Average B, all atoms (Å ²)	44.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.57% 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: NAG, CL, FUC, HEM, CA, XS1, MAN, 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	0.81	0/846	1.17	5/1154 (0.4%)
1	D	0.78	0/860	1.17	4/1172 (0.3%)
1	F	0.87	1/846 (0.1%)	1.15	1/1154 (0.1%)
1	H	0.83	0/834	1.17	1/1139 (0.1%)
2	B	0.82	0/3721	1.17	1/5063 (0.0%)
2	E	0.84	0/3736	1.14	6/5082 (0.1%)
2	G	0.84	0/3747	1.16	5/5093 (0.1%)
2	I	0.83	0/3722	1.14	3/5065 (0.1%)
All	All	0.83	1/18312 (0.0%)	1.16	26/24922 (0.1%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	F	87	MET	SD-CE	6.13	1.94	1.79

All (26) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	9	THR	N-CA-C	-6.04	102.74	110.53
1	D	9	THR	N-CA-C	-5.93	102.88	110.53
1	D	96	ASP	N-CA-C	-5.87	105.95	113.23
2	E	157	ASN	CA-CB-CG	5.75	118.35	112.60
2	E	114	ASN	CA-CB-CG	5.50	118.10	112.60
2	B	147	PHE	CA-CB-CG	5.50	119.30	113.80
1	H	9	THR	N-CA-C	-5.48	103.46	110.53
2	G	147	PHE	CA-CB-CG	5.47	119.28	113.80
2	G	318	ASP	CA-CB-CG	5.46	118.06	112.60
2	I	147	PHE	CA-CB-CG	5.42	119.22	113.80
1	A	92	LEU	CA-C-N	5.31	127.34	120.44
1	A	92	LEU	C-N-CA	5.31	127.34	120.44

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	I	272	ASP	CA-C-N	5.21	125.88	119.99
2	I	272	ASP	C-N-CA	5.21	125.88	119.99
1	A	37	TYR	CA-C-N	5.20	127.50	120.38
1	A	37	TYR	C-N-CA	5.20	127.50	120.38
1	D	37	TYR	CA-C-N	5.18	127.48	120.38
1	D	37	TYR	C-N-CA	5.18	127.48	120.38
2	E	284	VAL	CA-C-N	5.18	125.69	119.94
2	E	284	VAL	C-N-CA	5.18	125.69	119.94
2	E	205	ASP	CA-CB-CG	5.14	117.74	112.60
1	F	9	THR	N-CA-C	-5.09	103.96	110.53
2	G	157	ASN	CA-CB-CG	5.08	117.68	112.60
2	E	272	ASP	CA-CB-CG	5.03	117.63	112.60
2	G	432	GLY	CA-C-N	5.03	127.33	120.54
2	G	432	GLY	C-N-CA	5.03	127.33	120.54

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	821	0	776	13	0
1	D	835	0	793	10	0
1	F	821	0	776	14	0
1	H	809	0	761	11	0
2	B	3636	0	3562	19	0
2	E	3651	0	3598	11	0
2	G	3662	0	3617	20	0
2	I	3636	0	3566	12	0
3	C	28	0	25	0	0
3	K	28	0	25	0	0
3	M	28	0	25	0	0
4	J	71	0	61	0	0
4	L	71	0	61	0	0
4	N	71	0	61	0	0
4	O	71	0	61	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	A	1	0	0	0	0
5	B	1	0	0	0	0
5	F	1	0	0	0	0
5	G	1	0	0	0	0
6	B	43	0	30	9	0
6	E	43	0	30	9	0
6	G	43	0	30	9	0
6	I	43	0	30	9	0
7	B	14	0	13	0	0
7	E	14	0	13	0	0
7	G	14	0	13	0	0
7	I	28	0	26	0	0
8	B	1	0	0	0	0
8	E	1	0	0	0	0
8	G	1	0	0	0	0
8	I	1	0	0	0	0
9	B	33	27	0	0	0
9	E	39	32	0	0	0
9	G	39	32	0	0	0
9	I	39	32	0	0	0
10	A	59	0	0	0	0
10	B	162	0	0	0	0
10	D	61	0	0	0	0
10	E	213	0	0	0	0
10	F	56	0	0	0	0
10	G	188	0	0	0	0
10	H	47	0	0	0	0
10	I	180	0	0	0	0
All	All	19605	123	17953	102	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 (102) 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:F:94:ASP:CG	6:G:601:HEM:HMD3	1.41	1.45
1:D:94:ASP:CG	6:E:601:HEM:HMD3	1.41	1.44
1:F:94:ASP:CG	6:G:601:HEM:CMD	1.92	1.42
1:D:94:ASP:CG	6:E:601:HEM:CMD	1.96	1.38
1:D:94:ASP:OD2	6:E:601:HEM:CMD	1.71	1.38

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:94:ASP:CG	6:I:601:HEM:CMD	1.97	1.37
1:F:94:ASP:OD2	6:G:601:HEM:CMD	1.70	1.36
1:F:94:ASP:OD1	6:G:601:HEM:CMD	1.73	1.34
1:H:94:ASP:OD1	6:I:601:HEM:CMD	1.75	1.34
1:H:94:ASP:CG	6:I:601:HEM:HMD3	1.47	1.33
1:A:94:ASP:CG	6:B:601:HEM:HMD3	1.52	1.31
1:H:94:ASP:OD2	6:I:601:HEM:CMD	1.78	1.31
1:A:94:ASP:OD1	6:B:601:HEM:CMD	1.79	1.29
1:A:94:ASP:CG	6:B:601:HEM:CMD	2.05	1.29
1:A:94:ASP:OD2	6:B:601:HEM:HMD2	1.13	1.25
1:D:94:ASP:OD2	6:E:601:HEM:HMD2	1.06	1.22
1:A:94:ASP:OD2	6:B:601:HEM:CMD	1.86	1.22
1:D:94:ASP:OD1	6:E:601:HEM:CMD	1.83	1.21
1:H:94:ASP:OD2	6:I:601:HEM:HMD2	1.02	1.19
1:F:94:ASP:OD2	6:G:601:HEM:HMD2	0.96	1.14
1:D:94:ASP:OD1	6:E:601:HEM:HMD3	0.93	1.08
1:A:94:ASP:OD1	6:B:601:HEM:HMD3	0.89	1.05
1:H:94:ASP:OD1	6:I:601:HEM:HMD3	0.82	1.00
1:F:94:ASP:OD1	6:G:601:HEM:HMD3	0.81	0.99
1:F:94:ASP:CG	6:G:601:HEM:HMD2	1.75	0.95
1:H:94:ASP:CG	6:I:601:HEM:HMD2	1.76	0.91
1:D:94:ASP:CG	6:E:601:HEM:HMD2	1.81	0.86
2:E:313:TYR:HD1	2:E:507:ARG:HD3	1.44	0.82
1:A:94:ASP:CG	6:B:601:HEM:HMD2	1.88	0.80
2:E:313:TYR:CD1	2:E:507:ARG:HD3	2.22	0.74
2:E:113:VAL:HG21	2:E:122:GLN:HB3	1.73	0.71
2:B:284:VAL:HA	2:B:287:MET:HE2	1.74	0.69
2:G:284:VAL:HA	2:G:287:MET:HE2	1.75	0.68
1:A:94:ASP:OD1	6:B:601:HEM:C2D	2.50	0.64
2:B:313:TYR:HD1	2:B:507:ARG:HD3	1.62	0.64
2:G:367:ALA:HB1	2:G:370:ARG:HG3	1.80	0.63
1:A:94:ASP:OD2	2:B:332:PHE:HD2	1.85	0.60
1:F:94:ASP:OD1	6:G:601:HEM:C2D	2.52	0.57
2:B:241:SER:O	2:B:366:PHE:HA	2.04	0.57
1:H:94:ASP:OD2	2:I:332:PHE:HD2	1.89	0.56
2:G:256:GLU:HG2	2:G:287:MET:HE1	1.87	0.56
2:B:256:GLU:HG2	2:B:287:MET:HE1	1.86	0.56
1:D:94:ASP:OD1	6:E:601:HEM:C2D	2.57	0.53
1:H:92:LEU:HD22	2:I:249:MET:HB3	1.90	0.53
6:I:601:HEM:HMC2	6:I:601:HEM:HBC2	1.93	0.51
1:H:94:ASP:OD1	6:I:601:HEM:C2D	2.60	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:92:LEU:HD22	2:G:249:MET:HB3	1.94	0.50
2:G:313:TYR:HD1	2:G:507:ARG:HD3	1.74	0.50
2:E:333:ARG:HH11	2:E:421:ASN:ND2	2.09	0.50
2:I:333:ARG:HH11	2:I:421:ASN:ND2	2.10	0.50
2:B:120:VAL:HG11	2:G:443:PRO:HG2	1.95	0.49
2:G:333:ARG:HH11	2:G:421:ASN:ND2	2.10	0.49
2:B:367:ALA:HB1	2:B:370:ARG:HG3	1.94	0.49
1:F:94:ASP:OD2	2:G:332:PHE:HD2	1.95	0.49
2:G:256:GLU:CG	2:G:287:MET:HE1	2.43	0.49
6:B:601:HEM:HBC2	6:B:601:HEM:HMC2	1.95	0.48
1:F:1:CYS:SG	1:F:14:CYS:SG	3.08	0.48
2:I:241:SER:O	2:I:366:PHE:HA	2.14	0.48
2:B:333:ARG:HH11	2:B:421:ASN:ND2	2.11	0.48
2:G:333:ARG:HH11	2:G:421:ASN:HD22	1.61	0.47
6:G:601:HEM:HBC2	6:G:601:HEM:HMC2	1.96	0.47
1:A:68:ILE:CD1	2:B:463:LYS:HB3	2.45	0.47
2:B:256:GLU:CG	2:B:287:MET:HE1	2.45	0.46
2:B:313:TYR:CD1	2:B:507:ARG:HD3	2.47	0.46
2:I:333:ARG:HH11	2:I:421:ASN:HD22	1.63	0.46
2:B:440:CYS:CB	2:B:497:CYS:HG	2.25	0.46
1:H:16:ASN:O	1:H:20:PRO:HA	2.16	0.46
2:E:333:ARG:HH11	2:E:421:ASN:HD22	1.63	0.46
1:F:16:ASN:O	1:F:20:PRO:HA	2.16	0.46
1:A:68:ILE:HD13	2:B:463:LYS:HB3	1.97	0.45
2:I:416:ASP:OD1	2:I:418:PRO:HD2	2.16	0.45
1:A:16:ASN:O	1:A:20:PRO:HA	2.16	0.45
2:G:434:ASN:HB2	2:G:472:ASN:HA	1.99	0.45
2:B:333:ARG:HH11	2:B:421:ASN:HD22	1.64	0.45
1:D:16:ASN:O	1:D:20:PRO:HA	2.17	0.45
2:B:445:PRO:O	2:B:471:PRO:HG2	2.17	0.45
2:E:153:CYS:HB3	2:E:159:THR:HG21	2.00	0.44
2:I:153:CYS:SG	2:I:156:SER:HB2	2.58	0.44
2:G:137:ILE:HG12	2:G:413:ILE:HD11	1.99	0.44
2:E:119:CYS:HG	2:E:143:CYS:CB	2.23	0.44
2:G:259:ARG:HD2	2:G:539:ASP:HB3	2.00	0.43
2:E:137:ILE:HG12	2:E:413:ILE:HD11	2.01	0.43
2:G:416:ASP:OD1	2:G:418:PRO:HD2	2.19	0.43
2:G:491:VAL:HB	2:G:495:LEU:HB2	2.01	0.43
2:I:491:VAL:HB	2:I:495:LEU:HB2	2.01	0.43
2:B:491:VAL:HB	2:B:495:LEU:HB2	2.01	0.42
2:E:416:ASP:OD1	2:E:418:PRO:HD2	2.20	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:416:ASP:OD1	2:B:418:PRO:HD2	2.20	0.42
2:G:507:ARG:HD3	2:G:508:ASP:OD1	2.19	0.42
2:I:153:CYS:HB3	2:I:159:THR:HG21	2.02	0.42
6:E:601:HEM:HMC2	6:E:601:HEM:HBC2	2.02	0.42
2:G:153:CYS:SG	2:G:156:SER:HB2	2.60	0.41
1:A:1:CYS:SG	1:A:14:CYS:SG	3.02	0.41
2:B:119:CYS:SG	2:G:456:ARG:HG2	2.61	0.41
1:F:29:PHE:CE1	2:G:165:ASN:HB2	2.56	0.41
2:E:491:VAL:HB	2:E:495:LEU:HB2	2.03	0.41
2:I:410:VAL:HG23	2:I:411:MET:HG3	2.03	0.41
2:B:153:CYS:SG	2:B:156:SER:HB2	2.61	0.41
2:I:287:MET:HE3	2:I:570:LEU:HD22	2.03	0.41
1:D:94:ASP:OD2	2:E:332:PHE:HD2	2.03	0.40
1:F:83:SER:HB3	2:G:554:SER:O	2.22	0.40
2:I:137:ILE:HG12	2:I:413:ILE:HD11	2.03	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	101/105 (96%)	98 (97%)	3 (3%)	0	100	100
1	D	102/105 (97%)	98 (96%)	4 (4%)	0	100	100
1	F	101/105 (96%)	99 (98%)	2 (2%)	0	100	100
1	H	101/105 (96%)	98 (97%)	3 (3%)	0	100	100
2	B	462/466 (99%)	453 (98%)	9 (2%)	0	100	100
2	E	463/466 (99%)	452 (98%)	11 (2%)	0	100	100
2	G	462/466 (99%)	451 (98%)	11 (2%)	0	100	100
2	I	462/466 (99%)	453 (98%)	9 (2%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
All	All	2254/2284 (99%)	2202 (98%)	52 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	87/90 (97%)	86 (99%)	1 (1%)	65	79
1	D	89/90 (99%)	88 (99%)	1 (1%)	65	79
1	F	87/90 (97%)	86 (99%)	1 (1%)	65	79
1	H	83/90 (92%)	82 (99%)	1 (1%)	63	77
2	B	390/411 (95%)	383 (98%)	7 (2%)	51	68
2	E	393/411 (96%)	384 (98%)	9 (2%)	44	60
2	G	395/411 (96%)	384 (97%)	11 (3%)	38	53
2	I	389/411 (95%)	381 (98%)	8 (2%)	47	63
All	All	1913/2004 (96%)	1874 (98%)	39 (2%)	48	65

All (39) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	CYS
2	B	175	MET
2	B	181	GLU
2	B	254	LEU
2	B	259	ARG
2	B	359	VAL
2	B	361	LEU
2	B	507	ARG
1	D	1	CYS
2	E	159	THR
2	E	175	MET
2	E	218	ASP

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Mol	Chain	Res	Type
2	E	223	LEU
2	E	254	LEU
2	E	267	LEU
2	E	462	ARG
2	E	523	GLN
2	E	545	THR
1	F	1	CYS
2	G	122	GLN
2	G	175	MET
2	G	181	GLU
2	G	229	ARG
2	G	254	LEU
2	G	259	ARG
2	G	275	ARG
2	G	359	VAL
2	G	361	LEU
2	G	507	ARG
2	G	523	GLN
1	H	1	CYS
2	I	159	THR
2	I	175	MET
2	I	254	LEU
2	I	259	ARG
2	I	267	LEU
2	I	507	ARG
2	I	526	GLN
2	I	563	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (32) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	54	ASN
2	B	121	GLN
2	B	204	GLN
2	B	356	ASN
2	B	396	GLN
2	B	467	GLN
2	B	516	ASN
2	B	549	ASN
2	B	563	ASN
1	D	54	ASN
2	E	114	ASN

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Mol	Chain	Res	Type
2	E	121	GLN
2	E	204	GLN
2	E	396	GLN
2	E	409	GLN
2	E	467	GLN
2	E	516	ASN
1	F	54	ASN
2	G	204	GLN
2	G	356	ASN
2	G	396	GLN
2	G	409	GLN
2	G	516	ASN
2	G	540	ASN
2	G	549	ASN
1	H	54	ASN
2	I	121	GLN
2	I	396	GLN
2	I	473	ASN
2	I	516	ASN
2	I	540	ASN
2	I	549	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 ⓘ

