



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 12:18 PM UTC

PDB ID : 3CZH / pdb_00003czh
Title : Crystal structure of CYP2R1 in complex with vitamin D2
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Edwards, A.M.; Bountra, C.; Wilkstrom, M.; Bochkarev, A.; Park, H.; Struc-
tural Genomics Consortium (SGC)
Deposited on : 2008-04-29
Resolution : 2.30 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

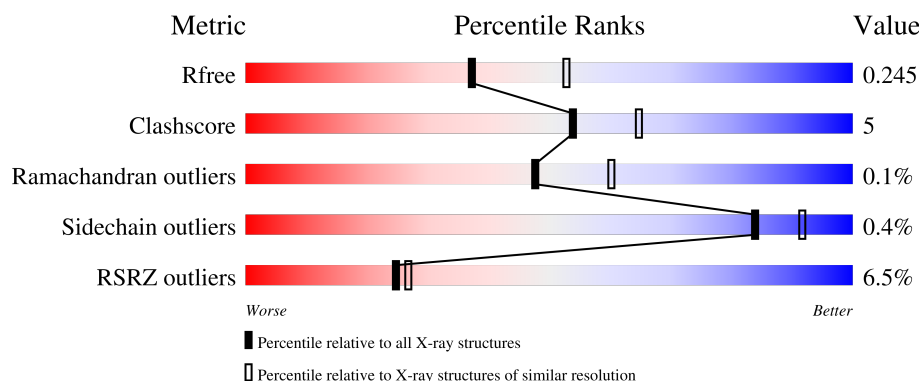
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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	6319 (2.30-2.30)
Clashscore	190562	6919 (2.30-2.30)
Ramachandran outliers	187476	6854 (2.30-2.30)
Sidechain outliers	187428	6854 (2.30-2.30)
RSRZ outliers	180081	6325 (2.30-2.30)

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	481	5% 91% 6%
1	B	481	8% 89% 7%
2	C	7	14% 86%
2	D	7	100%

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	UNX	A	1	-	-	-	X
3	UNX	A	16	-	-	-	X
3	UNX	A	17	-	-	-	X
3	UNX	A	19	-	-	-	X
3	UNX	A	3	-	-	-	X
3	UNX	A	4	-	-	-	X
3	UNX	A	5	-	-	-	X
3	UNX	A	6	-	-	-	X
3	UNX	A	7	-	-	-	X
3	UNX	A	8	-	-	-	X
3	UNX	B	10	-	-	-	X
3	UNX	B	11	-	-	-	X
3	UNX	B	12	-	-	-	X
3	UNX	B	13	-	-	-	X
3	UNX	B	14	-	-	-	X
3	UNX	B	15	-	-	-	X
3	UNX	B	18	-	-	-	X
3	UNX	B	2	-	-	-	X
3	UNX	B	20	-	-	-	X
3	UNX	B	21	-	-	-	X
3	UNX	B	9	-	-	-	X

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 7993 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.

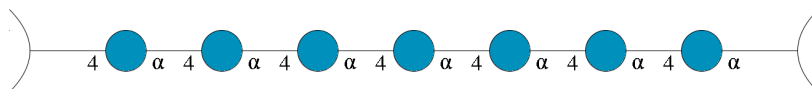
- Molecule 1 is a protein called Cytochrome P450 2R1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	465	Total	C	N	O	S	0	0	0
			3712	2419	611	664	18			
1	B	467	Total	C	N	O	S	0	0	0
			3732	2429	625	660	18			

There are 22 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	27	MET	-	expression tag	UNP Q6VVX0
A	28	ALA	-	expression tag	UNP Q6VVX0
A	29	LYS	-	expression tag	UNP Q6VVX0
A	30	LYS	-	expression tag	UNP Q6VVX0
A	31	THR	-	expression tag	UNP Q6VVX0
A	502	HIS	-	expression tag	UNP Q6VVX0
A	503	HIS	-	expression tag	UNP Q6VVX0
A	504	HIS	-	expression tag	UNP Q6VVX0
A	505	HIS	-	expression tag	UNP Q6VVX0
A	506	HIS	-	expression tag	UNP Q6VVX0
A	507	HIS	-	expression tag	UNP Q6VVX0
B	27	MET	-	expression tag	UNP Q6VVX0
B	28	ALA	-	expression tag	UNP Q6VVX0
B	29	LYS	-	expression tag	UNP Q6VVX0
B	30	LYS	-	expression tag	UNP Q6VVX0
B	31	THR	-	expression tag	UNP Q6VVX0
B	502	HIS	-	expression tag	UNP Q6VVX0
B	503	HIS	-	expression tag	UNP Q6VVX0
B	504	HIS	-	expression tag	UNP Q6VVX0
B	505	HIS	-	expression tag	UNP Q6VVX0
B	506	HIS	-	expression tag	UNP Q6VVX0
B	507	HIS	-	expression tag	UNP Q6VVX0

- Molecule 2 is an oligosaccharide called Cycloheptakis-(1-4)-(alpha-D-glucopyranose).

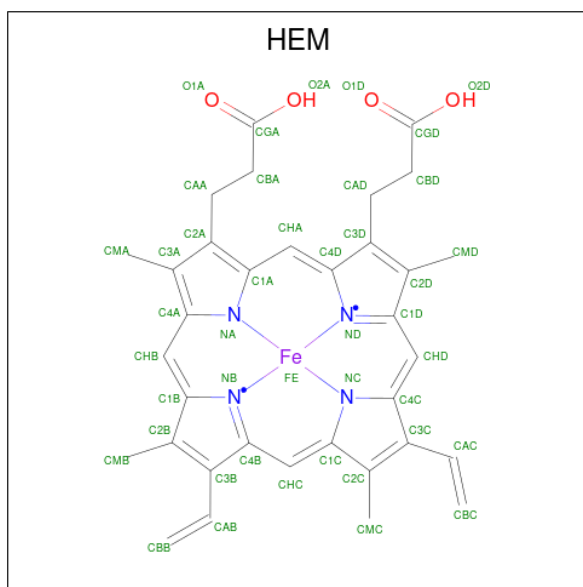


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace
2	C	7	Total	C	O	0	0	0
			77	42	35			
2	D	7	Total	C	O	0	0	0
			77	42	35			

- Molecule 3 is UNKNOWN ATOM OR ION (CCD ID: UNX) (formula: X).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	10	Total	X	0	0
			10	10		
3	B	11	Total	X	0	0
			11	11		

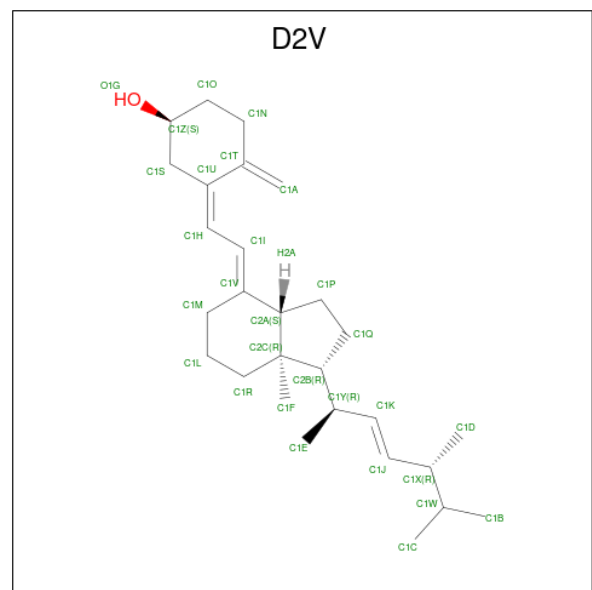
- Molecule 4 is PROTOPORPHYRIN IX CONTAINING FE (CCD ID: HEM) (formula: $C_{34}H_{32}FeN_4O_4$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total	C	Fe	N	O	0	1
			86	68	2	8	8		
4	B	1	Total	C	Fe	N	O	0	1
			86	68	2	8	8		

- Molecule 5 is (3S,5Z,7E,22E)-9,10-secoergosta-5,7,10,22-tetraen-3-ol (CCD ID: D2V)

(formula: C₂₈H₄₄O).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			29	28	1		
5	B	1	Total	C	O	0	0
			29	28	1		

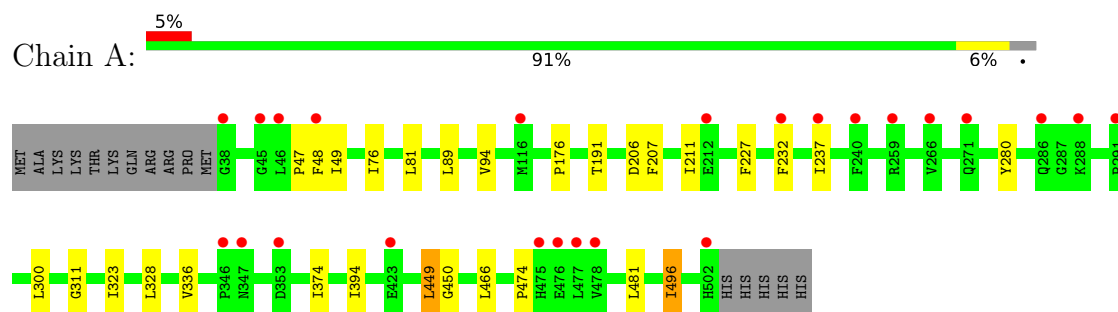
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	86	Total	O	0	0
			86	86		
6	B	58	Total	O	0	0
			58	58		