30 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
3	NAG	C	1	2,3	14,14,15	0.34	0	17,19,21	1.19	1 (5%)
3	NAG	C	2	3	14,14,15	0.30	0	17,19,21	1.02	2 (11%)
4	NAG	J	1	4,2	14,14,15	0.33	0	17,19,21	0.76	1 (5%)
4	NAG	J	2	4	14,14,15	0.27	0	17,19,21	0.48	0
4	BMA	J	3	4	11,11,12	0.31	0	15,15,17	0.52	0
4	MAN	J	4	4	11,11,12	0.33	0	15,15,17	0.93	1 (6%)
4	MAN	J	5	4	11,11,12	0.37	0	15,15,17	0.92	1 (6%)
4	FUC	J	6	4	10,10,11	0.33	0	14,14,16	0.79	1 (7%)
3	NAG	K	1	2,3	14,14,15	0.28	0	17,19,21	0.86	1 (5%)
3	NAG	K	2	3	14,14,15	0.29	0	17,19,21	1.12	2 (11%)
4	NAG	L	1	4,2	14,14,15	0.34	0	17,19,21	0.84	1 (5%)
4	NAG	L	2	4	14,14,15	0.29	0	17,19,21	0.53	0
4	BMA	L	3	4	11,11,12	0.35	0	15,15,17	0.63	0
4	MAN	L	4	4	11,11,12	0.30	0	15,15,17	1.02	1 (6%)
4	MAN	L	5	4	11,11,12	0.30	0	15,15,17	0.87	1 (6%)
4	FUC	L	6	4	10,10,11	0.41	0	14,14,16	0.73	1 (7%)
3	NAG	M	1	2,3	14,14,15	0.30	0	17,19,21	1.15	2 (11%)
3	NAG	M	2	3	14,14,15	0.29	0	17,19,21	0.62	0
4	NAG	N	1	4,2	14,14,15	0.40	0	17,19,21	1.08	1 (5%)
4	NAG	N	2	4	14,14,15	0.37	0	17,19,21	0.64	0
4	BMA	N	3	4	11,11,12	0.43	0	15,15,17	0.66	0
4	MAN	N	4	4	11,11,12	0.30	0	15,15,17	0.89	1 (6%)
4	MAN	N	5	4	11,11,12	0.30	0	15,15,17	0.97	1 (6%)
4	FUC	N	6	4	10,10,11	0.33	0	14,14,16	0.48	0
4	NAG	O	1	4,2	14,14,15	0.30	0	17,19,21	0.70	0
4	NAG	O	2	4	14,14,15	0.29	0	17,19,21	0.50	0
4	BMA	O	3	4	11,11,12	0.27	0	15,15,17	0.67	0
4	MAN	O	4	4	11,11,12	0.32	0	15,15,17	1.11	1 (6%)
4	MAN	O	5	4	11,11,12	0.38	0	15,15,17	0.77	1 (6%)
4	FUC	O	6	4	10,10,11	0.37	0	14,14,16	0.62	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
3	NAG	C	1	2,3	-	0/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	NAG	C	2	3	-	0/6/23/26	0/1/1/1
4	NAG	J	1	4,2	-	0/6/23/26	0/1/1/1
4	NAG	J	2	4	-	0/6/23/26	0/1/1/1
4	BMA	J	3	4	-	0/2/19/22	0/1/1/1
4	MAN	J	4	4	-	1/2/19/22	0/1/1/1
4	MAN	J	5	4	-	0/2/19/22	0/1/1/1
4	FUC	J	6	4	-	-	0/1/1/1
3	NAG	K	1	2,3	-	0/6/23/26	0/1/1/1
3	NAG	K	2	3	-	1/6/23/26	0/1/1/1
4	NAG	L	1	4,2	-	0/6/23/26	0/1/1/1
4	NAG	L	2	4	-	0/6/23/26	0/1/1/1
4	BMA	L	3	4	-	0/2/19/22	0/1/1/1
4	MAN	L	4	4	-	1/2/19/22	0/1/1/1
4	MAN	L	5	4	-	0/2/19/22	0/1/1/1
4	FUC	L	6	4	-	-	0/1/1/1
3	NAG	M	1	2,3	-	0/6/23/26	0/1/1/1
3	NAG	M	2	3	-	0/6/23/26	0/1/1/1
4	NAG	N	1	4,2	-	1/6/23/26	0/1/1/1
4	NAG	N	2	4	-	0/6/23/26	0/1/1/1
4	BMA	N	3	4	-	0/2/19/22	0/1/1/1
4	MAN	N	4	4	-	1/2/19/22	0/1/1/1
4	MAN	N	5	4	-	0/2/19/22	0/1/1/1
4	FUC	N	6	4	-	-	0/1/1/1
4	NAG	O	1	4,2	-	0/6/23/26	0/1/1/1
4	NAG	O	2	4	-	0/6/23/26	0/1/1/1
4	BMA	O	3	4	-	0/2/19/22	0/1/1/1
4	MAN	O	4	4	-	1/2/19/22	0/1/1/1
4	MAN	O	5	4	-	0/2/19/22	0/1/1/1
4	FUC	O	6	4	-	-	0/1/1/1

There are no bond length outliers.

All (21) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	O	4	MAN	C1-O5-C5	3.94	117.47	112.19
3	C	1	NAG	C1-C2-N2	-3.76	104.51	110.43
4	N	1	NAG	O5-C1-C2	-3.64	105.66	111.29
4	L	4	MAN	C1-O5-C5	3.59	117.00	112.19
4	N	5	MAN	C1-O5-C5	3.47	116.83	112.19
3	M	1	NAG	O5-C1-C2	-3.39	106.04	111.29
4	J	5	MAN	C1-O5-C5	3.32	116.63	112.19
4	N	4	MAN	C1-O5-C5	3.21	116.49	112.19

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	J	4	MAN	C1-O5-C5	3.16	116.42	112.19
3	K	2	NAG	O5-C1-C2	-3.05	106.58	111.29
4	L	5	MAN	C1-O5-C5	3.00	116.20	112.19
3	C	2	NAG	O5-C1-C2	-2.77	107.00	111.29
4	O	5	MAN	C1-O5-C5	2.69	115.79	112.19
3	K	2	NAG	C1-O5-C5	2.58	115.64	112.19
3	K	1	NAG	O5-C1-C2	-2.57	107.31	111.29
3	M	1	NAG	C1-C2-N2	-2.57	106.38	110.43
4	J	6	FUC	C1-O5-C5	2.44	118.73	112.97
3	C	2	NAG	C1-O5-C5	2.40	115.40	112.19
4	L	6	FUC	C1-O5-C5	2.34	118.49	112.97
4	L	1	NAG	O5-C1-C2	-2.32	107.70	111.29
4	J	1	NAG	O5-C1-C2	-2.13	107.99	111.29

There are no chirality outliers.

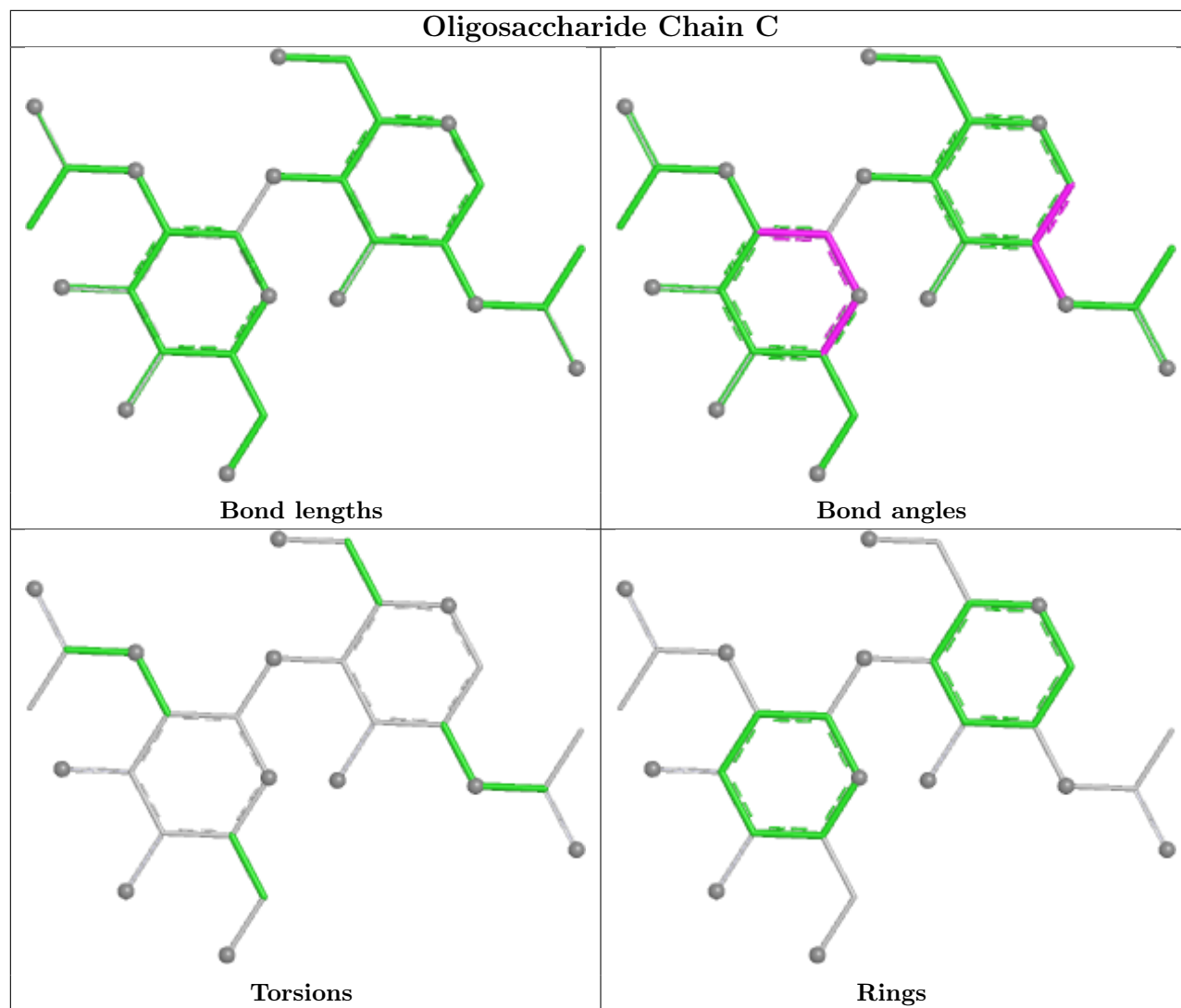
All (6) torsion outliers are listed below:

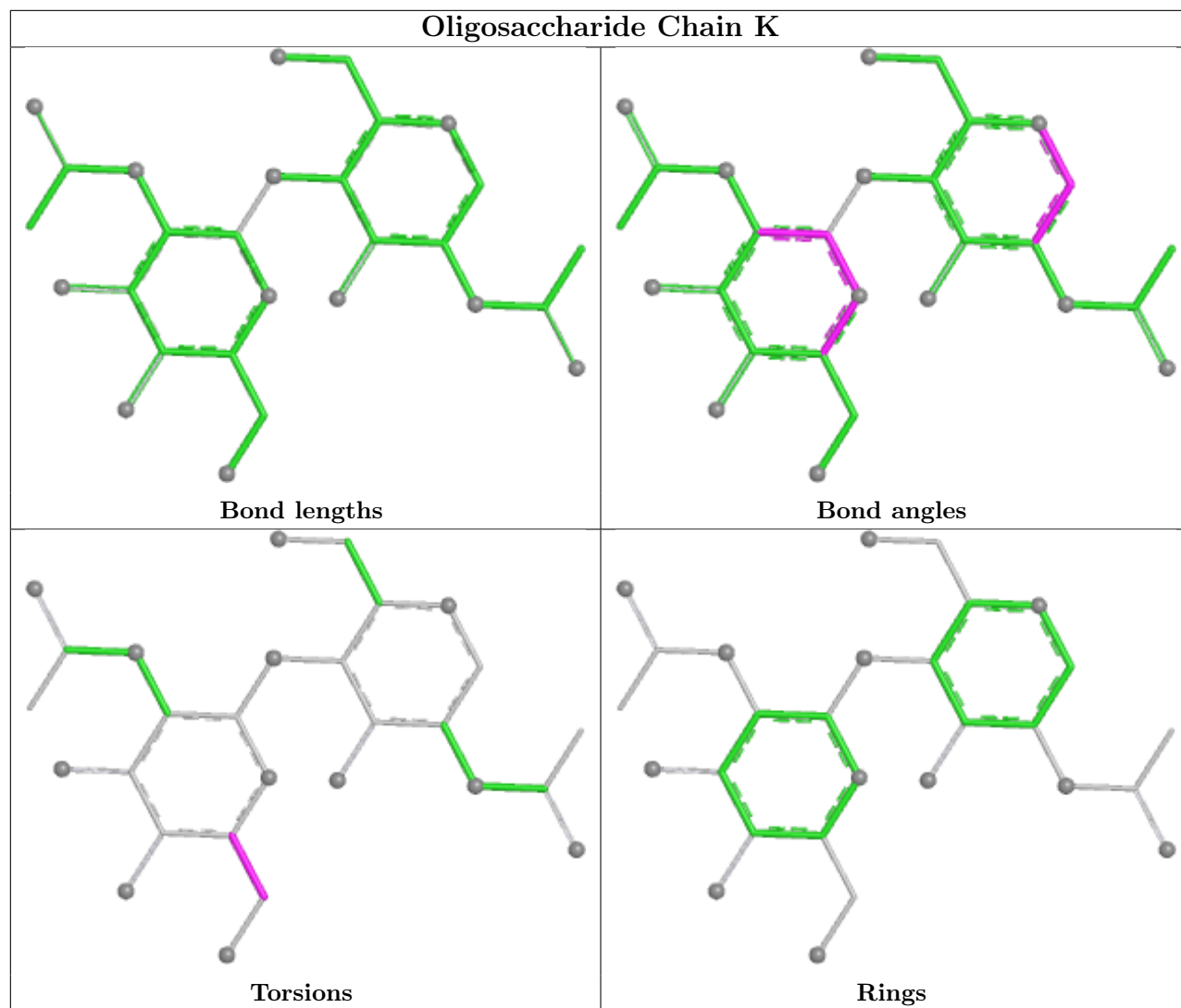
Mol	Chain	Res	Type	Atoms
4	J	4	MAN	O5-C5-C6-O6
4	L	4	MAN	O5-C5-C6-O6
4	N	4	MAN	O5-C5-C6-O6
4	O	4	MAN	O5-C5-C6-O6
4	N	1	NAG	C4-C5-C6-O6
3	K	2	NAG	O5-C5-C6-O6

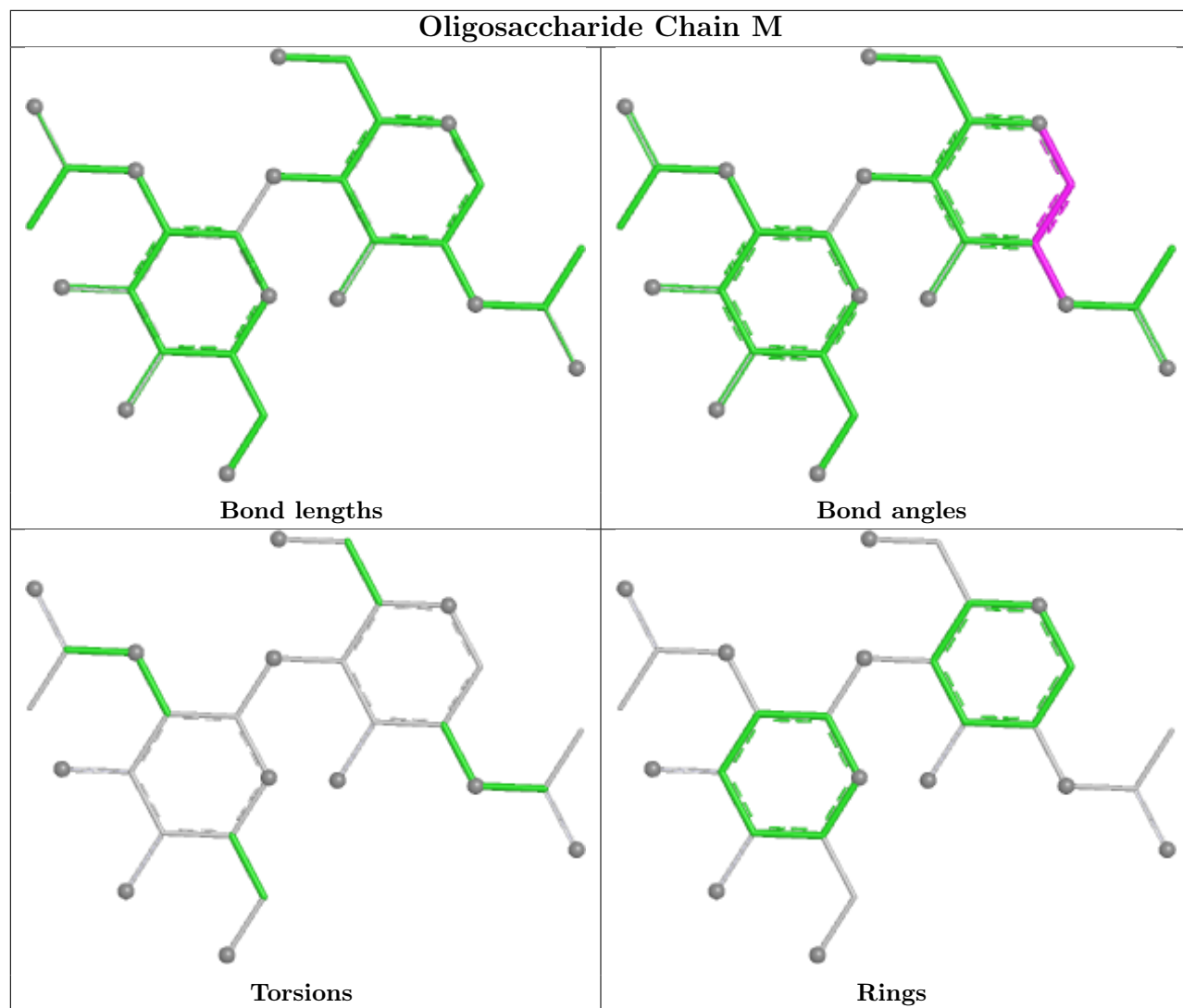
There are no ring outliers.