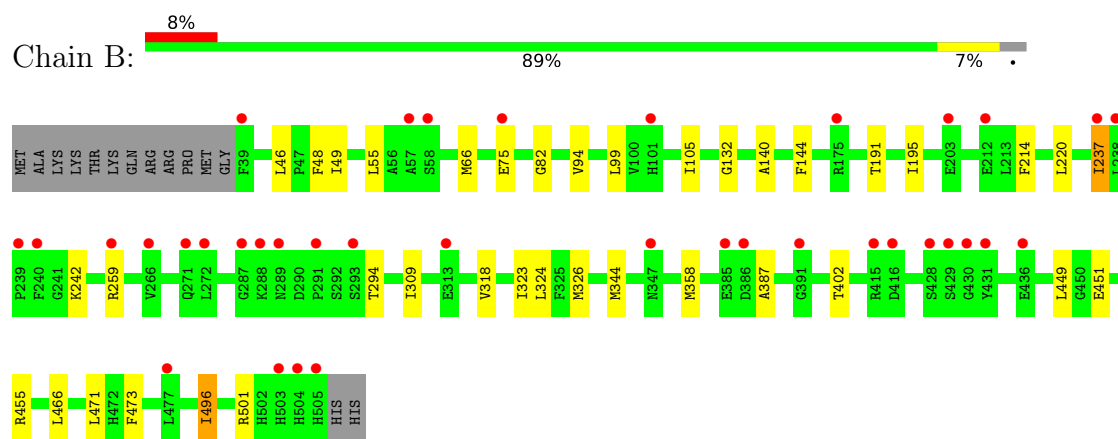
3 Residue-property plots [i](#)

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: Cytochrome P450 2R1



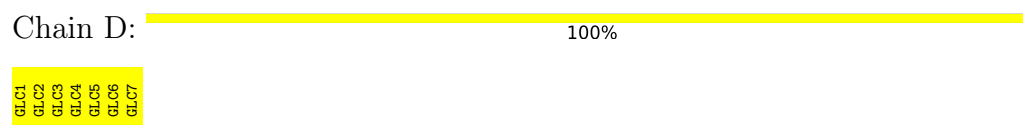
- Molecule 1: Cytochrome P450 2R1



- Molecule 2: Cycloheptakis-(1-4)-(alpha-D-glucopyranose)



- Molecule 2: Cycloheptakis-(1-4)-(alpha-D-glucopyranose)



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	137.36Å 163.39Å 152.29Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.00 – 2.30 30.00 – 2.30	Depositor EDS
% Data completeness (in resolution range)	96.7 (30.00-2.30) 96.6 (30.00-2.30)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.04 (at 2.31Å)	Xtriage
Refinement program	REFMAC 5.3.0037	Depositor
R, R_{free}	0.205 , 0.231 0.216 , 0.245	Depositor DCC
R_{free} test set	2215 reflections (2.92%)	wwPDB-VP
Wilson B-factor (Å ²)	41.5	Xtriage
Anisotropy	0.118	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 54.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	7993	wwPDB-VP
Average B, all atoms (Å ²)	45.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.94% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: HEM, D2V, UNX, GLC

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	0.94	0/3821	0.84	0/5189
1	B	0.85	0/3844	0.80	1/5223 (0.0%)
All	All	0.90	0/7665	0.82	1/10412 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	237	ILE	CB-CA-C	-6.01	104.19	111.88

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	3712	0	3588	27	0
1	B	3732	0	3590	33	0
2	C	77	0	63	0	0
2	D	77	0	63	0	0
3	A	10	0	0	0	0
3	B	11	0	0	0	0
4	A	86	0	60	12	0
4	B	86	0	60	13	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	A	29	0	44	1	0
5	B	29	0	44	1	0
6	A	86	0	0	1	0
6	B	58	0	0	0	0
All	All	7993	0	7512	75	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 5.

All (75) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:55:LEU:HD11	1:B:66:MET:HE3	1.65	0.78
1:A:237:ILE:HD11	1:B:237:ILE:HD11	1.66	0.78
4:B:601[A]:HEM:HBB2	4:B:601[A]:HEM:HMB1	1.68	0.75
4:B:601[B]:HEM:HBC2	4:B:601[B]:HEM:HMC1	1.69	0.74
1:A:323:ILE:HG22	1:A:496:ILE:HD12	1.72	0.70
4:A:601[B]:HEM:HMC1	4:A:601[B]:HEM:HBC2	1.75	0.69
4:B:601[B]:HEM:HMB1	4:B:601[B]:HEM:HBB2	1.74	0.68
4:B:601[A]:HEM:HBC2	4:B:601[A]:HEM:HMC2	1.76	0.68
4:A:601[A]:HEM:HBC2	4:A:601[A]:HEM:HMC2	1.78	0.66
1:B:451:GLU:O	1:B:455:ARG:HG3	1.95	0.66
4:A:601[A]:HEM:HMB1	4:A:601[A]:HEM:HBB2	1.77	0.65
1:B:75:GLU:HB3	1:B:94:VAL:HG21	1.78	0.65
1:A:237:ILE:HD11	1:B:237:ILE:CD1	2.26	0.64
4:A:601[B]:HEM:HBB2	4:A:601[B]:HEM:HMB2	1.81	0.63
1:B:220:LEU:HD22	1:B:242:LYS:HB3	1.81	0.61
5:A:602:D2V:H1RA	5:A:602:D2V:H1E	1.80	0.61
1:A:48:PHE:CE1	1:A:232:PHE:CE2	2.88	0.61
1:A:323:ILE:CG2	1:A:496:ILE:HD12	2.30	0.60
1:A:237:ILE:CD1	1:B:237:ILE:HD11	2.31	0.58
1:B:140:ALA:HB3	1:B:449:LEU:HD11	1.84	0.58
1:B:214:PHE:HB3	1:B:309:ILE:HD13	1.85	0.57
5:B:602:D2V:H1E	5:B:602:D2V:H1RA	1.86	0.57
1:A:323:ILE:HG22	1:A:496:ILE:CD1	2.34	0.57
1:A:450:GLY:HA3	4:A:601[B]:HEM:C2B	2.40	0.57
4:B:601[A]:HEM:HHA	4:B:601[A]:HEM:HBA2	1.88	0.56
1:B:144:PHE:CE2	4:B:601[A]:HEM:HBC1	2.41	0.56
1:B:318:VAL:HG21	4:B:601[A]:HEM:CBB	2.38	0.54
4:B:601[A]:HEM:HBB2	4:B:601[A]:HEM:CMB	2.38	0.54
1:B:323:ILE:HG22	1:B:496:ILE:HD12	1.89	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:176:PRO:HB3	1:A:474:PRO:HG3	1.91	0.53
1:B:324:LEU:HD12	1:B:473:PHE:CZ	2.44	0.53
1:B:323:ILE:CG2	1:B:496:ILE:HD12	2.39	0.53
1:B:132:GLY:HA2	1:B:294:THR:HG21	1.91	0.51
4:B:601[B]:HEM:HBC2	4:B:601[B]:HEM:CMC	2.39	0.50
1:B:99:LEU:HD11	1:B:402:THR:HG21	1.92	0.50
1:A:47:PRO:O	1:A:48:PHE:HB3	2.11	0.50
4:B:601[B]:HEM:HBB2	4:B:601[B]:HEM:CMB	2.42	0.50
1:B:191:THR:CG2	1:B:195:ILE:HD12	2.43	0.48
1:B:105:ILE:HG21	1:B:387:ALA:HB2	1.95	0.48
1:B:318:VAL:HG21	4:B:601[A]:HEM:HBB1	1.96	0.48
1:A:48:PHE:CD2	1:A:49:ILE:HG23	2.48	0.48
1:B:75:GLU:CB	1:B:94:VAL:HG21	2.44	0.46
4:A:601[B]:HEM:HBC2	4:A:601[B]:HEM:CMC	2.44	0.46
4:B:601[A]:HEM:HBC2	4:B:601[A]:HEM:CMC	2.45	0.45
4:A:601[A]:HEM:HBB2	4:A:601[A]:HEM:CMB	2.47	0.45
1:A:311:GLY:HA2	4:A:601[A]:HEM:CMC	2.46	0.45
1:B:140:ALA:CB	1:B:449:LEU:HD11	2.47	0.44
1:B:323:ILE:HG22	1:B:496:ILE:CD1	2.47	0.44
1:B:46:LEU:CD1	1:B:49:ILE:HD11	2.48	0.44
1:B:55:LEU:HD11	1:B:66:MET:CE	2.43	0.43
1:A:280:TYR:CG	1:A:300:LEU:HD13	2.53	0.43
1:A:237:ILE:CG1	1:B:237:ILE:HG13	2.48	0.43
1:A:81:LEU:HD13	1:A:227:PHE:CE2	2.53	0.43
1:A:449:LEU:HD13	4:A:601[A]:HEM:CMD	2.49	0.43
4:A:601[A]:HEM:HBC2	4:A:601[A]:HEM:CMC	2.47	0.43
1:B:323:ILE:HG23	1:B:471:LEU:HD13	2.00	0.43
1:A:336:VAL:HG11	1:A:466:LEU:HD12	2.01	0.42
1:B:48:PHE:O	1:B:82:GLY:HA2	2.19	0.42
1:A:374:ILE:HD12	1:A:374:ILE:N	2.34	0.42
1:B:344:MET:HE3	1:B:358:MET:SD	2.60	0.42
4:B:601[B]:HEM:HHA	4:B:601[B]:HEM:HBD2	2.02	0.41
1:B:46:LEU:HD12	1:B:49:ILE:HD11	2.01	0.41
1:A:76:ILE:HD11	1:A:394:ILE:HD11	2.03	0.41
1:A:48:PHE:HD2	1:A:49:ILE:HG23	1.85	0.41
1:A:191:THR:HG23	6:A:642:HOH:O	2.20	0.41
1:A:207:PHE:CE2	1:A:211:ILE:HD12	2.56	0.41
1:B:326:MET:HG3	1:B:466:LEU:HD11	2.03	0.41
1:A:206:ASP:OD1	1:B:259:ARG:NH2	2.52	0.41
4:A:601[B]:HEM:HBB2	4:A:601[B]:HEM:CMB	2.50	0.41
1:A:449:LEU:HD13	4:A:601[A]:HEM:HMD2	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:323:ILE:CG2	1:A:496:ILE:CD1	2.95	0.40
1:B:324:LEU:HD12	1:B:473:PHE:HZ	1.86	0.40
1:A:89:LEU:HD22	1:A:94:VAL:HG12	2.03	0.40
1:A:328:LEU:HD22	1:A:481:LEU:HD11	2.04	0.40
1:B:324:LEU:HD12	1:B:473:PHE:CE1	2.57	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	463/481 (96%)	451 (97%)	12 (3%)	0	100	100
1	B	465/481 (97%)	454 (98%)	10 (2%)	1 (0%)	43	55
All	All	928/962 (96%)	905 (98%)	22 (2%)	1 (0%)	48	60