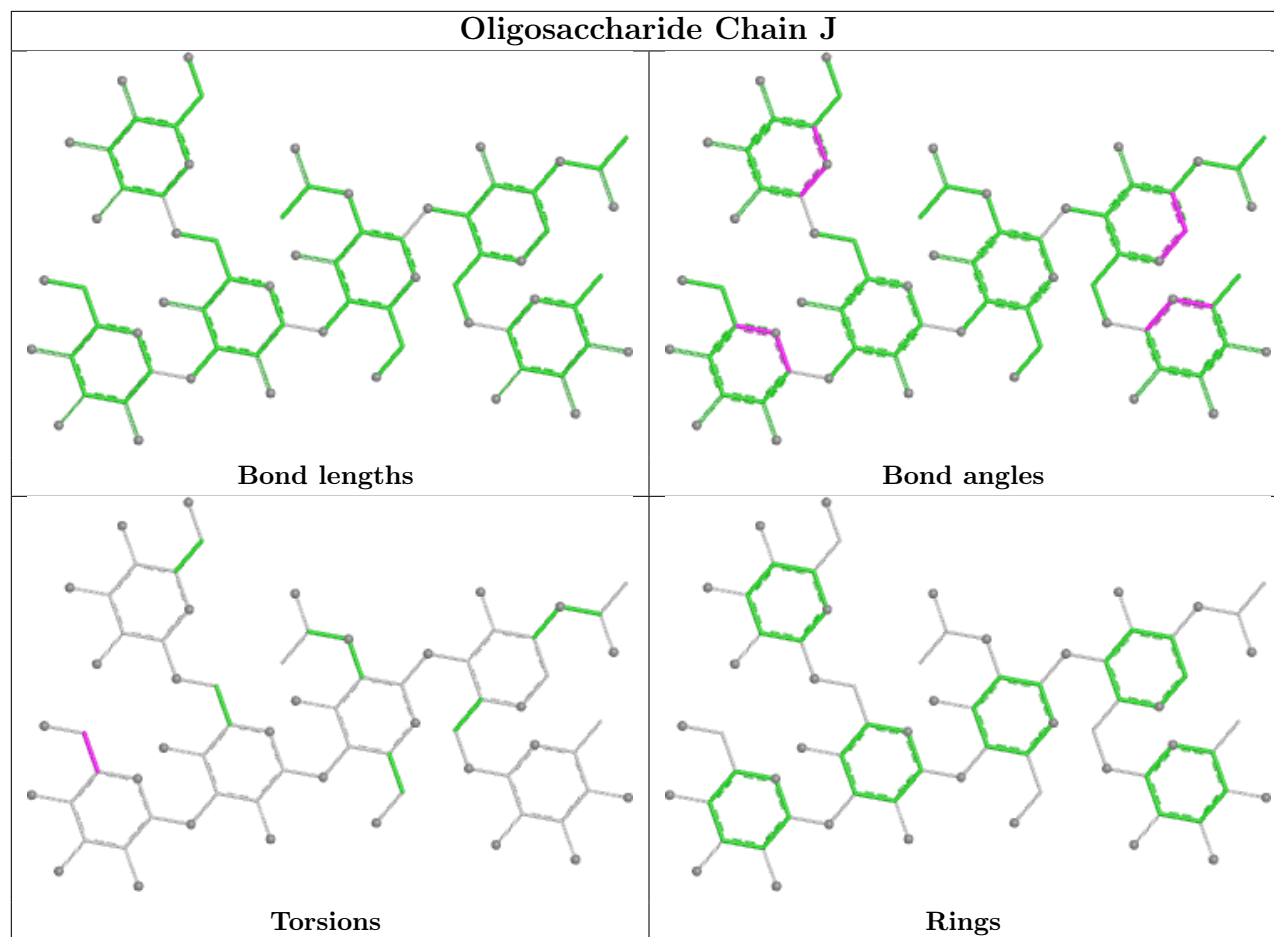
No monomer is involved in short contacts.

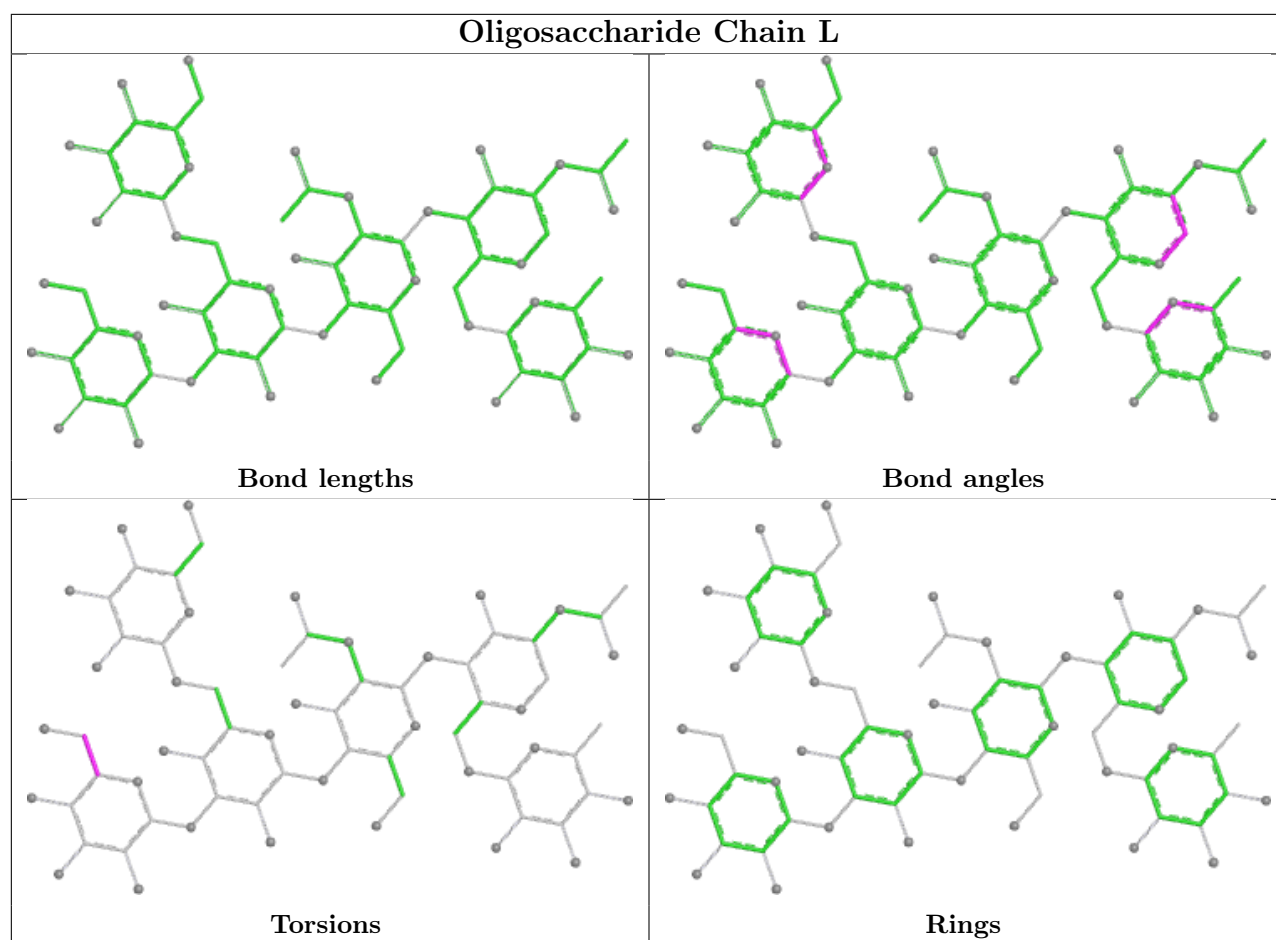
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.