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	501	ARG

5.3.2 Protein sidechains [i](#)

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

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

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	393/424 (93%)	391 (100%)	2 (0%)	81	90
1	B	392/424 (92%)	391 (100%)	1 (0%)	86	93
All	All	785/848 (93%)	782 (100%)	3 (0%)	84	92

All (3) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	449	LEU
1	A	496	ILE
1	B	496	ILE

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

Mol	Chain	Res	Type
1	A	69	GLN
1	A	181	GLN
1	A	492	GLN
1	B	503	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

14 monosaccharides are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the

expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	GLC	C	1	2	11,11,12	0.73	0	15,15,17	1.07	1 (6%)
2	GLC	C	2	2	11,11,12	0.46	0	15,15,17	0.74	0
2	GLC	C	3	2	11,11,12	0.55	0	15,15,17	1.32	2 (13%)
2	GLC	C	4	2	11,11,12	0.38	0	15,15,17	1.48	3 (20%)
2	GLC	C	5	2	11,11,12	0.65	0	15,15,17	1.71	2 (13%)
2	GLC	C	6	2	11,11,12	0.57	0	15,15,17	0.78	1 (6%)
2	GLC	C	7	2	11,11,12	0.43	0	15,15,17	1.45	3 (20%)
2	GLC	D	1	2	11,11,12	0.90	0	15,15,17	1.61	4 (26%)
2	GLC	D	2	2	11,11,12	0.63	0	15,15,17	1.16	2 (13%)
2	GLC	D	3	2	11,11,12	0.65	0	15,15,17	0.99	1 (6%)
2	GLC	D	4	2	11,11,12	0.47	0	15,15,17	1.53	4 (26%)
2	GLC	D	5	2	11,11,12	0.63	0	15,15,17	1.09	1 (6%)
2	GLC	D	6	2	11,11,12	0.72	0	15,15,17	1.19	1 (6%)
2	GLC	D	7	2	11,11,12	0.37	0	15,15,17	1.10	2 (13%)

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	GLC	C	1	2	-	2/2/19/22	0/1/1/1
2	GLC	C	2	2	-	2/2/19/22	0/1/1/1
2	GLC	C	3	2	-	1/2/19/22	0/1/1/1
2	GLC	C	4	2	-	2/2/19/22	0/1/1/1
2	GLC	C	5	2	-	2/2/19/22	0/1/1/1
2	GLC	C	6	2	-	2/2/19/22	0/1/1/1
2	GLC	C	7	2	-	1/2/19/22	0/1/1/1
2	GLC	D	1	2	-	2/2/19/22	0/1/1/1
2	GLC	D	2	2	-	2/2/19/22	0/1/1/1
2	GLC	D	3	2	-	2/2/19/22	0/1/1/1
2	GLC	D	4	2	-	0/2/19/22	0/1/1/1
2	GLC	D	5	2	-	2/2/19/22	0/1/1/1
2	GLC	D	6	2	-	2/2/19/22	0/1/1/1
2	GLC	D	7	2	-	0/2/19/22	0/1/1/1

There are no bond length outliers.