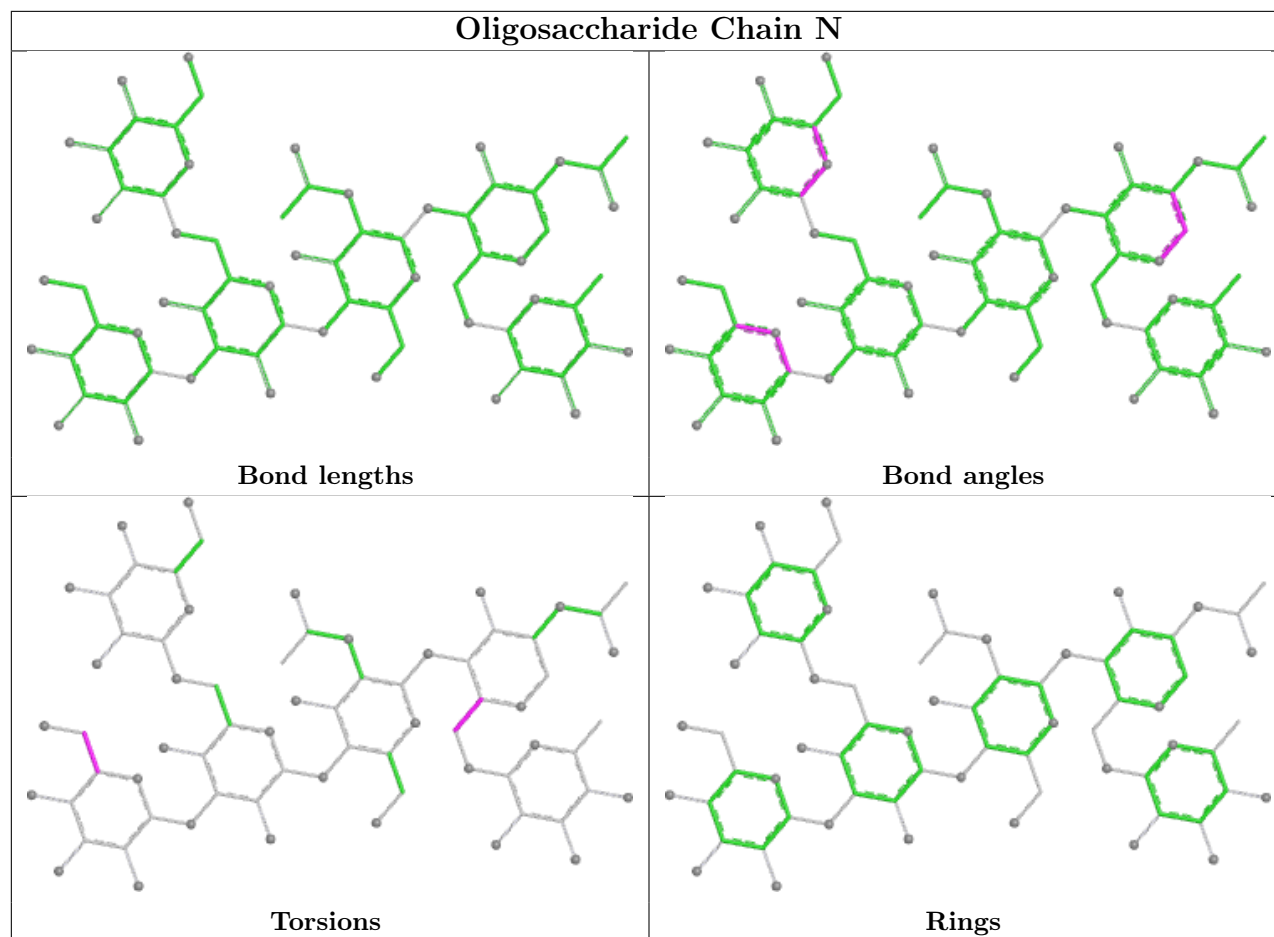


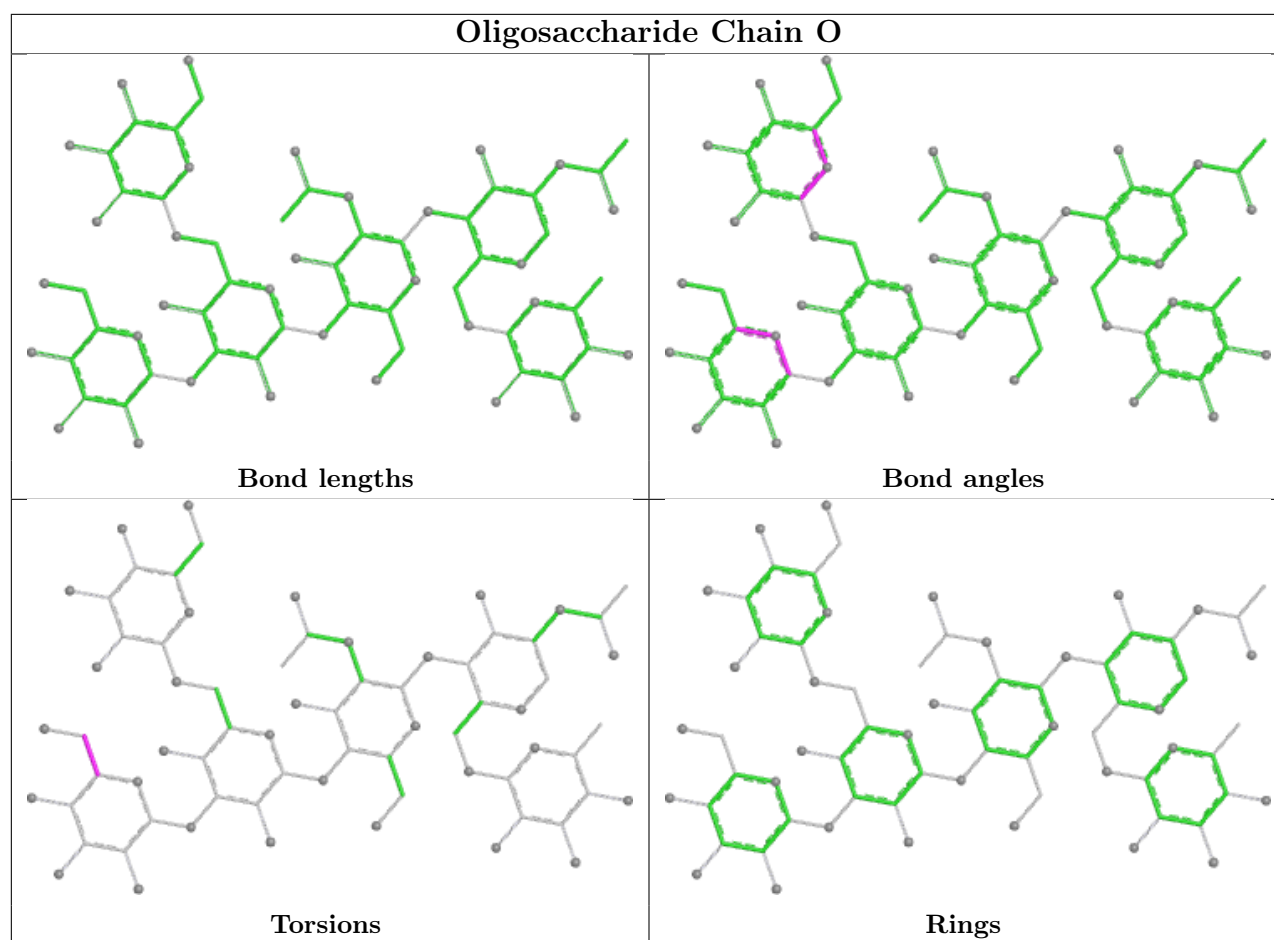












5.6 Ligand geometry [i](#)

Of 21 ligands modelled in this entry, 8 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	$\# Z > 2$
9	XS1	I	605	-	45,45,45	0.76	1 (2%)	46,64,64	1.09	2 (4%)
6	HEM	G	601	2,10	50,50,50	1.47	10 (20%)	67,82,82	1.47	15 (22%)
6	HEM	B	601	2,10	50,50,50	1.53	8 (16%)	67,82,82	1.46	10 (14%)
7	NAG	I	602	2	14,14,15	0.28	0	17,19,21	0.85	1 (5%)
9	XS1	G	605	-	45,45,45	0.78	1 (2%)	46,64,64	1.07	2 (4%)
7	NAG	I	603	2	14,14,15	0.26	0	17,19,21	0.77	1 (5%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
6	HEM	E	601	2	50,50,50	1.43	9 (18%)	67,82,82	1.41	9 (13%)
7	NAG	G	602	2	14,14,15	0.30	0	17,19,21	0.83	1 (5%)
9	XS1	E	604	-	45,45,45	0.82	1 (2%)	46,64,64	1.13	2 (4%)
6	HEM	I	601	2	50,50,50	1.47	8 (16%)	67,82,82	1.52	12 (17%)
9	XS1	B	605	-	37,38,45	0.82	1 (2%)	37,55,64	1.25	2 (5%)
7	NAG	B	602	2	14,14,15	0.31	0	17,19,21	0.90	1 (5%)
7	NAG	E	602	2	14,14,15	0.27	0	17,19,21	0.88	1 (5%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
9	XS1	I	605	-	-	5/26/38/38	0/6/7/7
6	HEM	G	601	2,10	-	4/14/54/54	-
6	HEM	B	601	2,10	-	4/14/54/54	-
7	NAG	I	602	2	-	0/6/23/26	0/1/1/1
9	XS1	G	605	-	-	4/26/38/38	0/6/7/7
7	NAG	I	603	2	-	0/6/23/26	0/1/1/1
6	HEM	E	601	2	-	4/14/54/54	-
7	NAG	G	602	2	-	0/6/23/26	0/1/1/1
9	XS1	E	604	-	-	5/26/38/38	0/6/7/7
6	HEM	I	601	2	-	4/14/54/54	-
9	XS1	B	605	-	-	3/22/34/38	0/5/6/7
7	NAG	B	602	2	-	0/6/23/26	0/1/1/1
7	NAG	E	602	2	-	0/6/23/26	0/1/1/1

All (39) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	E	601	HEM	C4D-ND	-4.73	1.32	1.40
6	I	601	HEM	C1D-ND	-4.14	1.30	1.38
6	G	601	HEM	C1D-ND	-3.68	1.31	1.38
6	B	601	HEM	C4D-ND	-3.27	1.34	1.40
6	I	601	HEM	FE-ND	3.23	2.04	1.94
6	I	601	HEM	C4B-NB	-3.19	1.32	1.38
6	G	601	HEM	C3B-C4B	3.14	1.51	1.44
6	B	601	HEM	C1D-ND	-3.13	1.32	1.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	B	601	HEM	FE-NB	3.13	2.04	1.94
6	I	601	HEM	C4D-ND	-3.09	1.34	1.40
6	I	601	HEM	C3B-C4B	3.07	1.50	1.44
6	E	601	HEM	C3B-C4B	3.00	1.50	1.44
6	G	601	HEM	FE-NB	2.92	2.03	1.94
6	B	601	HEM	C1D-C2D	2.88	1.50	1.44
6	G	601	HEM	C4B-NB	-2.87	1.33	1.38
6	B	601	HEM	C3B-C4B	2.86	1.50	1.44
6	B	601	HEM	CHA-C4D	-2.76	1.33	1.38
6	E	601	HEM	C3D-C2D	-2.75	1.30	1.36
6	E	601	HEM	C1D-ND	-2.65	1.33	1.38
6	I	601	HEM	FE-NB	2.51	2.02	1.94
6	E	601	HEM	C2A-C3A	-2.42	1.32	1.38
9	E	604	XS1	C2-N2	2.35	1.41	1.35
6	E	601	HEM	C1C-NC	-2.34	1.35	1.39
6	G	601	HEM	C1B-NB	-2.33	1.36	1.40
6	B	601	HEM	C4B-NB	-2.32	1.34	1.38
6	I	601	HEM	C4C-NC	-2.30	1.35	1.39
9	B	605	XS1	C2-N2	2.29	1.41	1.35
9	I	605	XS1	C2-N2	2.28	1.41	1.35
9	G	605	XS1	C2-N2	2.25	1.41	1.35
6	G	601	HEM	C4D-ND	-2.21	1.36	1.40
6	G	601	HEM	CHB-C1B	2.20	1.42	1.38
6	E	601	HEM	C4B-NB	-2.19	1.34	1.38
6	E	601	HEM	O2A-CGA	2.18	1.38	1.30
6	G	601	HEM	C3D-C2D	-2.14	1.32	1.36
6	I	601	HEM	C1B-NB	-2.11	1.36	1.40
6	E	601	HEM	FE-ND	2.11	2.01	1.94
6	G	601	HEM	C3C-C2C	2.10	1.41	1.37
6	B	601	HEM	C2A-C3A	-2.09	1.33	1.38
6	G	601	HEM	C1C-NC	-2.02	1.35	1.39

All (59) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
9	B	605	XS1	C1-C5-C4	-3.78	114.21	118.25
9	I	605	XS1	C1-C5-C4	-3.63	114.37	118.25
9	E	604	XS1	C1-C5-C4	-3.61	114.39	118.25
6	G	601	HEM	C2B-C1B-NB	3.43	113.78	109.84
9	G	605	XS1	C1-C5-C4	-3.42	114.59	118.25
6	I	601	HEM	C4A-NA-C1A	-3.41	100.25	105.82
6	G	601	HEM	CBA-CAA-C2A	-3.21	103.67	112.53

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	B	601	HEM	O2A-CGA-O1A	-3.15	115.24	123.33
6	B	601	HEM	O2D-CGD-CBD	3.11	123.84	114.00
6	B	601	HEM	O2A-CGA-CBA	3.08	123.75	114.00
6	I	601	HEM	O2A-CGA-CBA	3.07	123.69	114.00
6	I	601	HEM	O2A-CGA-O1A	-3.05	115.48	123.33
6	E	601	HEM	O2A-CGA-CBA	3.03	123.58	114.00
6	I	601	HEM	C2B-C1B-NB	3.00	113.29	109.84
7	B	602	NAG	C1-O5-C5	2.95	116.14	112.19
7	E	602	NAG	C1-O5-C5	2.94	116.13	112.19
6	B	601	HEM	C4A-NA-C1A	-2.93	101.04	105.82
6	B	601	HEM	O2D-CGD-O1D	-2.93	115.79	123.33
6	I	601	HEM	CBA-CAA-C2A	-2.93	104.43	112.53
7	I	602	NAG	C1-O5-C5	2.93	116.11	112.19
7	G	602	NAG	C1-O5-C5	2.92	116.10	112.19
6	E	601	HEM	C2B-C1B-NB	2.91	113.18	109.84
6	E	601	HEM	CHA-C4D-C3D	-2.84	120.00	125.23
6	E	601	HEM	CBA-CAA-C2A	-2.83	104.72	112.53
6	B	601	HEM	CBA-CAA-C2A	-2.71	105.04	112.53
7	I	603	NAG	C1-O5-C5	2.70	115.81	112.19
6	B	601	HEM	CAD-C3D-C4D	2.68	129.37	124.70
6	G	601	HEM	C4A-NA-C1A	-2.65	101.50	105.82
6	E	601	HEM	C3B-C4B-NB	2.59	111.33	109.47
6	B	601	HEM	CHA-C4D-C3D	-2.54	120.54	125.23
6	I	601	HEM	CHA-C4D-C3D	-2.52	120.57	125.23
6	G	601	HEM	CHA-C4D-C3D	-2.50	120.61	125.23
6	I	601	HEM	C3B-C4B-NB	2.44	111.22	109.47
6	B	601	HEM	C2B-C1B-NB	2.44	112.64	109.84
6	I	601	HEM	C3A-C4A-NA	2.44	114.05	110.14
6	I	601	HEM	C3B-C2B-C1B	-2.36	104.64	106.41
6	E	601	HEM	C4A-NA-C1A	-2.34	102.00	105.82
6	E	601	HEM	O2A-CGA-O1A	-2.34	117.31	123.33
6	B	601	HEM	CHD-C1D-ND	2.33	126.93	124.42
6	G	601	HEM	O2A-CGA-CBA	2.29	121.23	114.00
6	G	601	HEM	O2D-CGD-CBD	2.27	121.16	114.00
6	G	601	HEM	O2A-CGA-O1A	-2.25	117.55	123.33
9	I	605	XS1	N3-C3-N1	2.23	127.48	125.64
6	I	601	HEM	C4C-C3C-C2C	-2.23	104.88	106.81
6	G	601	HEM	CAD-C3D-C4D	2.21	128.55	124.70
9	G	605	XS1	N3-C3-N1	2.20	127.45	125.64
6	G	601	HEM	C3B-C4B-NB	2.20	111.05	109.47
6	G	601	HEM	O2D-CGD-O1D	-2.14	117.83	123.33
9	B	605	XS1	N3-C3-N1	2.12	127.38	125.64

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	G	601	HEM	C2A-C1A-NA	2.12	112.50	110.15
6	I	601	HEM	CAD-C3D-C4D	2.11	128.38	124.70
9	E	604	XS1	N3-C3-N1	2.10	127.37	125.64
6	E	601	HEM	O2D-CGD-CBD	2.09	120.61	114.00
6	G	601	HEM	CAA-C2A-C1A	2.07	128.98	124.94
6	I	601	HEM	C4C-CHD-C1D	-2.04	121.67	126.02
6	G	601	HEM	CHB-C1B-C2B	-2.04	121.14	126.95
6	E	601	HEM	C4C-C3C-C2C	-2.04	105.04	106.81
6	G	601	HEM	C3C-C2C-C1C	-2.01	105.14	107.05
6	G	601	HEM	C3B-C2B-C1B	-2.01	104.91	106.41

There are no chirality outliers.

All (33) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
9	B	605	XS1	C11-C13-C6-C7
9	E	604	XS1	C11-C13-C6-C7
9	E	604	XS1	C22-C23-C24-C25
9	E	604	XS1	C10-C23-C24-C25
9	G	605	XS1	C11-C13-C6-C7
9	G	605	XS1	C22-C23-C24-C25
9	G	605	XS1	C10-C23-C24-C25
9	I	605	XS1	C22-C23-C24-C25
9	I	605	XS1	C10-C23-C24-C25
9	I	605	XS1	C1-C5-C6-C7
9	I	605	XS1	C11-C13-C6-C7
9	B	605	XS1	C1-C5-C6-C7
9	E	604	XS1	C1-C5-C6-C7
9	B	605	XS1	C4-C5-C6-C7
9	E	604	XS1	C4-C5-C6-C7
9	G	605	XS1	C4-C5-C6-C7
9	I	605	XS1	C4-C5-C6-C7
6	G	601	HEM	CAD-CBD-CGD-O1D
6	B	601	HEM	CAD-CBD-CGD-O1D
6	E	601	HEM	CAA-CBA-CGA-O2A
6	I	601	HEM	CAA-CBA-CGA-O1A
6	I	601	HEM	CAD-CBD-CGD-O1D
6	G	601	HEM	CAD-CBD-CGD-O2D
6	E	601	HEM	CAD-CBD-CGD-O1D
6	E	601	HEM	CAD-CBD-CGD-O2D
6	B	601	HEM	CAD-CBD-CGD-O2D
6	I	601	HEM	CAA-CBA-CGA-O2A

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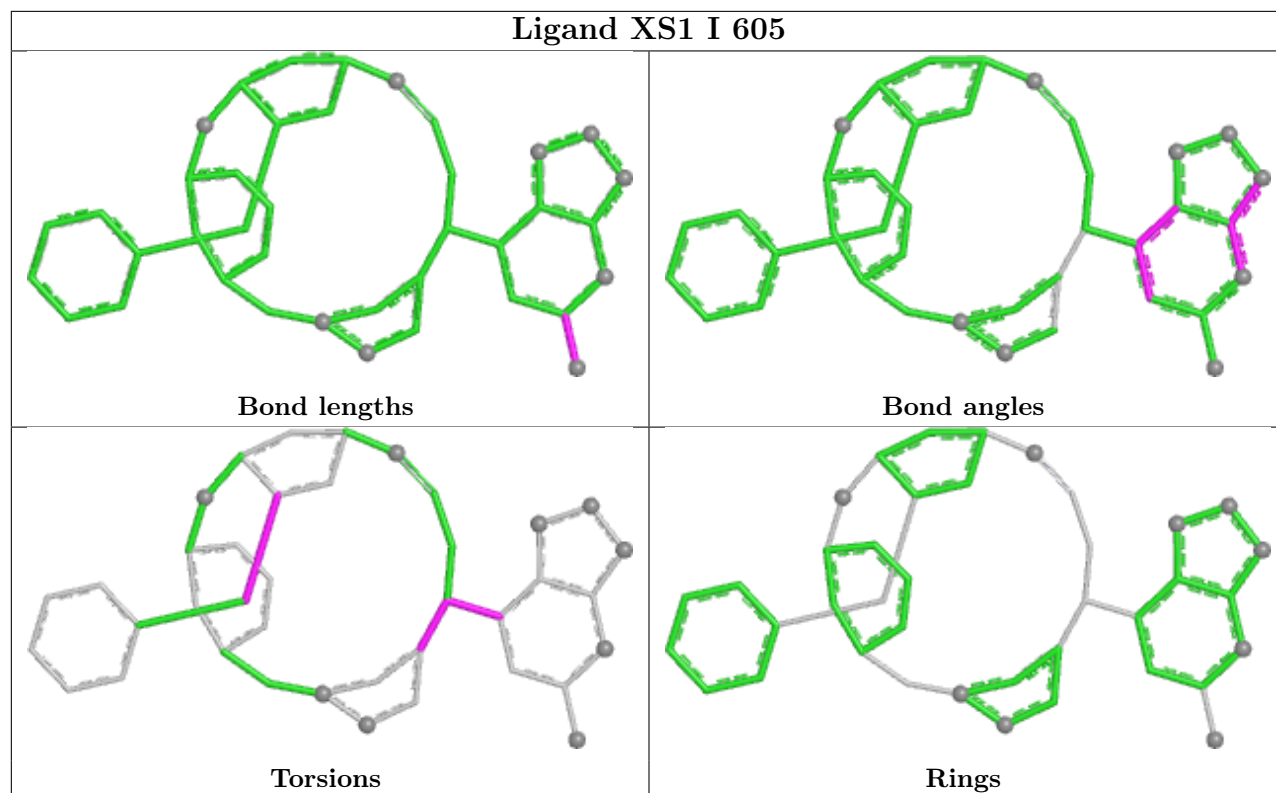
Mol	Chain	Res	Type	Atoms
6	I	601	HEM	CAD-CBD-CGD-O2D
6	B	601	HEM	CAA-CBA-CGA-O1A
6	E	601	HEM	CAA-CBA-CGA-O1A
6	B	601	HEM	CAA-CBA-CGA-O2A
6	G	601	HEM	CAA-CBA-CGA-O1A
6	G	601	HEM	CAA-CBA-CGA-O2A

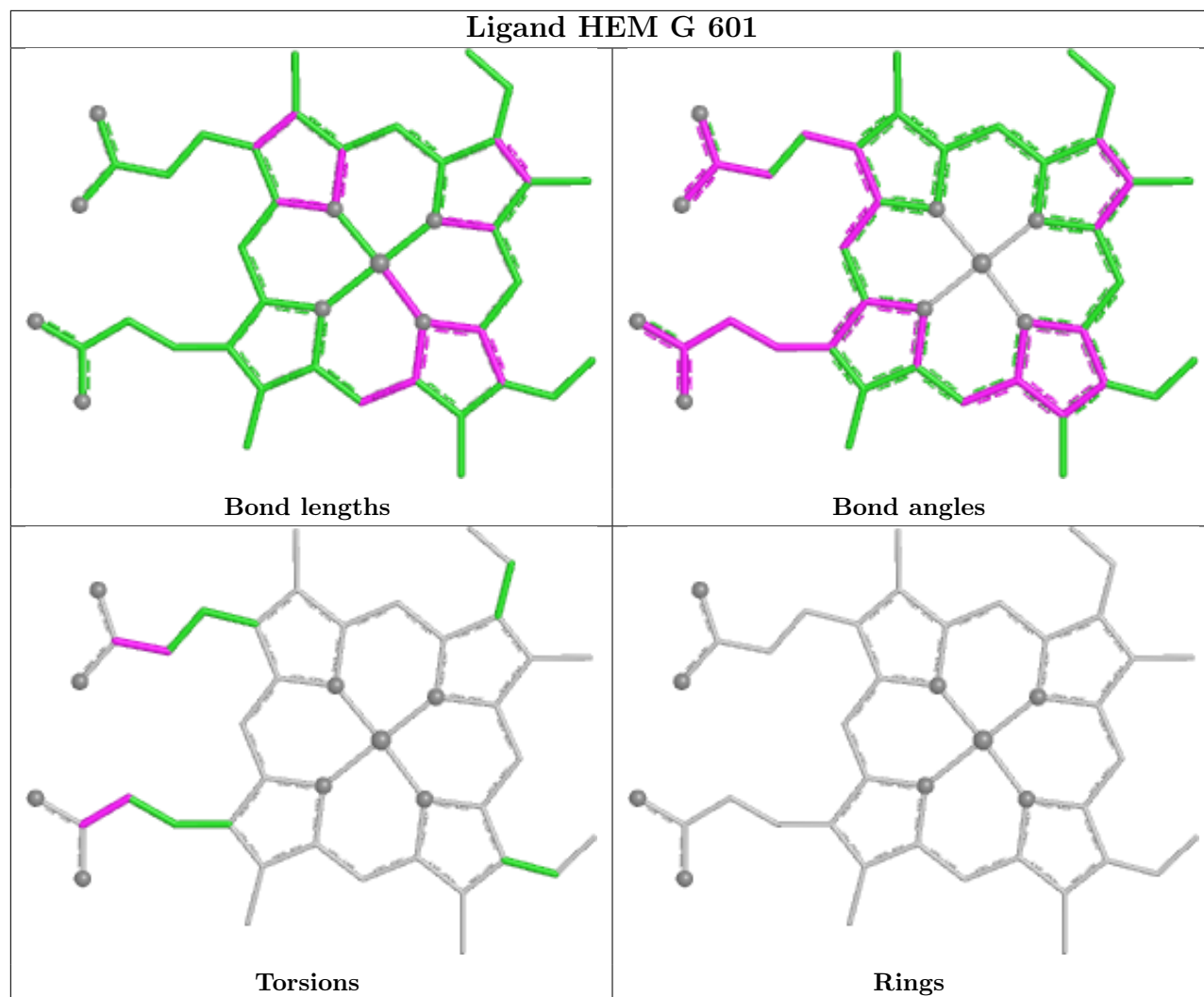
There are no ring outliers.

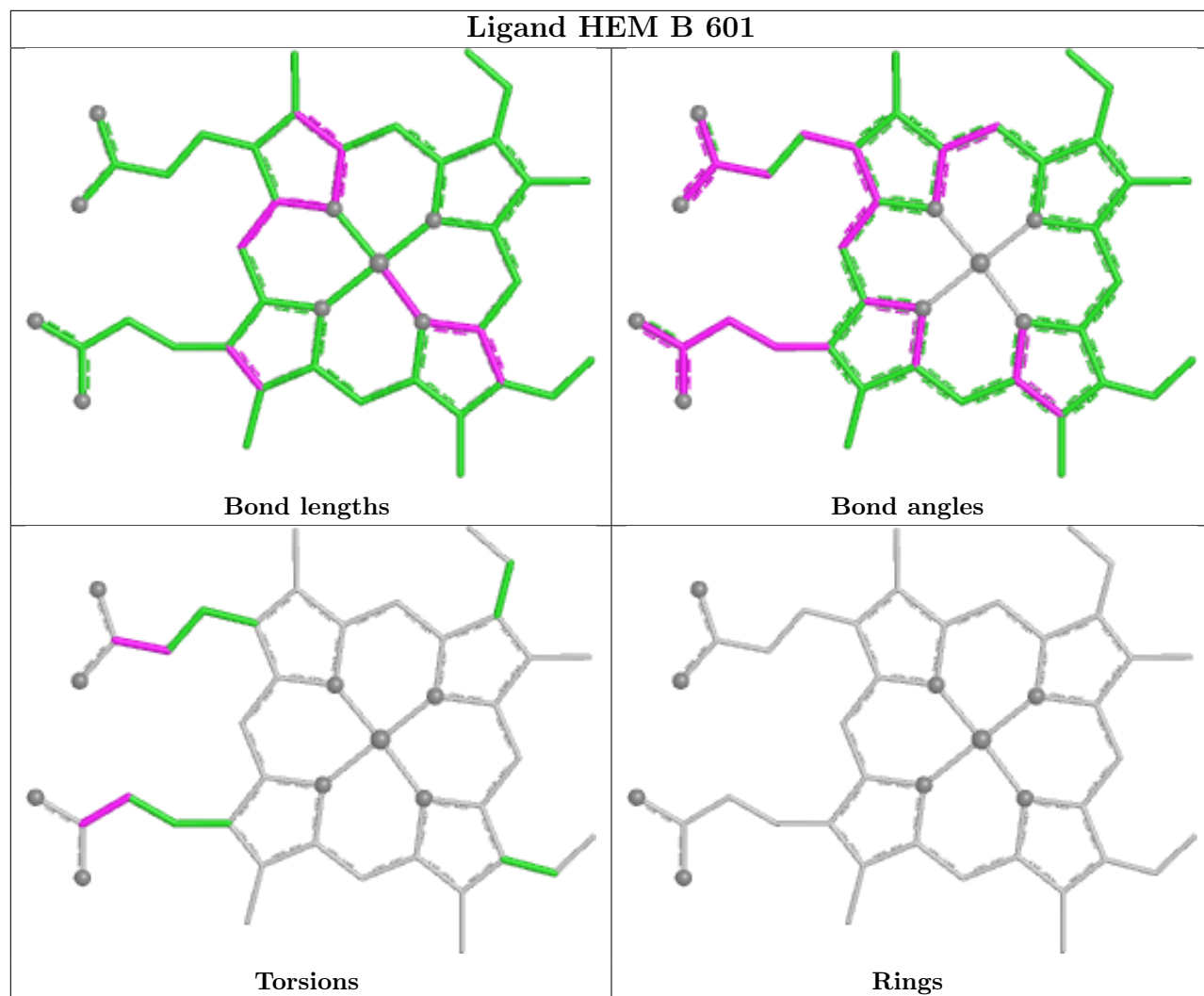
4 monomers are involved in 36 short contacts:

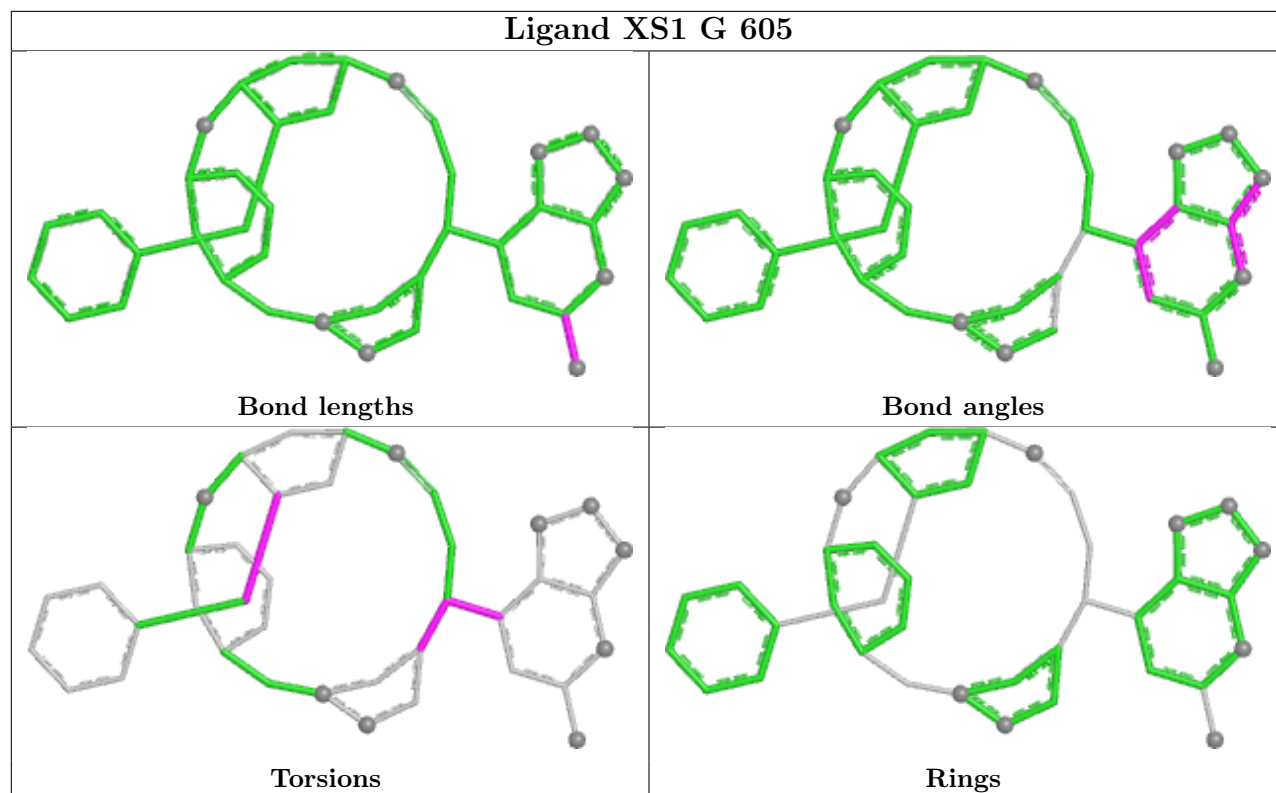
Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	G	601	HEM	9	0
6	B	601	HEM	9	0
6	E	601	HEM	9	0
6	I	601	HEM	9	0

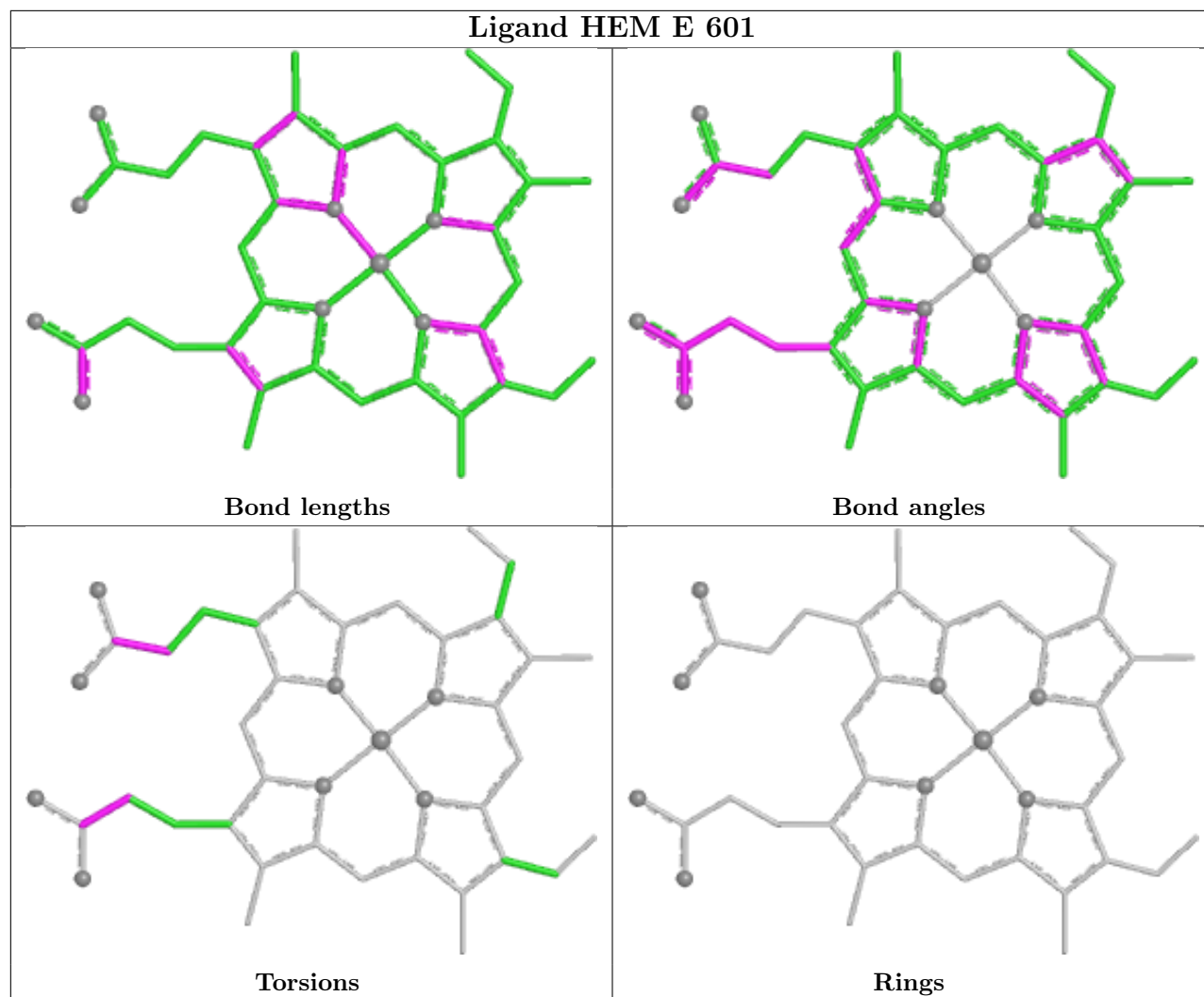
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