All (27) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	5	GLC	C1-C2-C3	4.58	116.31	109.64
2	D	4	GLC	C1-C2-C3	3.52	114.77	109.64
2	D	1	GLC	C3-C4-C5	3.48	116.54	110.23
2	C	5	GLC	C2-C3-C4	3.45	116.92	110.86
2	C	7	GLC	C3-C4-C5	3.21	116.05	110.23
2	D	2	GLC	C1-C2-C3	3.14	114.21	109.64
2	C	4	GLC	C1-C2-C3	3.01	114.03	109.64
2	C	3	GLC	C1-C2-C3	3.01	114.03	109.64
2	C	4	GLC	C1-O5-C5	2.90	116.07	112.19
2	D	6	GLC	C1-C2-C3	2.83	113.76	109.64
2	D	7	GLC	C1-O5-C5	2.79	115.92	112.19
2	D	1	GLC	O4-C4-C3	-2.72	103.97	110.38
2	D	7	GLC	O4-C4-C3	-2.69	104.04	110.38
2	D	5	GLC	C1-O5-C5	2.66	115.76	112.19
2	C	3	GLC	C2-C3-C4	2.49	115.24	110.86
2	D	1	GLC	O5-C1-C2	2.42	116.56	110.79
2	D	4	GLC	C2-C3-C4	2.35	114.99	110.86
2	C	6	GLC	C1-O5-C5	2.30	115.27	112.19
2	C	4	GLC	C2-C3-C4	2.28	114.88	110.86
2	D	4	GLC	O4-C4-C3	-2.24	105.11	110.38
2	D	1	GLC	C1-O5-C5	-2.22	109.20	112.19
2	C	7	GLC	C1-O5-C5	2.22	115.17	112.19
2	D	4	GLC	O5-C5-C4	-2.18	105.53	110.83
2	D	3	GLC	C1-O5-C5	2.14	115.06	112.19
2	C	1	GLC	O4-C4-C3	-2.05	105.55	110.38
2	C	7	GLC	O4-C4-C3	-2.04	105.56	110.38
2	D	2	GLC	C2-C3-C4	2.04	114.45	110.86

There are no chirality outliers.

All (22) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	C	4	GLC	O5-C5-C6-O6
2	D	3	GLC	O5-C5-C6-O6
2	D	1	GLC	C4-C5-C6-O6
2	C	4	GLC	C4-C5-C6-O6
2	D	2	GLC	O5-C5-C6-O6
2	D	2	GLC	C4-C5-C6-O6
2	D	1	GLC	O5-C5-C6-O6
2	D	5	GLC	C4-C5-C6-O6

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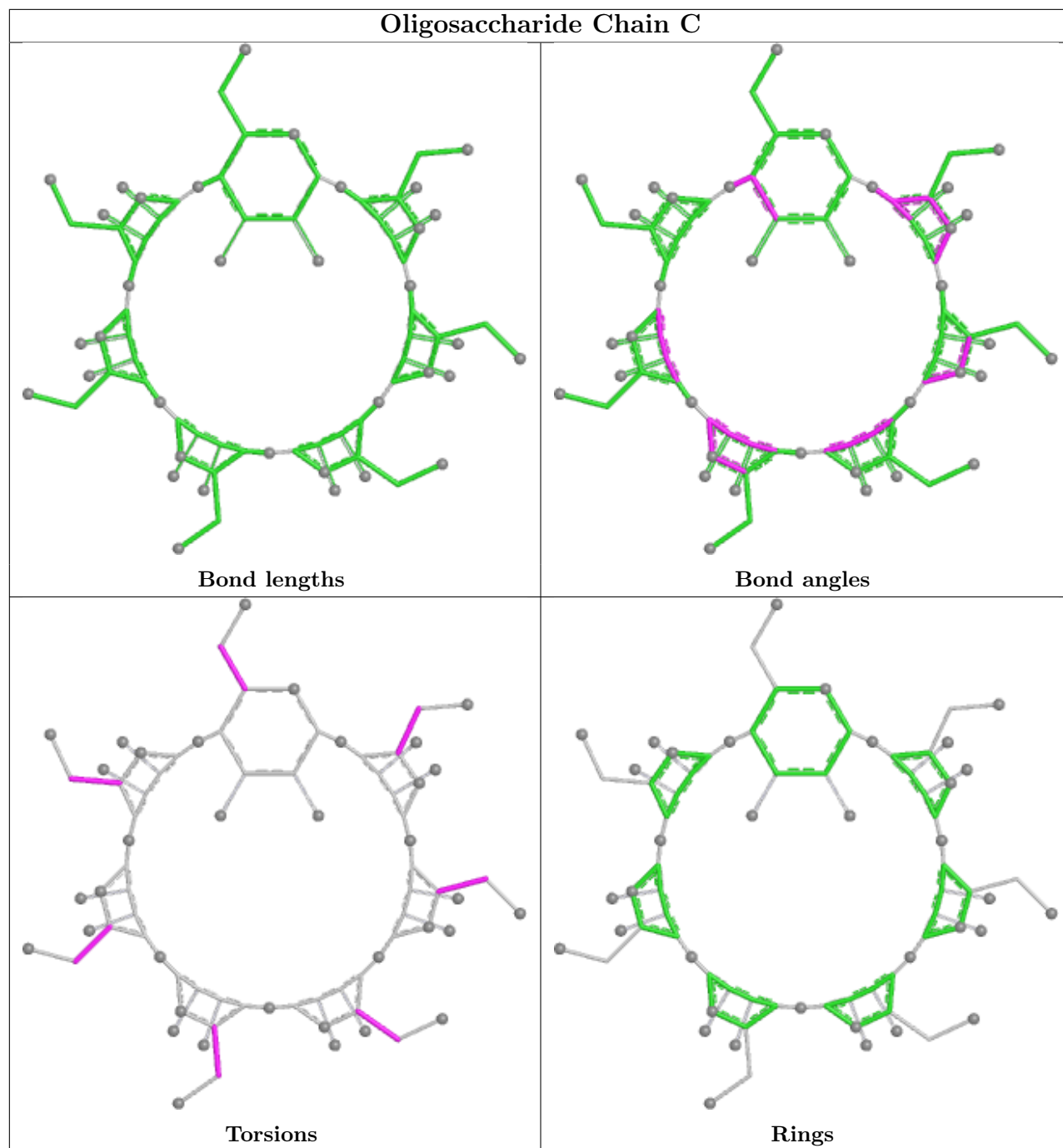
Mol	Chain	Res	Type	Atoms
2	D	3	GLC	C4-C5-C6-O6
2	D	5	GLC	O5-C5-C6-O6
2	C	1	GLC	O5-C5-C6-O6
2	C	5	GLC	C4-C5-C6-O6
2	C	2	GLC	O5-C5-C6-O6
2	C	6	GLC	O5-C5-C6-O6
2	C	1	GLC	C4-C5-C6-O6
2	C	5	GLC	O5-C5-C6-O6
2	D	6	GLC	O5-C5-C6-O6
2	C	3	GLC	O5-C5-C6-O6
2	C	2	GLC	C4-C5-C6-O6
2	C	7	GLC	O5-C5-C6-O6
2	C	6	GLC	C4-C5-C6-O6
2	D	6	GLC	C4-C5-C6-O6