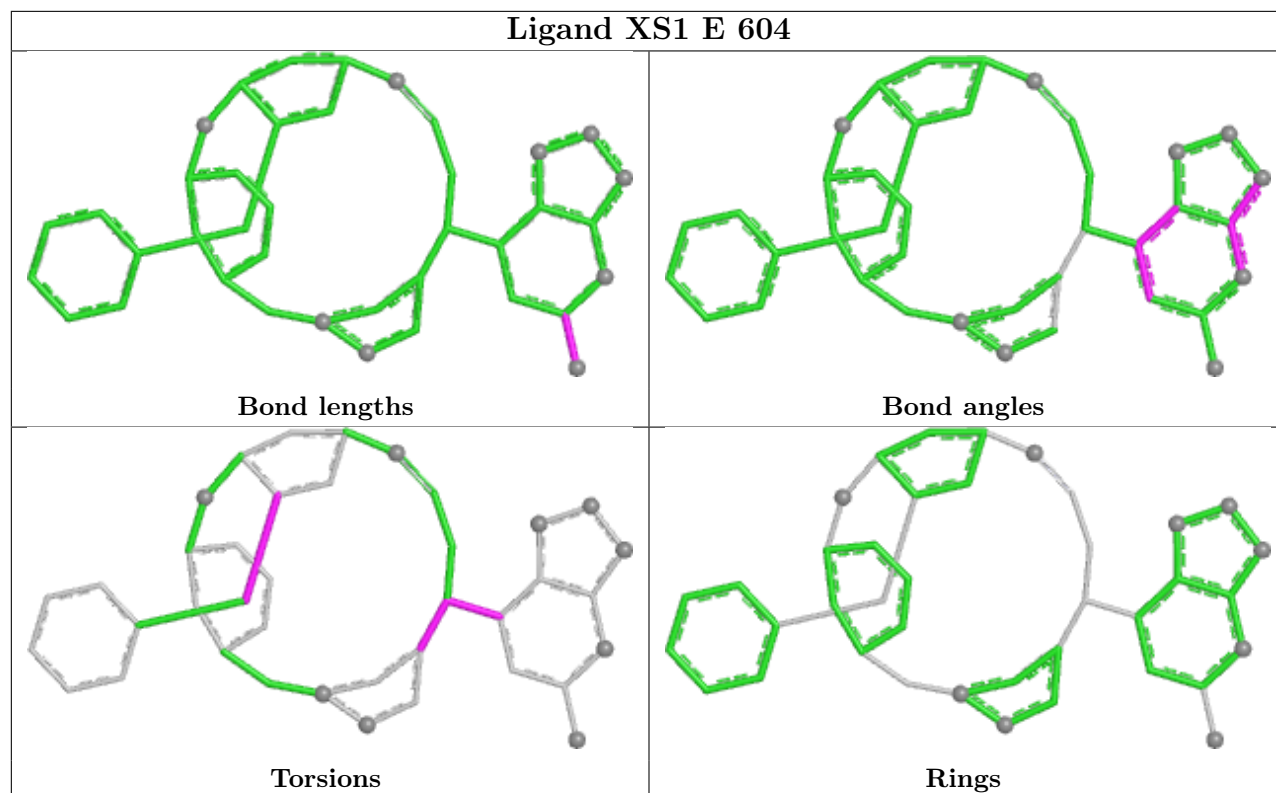




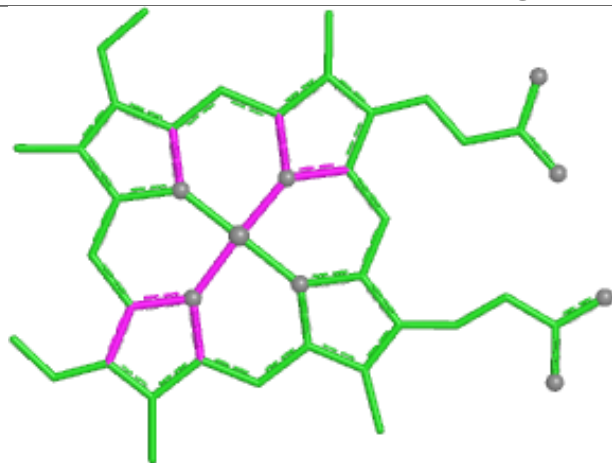




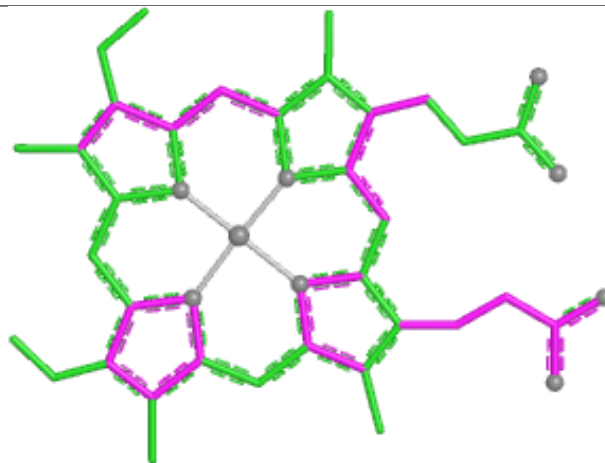




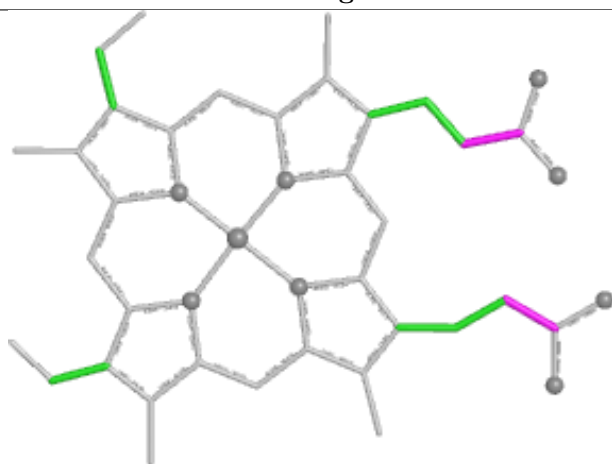
Ligand HEM I 601



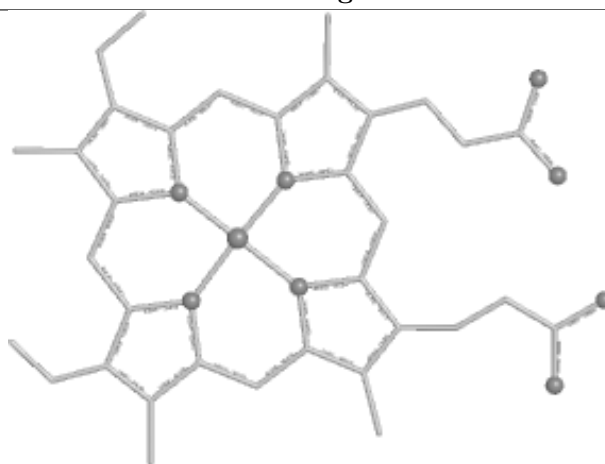
Bond lengths



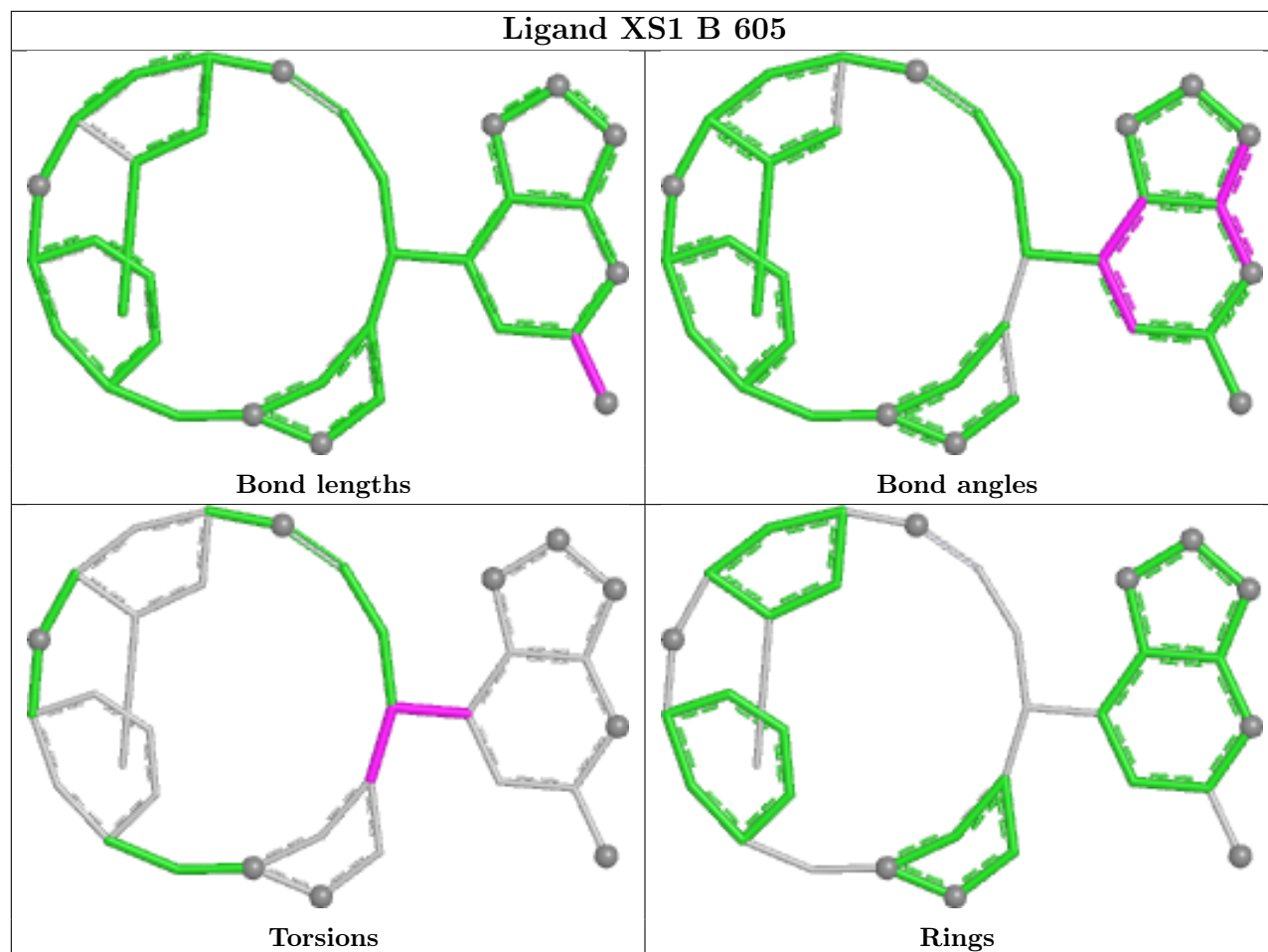
Bond angles



Torsions



Rings



5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 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	103/105 (98%)	0.01	2 (1%) 66 68	29, 41, 65, 86	1 (0%)
1	D	104/105 (99%)	-0.11	2 (1%) 66 68	25, 37, 60, 85	2 (1%)
1	F	103/105 (98%)	-0.11	1 (0%) 79 81	27, 37, 63, 81	1 (0%)
1	H	103/105 (98%)	-0.04	0 100 100	27, 38, 60, 70	0
2	B	464/466 (99%)	0.31	18 (3%) 43 45	25, 46, 69, 96	9 (1%)
2	E	465/466 (99%)	-0.06	5 (1%) 78 79	20, 40, 61, 82	12 (2%)
2	G	464/466 (99%)	-0.01	4 (0%) 81 82	21, 41, 63, 90	9 (1%)
2	I	464/466 (99%)	0.17	9 (1%) 66 68	23, 44, 65, 79	11 (2%)
All	All	2270/2284 (99%)	0.07	41 (1%) 67 69	20, 42, 66, 96	45 (1%)

All (41) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	195	GLY	3.4
2	I	348	ASN	3.2
2	E	567	LEU	3.2
2	B	114	ASN	3.1
2	B	190	MET	3.0
2	B	194	LEU	3.0
2	I	354	GLU	2.8
2	B	218	ASP	2.8
2	G	218	ASP	2.8
2	I	353	MET	2.6
2	B	267	LEU	2.6
2	I	355	PRO	2.6
2	G	217	HIS	2.6
1	A	4	GLN	2.5
2	B	191	SER	2.5
1	A	1	CYS	2.5

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Mol	Chain	Res	Type	RSRZ
2	E	577	GLU	2.5
1	F	1	CYS	2.4
2	B	372	VAL	2.4
2	I	577	GLU	2.4
1	D	104	ALA	2.4
2	I	218	ASP	2.3
2	B	542	GLY	2.3
2	B	522	MET	2.3
2	E	348	ASN	2.3
2	I	411	MET	2.3
2	B	202	ARG	2.3
2	E	113	VAL	2.3
1	D	1	CYS	2.3
2	G	522	MET	2.3
2	I	114	ASN	2.2
2	B	185	ARG	2.2
2	B	566	THR	2.2
2	B	264	LEU	2.1
2	B	263	GLU	2.1
2	I	217	HIS	2.1
2	B	181	GLU	2.1
2	B	572	LEU	2.1
2	E	206	ASN	2.0
2	G	348	ASN	2.0
2	B	270	ARG	2.0

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

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

6.3 Carbohydrates ⓘ

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 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	C	1	14/15	-	-	79,84,87,90	0
3	NAG	C	2	14/15	-	-	93,95,97,97	0
3	NAG	K	2	14/15	0.56	0.16	95,101,103,104	0

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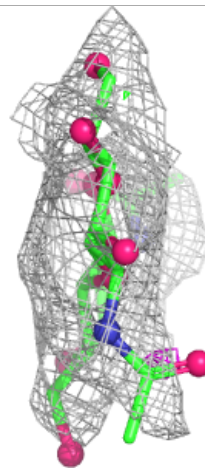
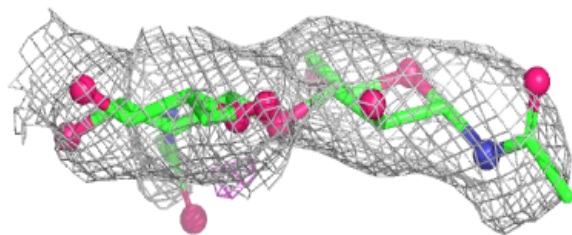
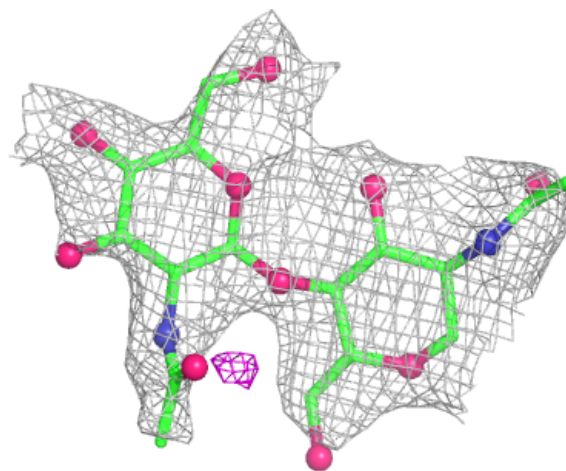
Continued from previous page...

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	NAG	M	2	14/15	0.75	0.11	75,79,84,85	0
4	MAN	L	4	11/12	0.76	0.12	61,65,75,79	0
4	MAN	O	4	11/12	0.79	0.11	54,61,65,67	0
4	MAN	J	4	11/12	0.82	0.11	53,58,64,68	0
4	MAN	N	4	11/12	0.85	0.10	53,59,63,68	0
3	NAG	K	1	14/15	0.89	0.12	75,80,84,88	0
3	NAG	M	1	14/15	0.90	0.09	59,63,69,76	0
4	MAN	N	5	11/12	0.93	0.08	38,43,45,49	0
4	MAN	L	5	11/12	0.93	0.07	40,42,45,48	0
4	MAN	J	5	11/12	0.94	0.06	36,38,41,42	0
4	BMA	J	3	11/12	0.94	0.07	32,36,40,46	0
4	NAG	J	1	14/15	0.94	0.07	32,34,39,46	0
4	MAN	O	5	11/12	0.94	0.06	38,40,45,45	0
4	BMA	N	3	11/12	0.95	0.06	36,41,46,47	0
4	NAG	L	1	14/15	0.95	0.06	31,39,44,44	0
4	BMA	L	3	11/12	0.95	0.06	35,38,44,53	0
4	FUC	N	6	10/11	0.95	0.07	34,35,38,38	0
4	FUC	L	6	10/11	0.95	0.07	38,42,44,44	0
4	NAG	N	1	14/15	0.95	0.07	32,35,42,44	0
4	FUC	O	6	10/11	0.95	0.06	37,40,41,42	0
4	FUC	J	6	10/11	0.96	0.07	37,43,44,44	0
4	NAG	N	2	14/15	0.96	0.06	23,29,34,37	0
4	NAG	J	2	14/15	0.96	0.05	25,33,38,40	0
4	BMA	O	3	11/12	0.97	0.05	34,38,42,48	0
4	NAG	L	2	14/15	0.97	0.05	25,31,34,37	0
4	NAG	O	1	14/15	0.97	0.06	35,38,44,48	0
4	NAG	O	2	14/15	0.97	0.05	24,33,37,38	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.

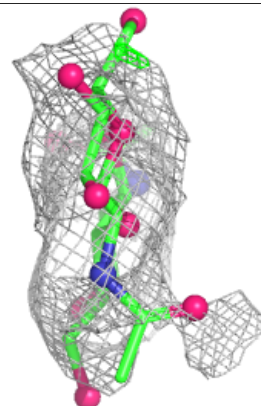
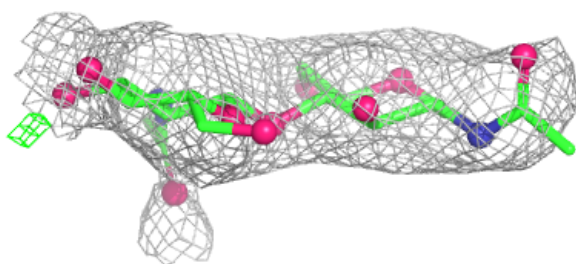
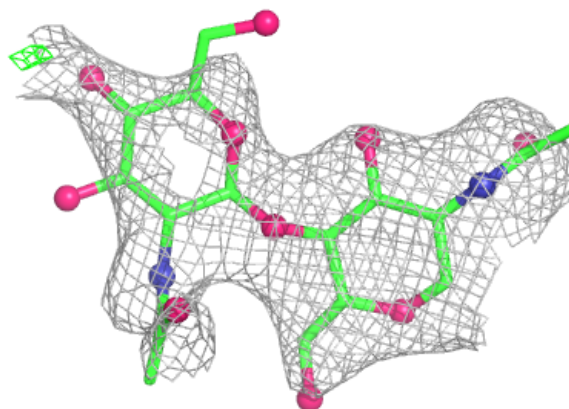
Electron density around Chain C:

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



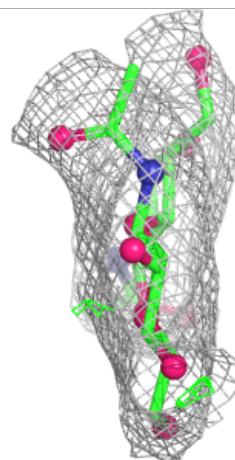
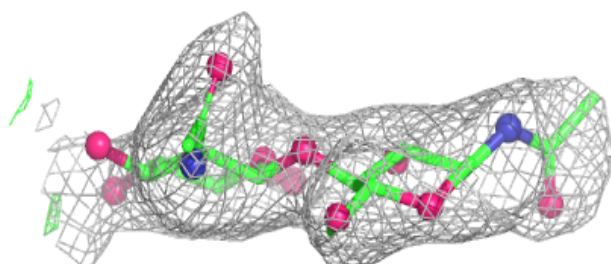
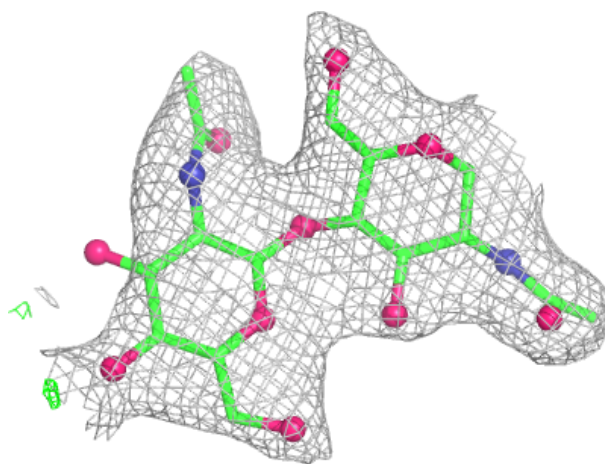
Electron density around Chain K:

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



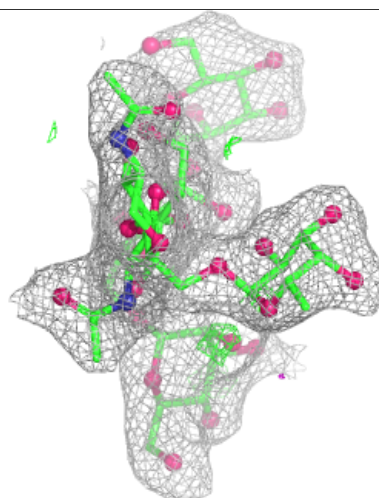
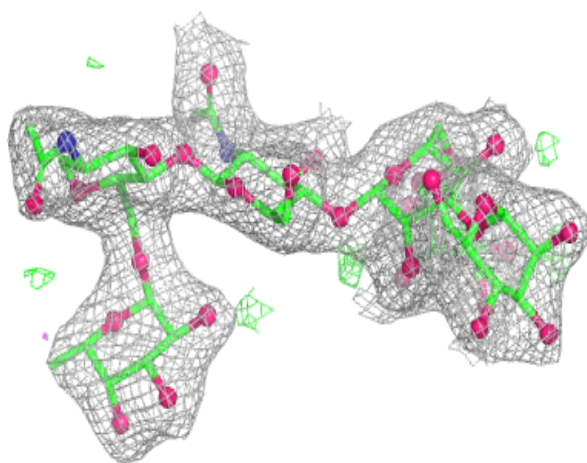
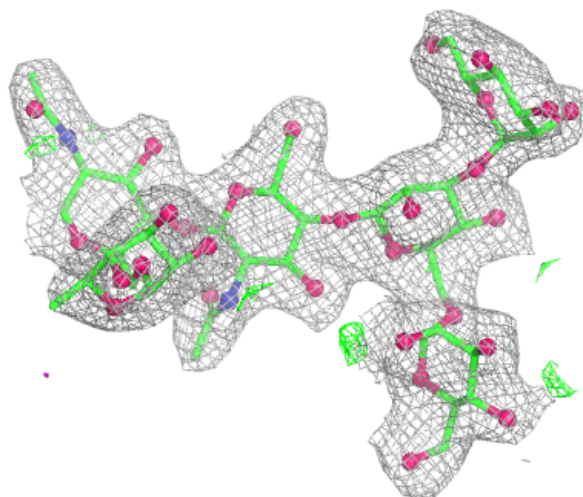
Electron density around Chain M:

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



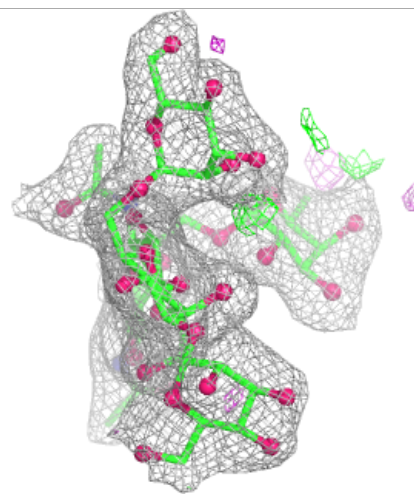
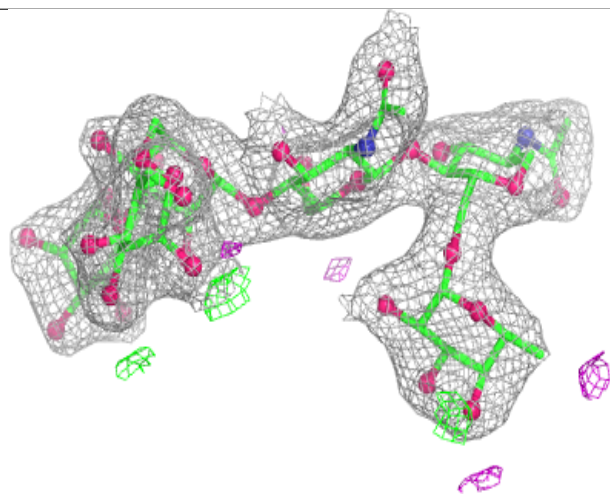
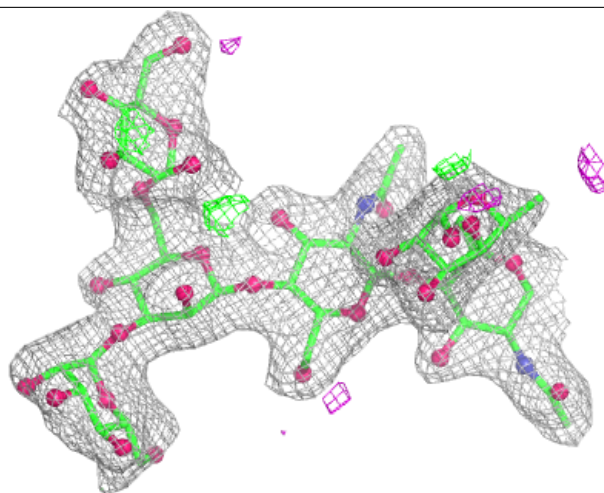
Electron density around Chain J:

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



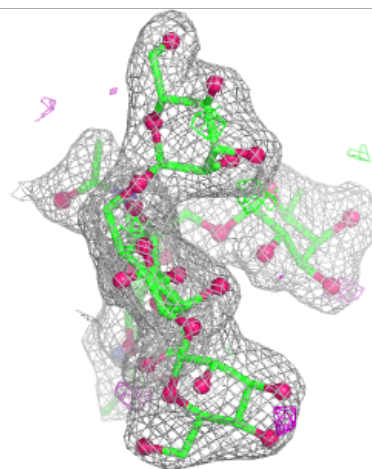
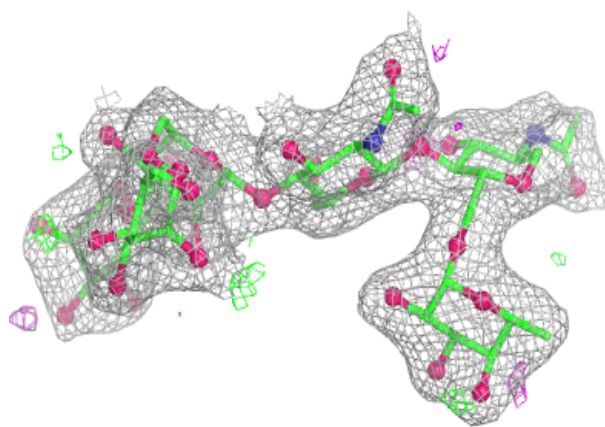
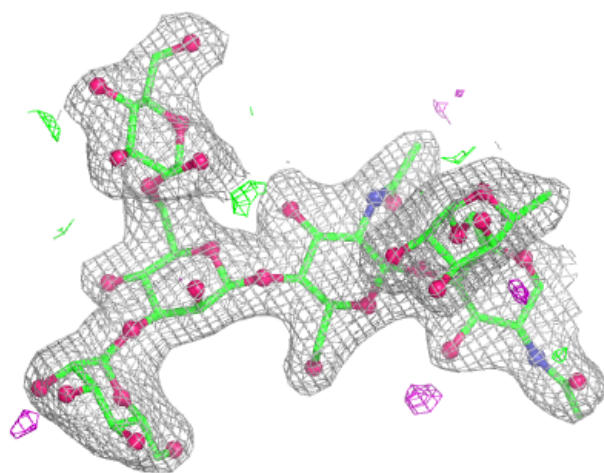
Electron density around Chain L:

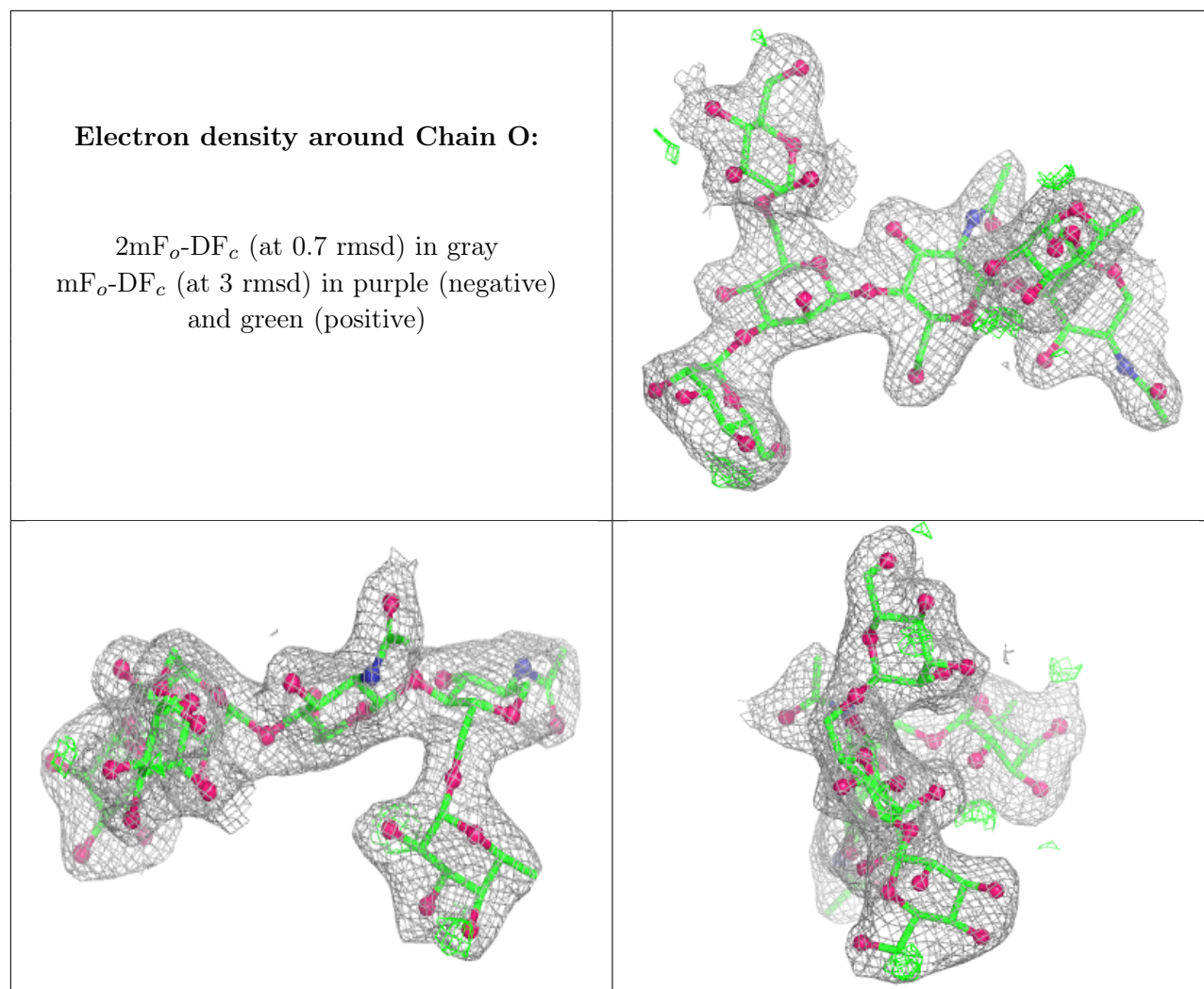
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around Chain N:

$2mF_o-DF_c$ (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(Å ²)	Q<0.9
7	NAG	G	602	14/15	0.74	0.12	55,65,70,70	0
7	NAG	I	603	14/15	0.74	0.15	89,91,93,93	0
7	NAG	B	602	14/15	0.75	0.12	49,69,72,73	0
7	NAG	E	602	14/15	0.82	0.11	47,57,64,64	0
7	NAG	I	602	14/15	0.83	0.11	57,61,63,65	0
5	CL	B	604	1/1	0.89	0.26	65,65,65,65	0
9	XS1	B	605	33/39	0.90	0.08	42,46,52,53	27
9	XS1	I	605	39/39	0.91	0.10	37,42,58,60	32
9	XS1	G	605	39/39	0.93	0.08	26,36,46,48	32