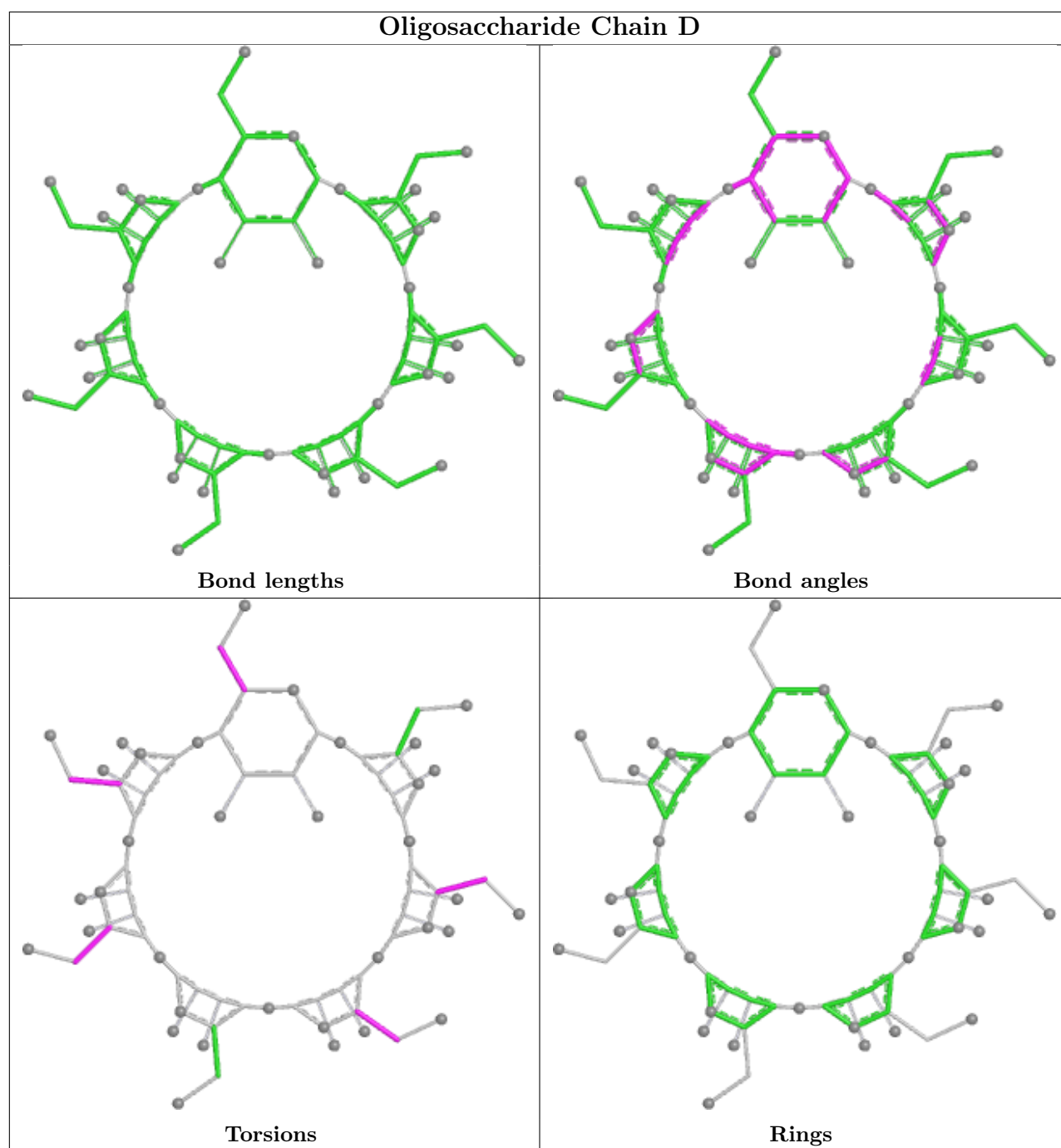
There are no ring outliers.