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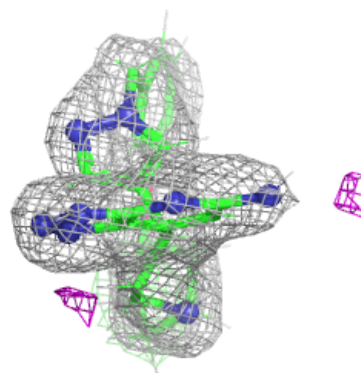
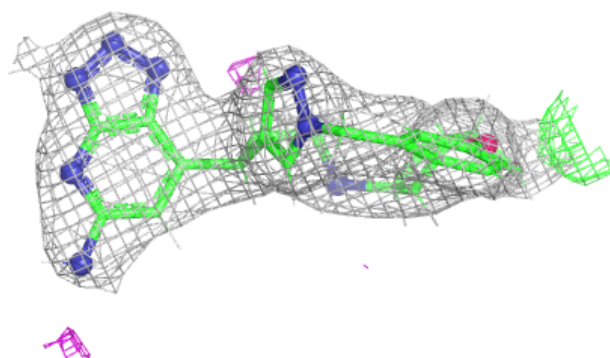
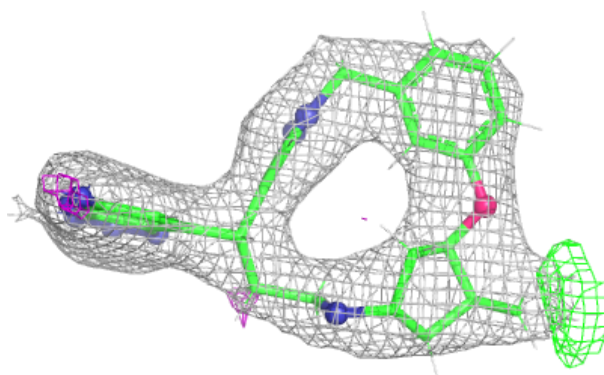
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
9	XS1	E	604	39/39	0.94	0.07	29,36,50,51	32
6	HEM	B	601	43/43	0.97	0.08	33,34,38,48	0
6	HEM	G	601	43/43	0.97	0.08	33,34,37,47	0
6	HEM	I	601	43/43	0.97	0.07	35,36,40,46	0
5	CL	G	604	1/1	0.97	0.26	58,58,58,58	0
6	HEM	E	601	43/43	0.98	0.06	30,31,36,45	0
5	CL	F	201	1/1	0.98	0.06	30,30,30,30	0
5	CL	A	201	1/1	0.99	0.03	30,30,30,30	0
8	CA	E	603	1/1	0.99	0.04	30,30,30,30	0
8	CA	G	603	1/1	0.99	0.06	30,30,30,30	0
8	CA	I	604	1/1	0.99	0.03	35,35,35,35	0
8	CA	B	603	1/1	1.00	0.02	38,38,38,38	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.

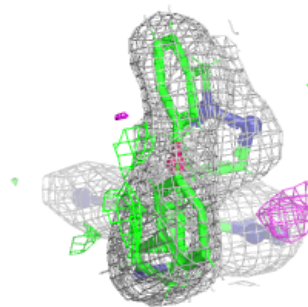
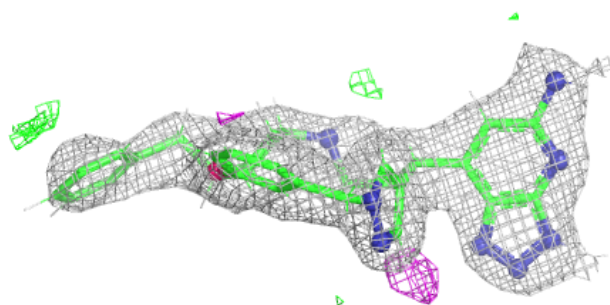
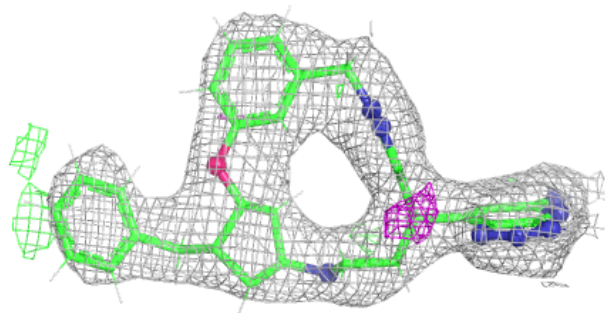
Electron density around XS1 B 605:

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

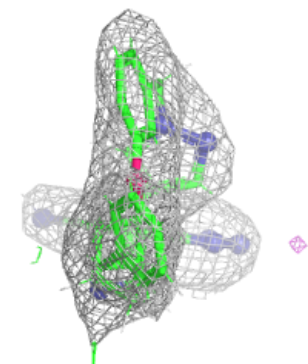
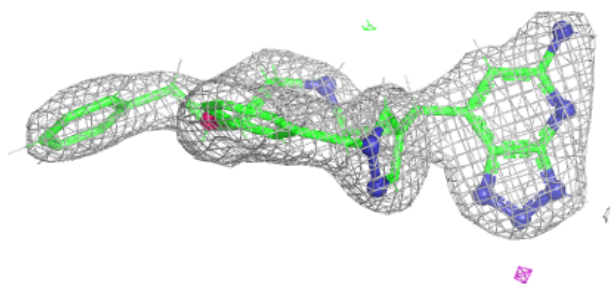
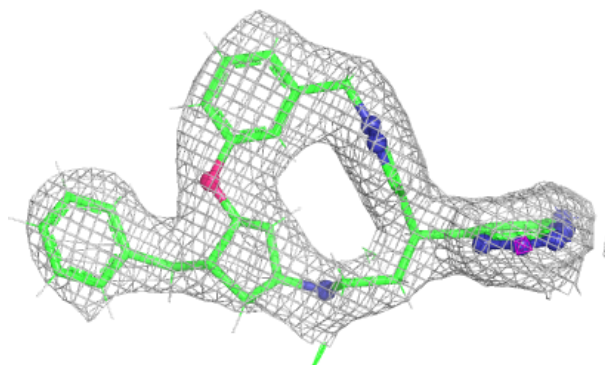


Electron density around XS1 I 605:

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

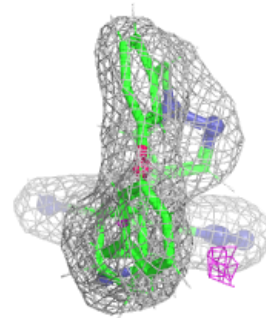
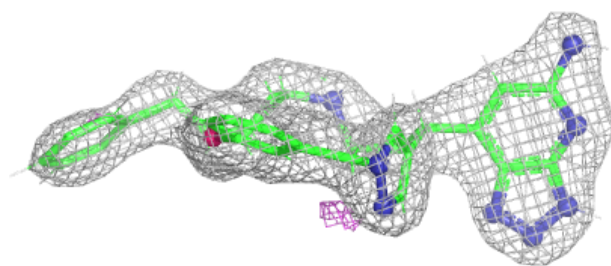
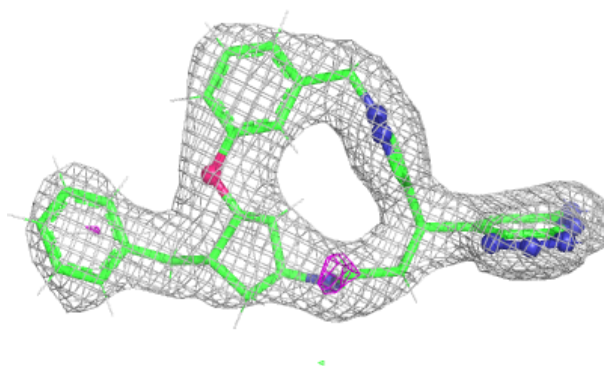
**Electron density around XS1 G 605:**

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



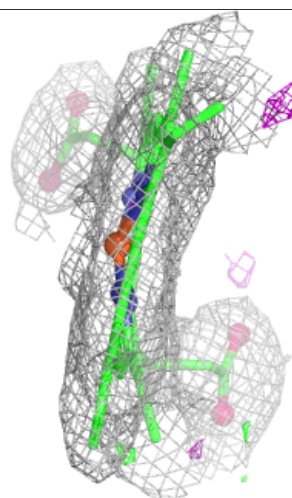
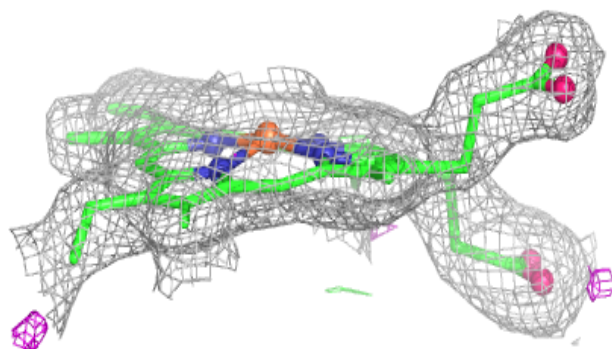
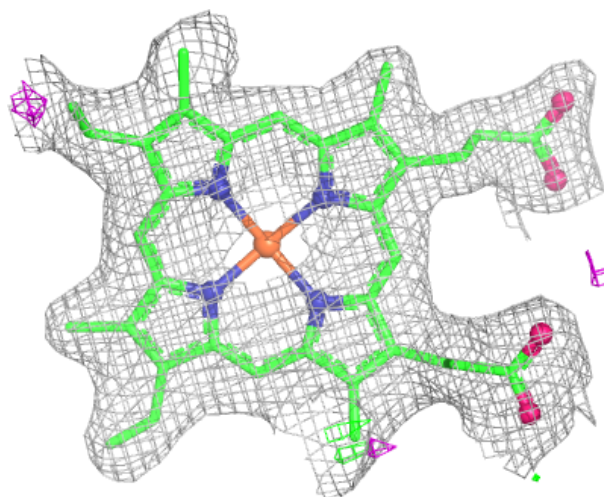
Electron density around XS1 E 604:

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



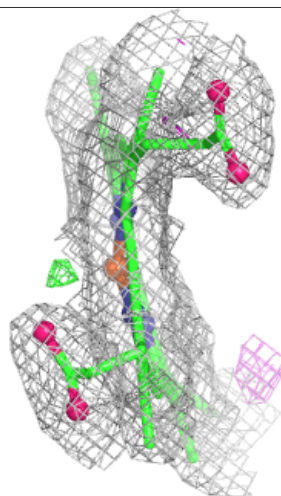
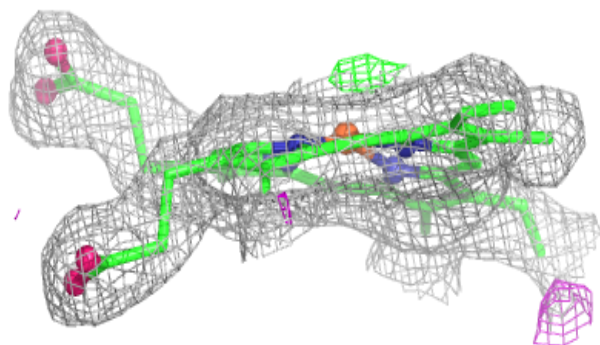
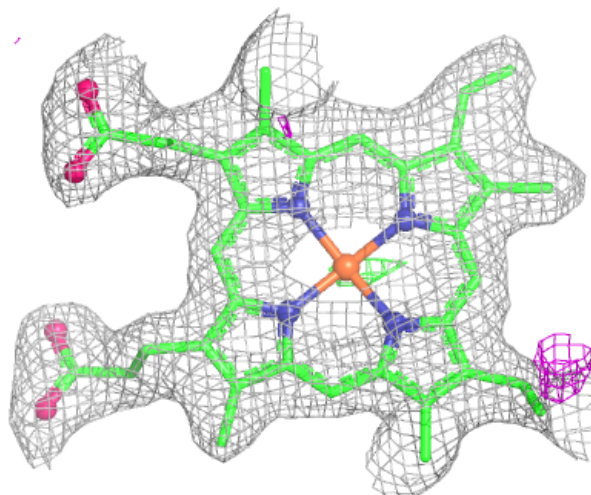
Electron density around HEM B 601:

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



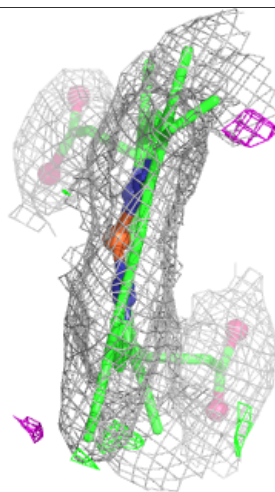
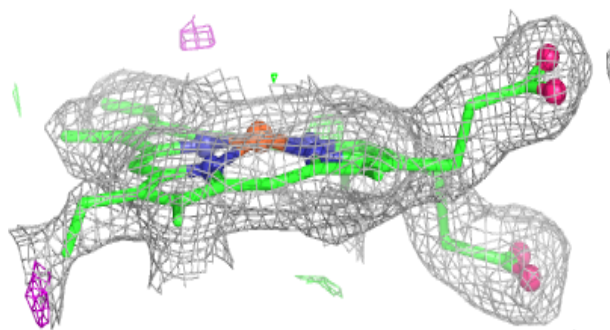
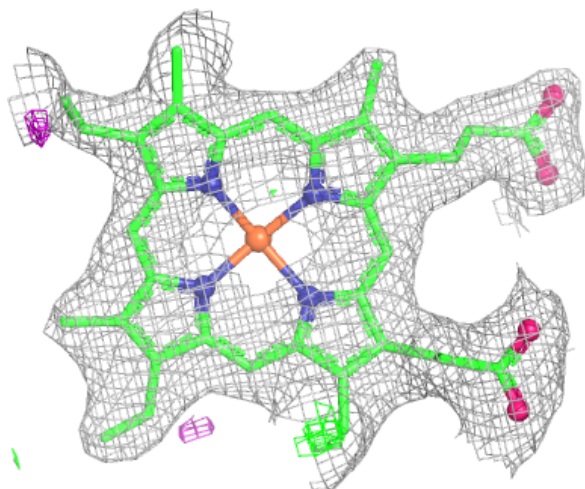
Electron density around HEM G 601:

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



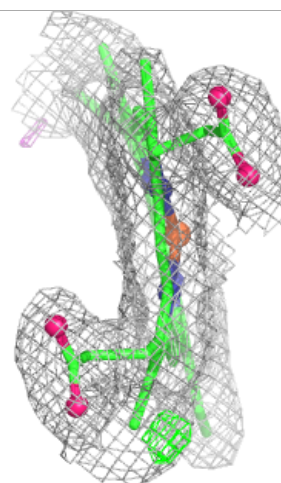
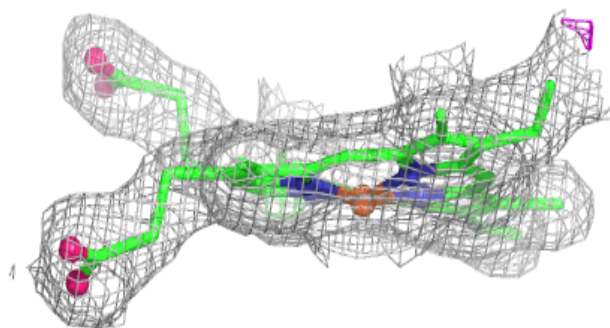
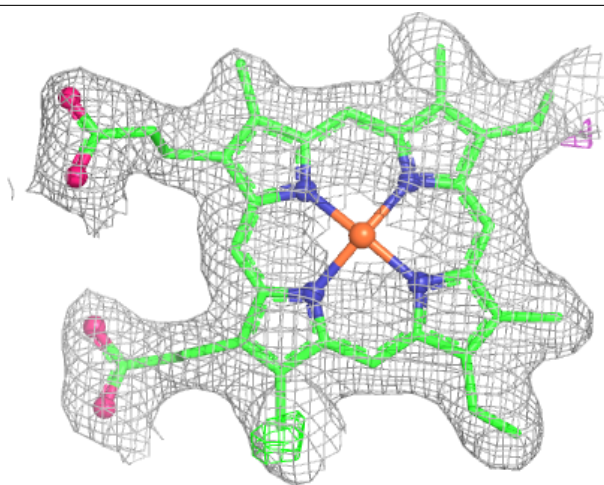
Electron density around HEM I 601:

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



Electron density around HEM E 601:

$2mF_o-DF_c$ (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.