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.

Oligosaccharide Chain C





5.6 Ligand geometry [i](#)

Of 27 ligands modelled in this entry, 21 are unknown - leaving 6 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
4	HEM	A	601[A]	1	50,50,50	2.06	9 (18%)	67,82,82	1.47	10 (14%)
4	HEM	B	601[A]	1	50,50,50	1.99	9 (18%)	67,82,82	1.34	4 (5%)
5	D2V	A	602	-	31,31,31	1.86	8 (25%)	45,45,45	1.92	10 (22%)
4	HEM	A	601[B]	1	50,50,50	2.15	11 (22%)	67,82,82	1.41	8 (11%)
5	D2V	B	602	-	31,31,31	1.79	8 (25%)	45,45,45	1.84	11 (24%)
4	HEM	B	601[B]	1	50,50,50	1.92	10 (20%)	67,82,82	1.29	5 (7%)

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
4	HEM	A	601[A]	1	-	5/14/54/54	-
4	HEM	B	601[A]	1	-	6/14/54/54	-
5	D2V	A	602	-	-	3/18/59/59	0/3/3/3
4	HEM	A	601[B]	1	-	6/14/54/54	-
5	D2V	B	602	-	-	4/18/59/59	0/3/3/3
4	HEM	B	601[B]	1	-	4/14/54/54	-

All (55) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	601[A]	HEM	C3D-C2D	7.80	1.53	1.36
4	A	601[A]	HEM	C3D-C2D	7.56	1.53	1.36
4	B	601[B]	HEM	C3D-C2D	7.35	1.52	1.36
4	A	601[B]	HEM	C3D-C2D	7.26	1.52	1.36
4	A	601[B]	HEM	FE-ND	6.88	2.16	1.94
4	A	601[A]	HEM	FE-ND	6.08	2.13	1.94
4	B	601[A]	HEM	FE-NB	6.07	2.13	1.94
4	B	601[B]	HEM	FE-ND	6.03	2.13	1.94
4	B	601[A]	HEM	FE-NA	5.57	2.13	1.95
4	A	601[A]	HEM	FE-NA	5.44	2.13	1.95
4	A	601[B]	HEM	FE-NC	4.88	2.11	1.95
4	A	601[B]	HEM	FE-NA	4.85	2.11	1.95
5	A	602	D2V	C1T-C1U	-4.83	1.40	1.48
4	A	601[A]	HEM	FE-NB	4.73	2.09	1.94
4	A	601[B]	HEM	FE-NB	4.58	2.09	1.94

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	602	D2V	C1I-C1V	4.47	1.41	1.34
5	B	602	D2V	C1T-C1U	-4.39	1.40	1.48
4	B	601[B]	HEM	FE-NC	4.01	2.08	1.95
5	B	602	D2V	C1I-C1V	3.97	1.40	1.34
4	B	601[B]	HEM	FE-NB	3.56	2.05	1.94
4	A	601[A]	HEM	CAB-C3B	3.39	1.56	1.47
4	A	601[B]	HEM	CAB-C3B	3.34	1.56	1.47
4	B	601[B]	HEM	CAC-C3C	3.28	1.56	1.47
4	B	601[A]	HEM	CAB-C3B	3.25	1.56	1.47
4	A	601[A]	HEM	CAC-C3C	3.17	1.55	1.47
5	A	602	D2V	C1X-C1J	3.16	1.58	1.51
4	B	601[A]	HEM	CAC-C3C	3.06	1.55	1.47
5	A	602	D2V	C1A-C1T	3.06	1.38	1.32
5	B	602	D2V	C2C-C2A	-3.03	1.51	1.56
4	A	601[B]	HEM	CAC-C3C	3.03	1.55	1.47
4	B	601[B]	HEM	CAB-C3B	2.91	1.55	1.47
5	B	602	D2V	C1A-C1T	2.86	1.38	1.32
4	A	601[A]	HEM	CMD-C2D	2.80	1.56	1.50
5	B	602	D2V	C1O-C1Z	2.78	1.58	1.51
4	A	601[B]	HEM	CMA-C3A	2.70	1.56	1.50
5	A	602	D2V	C1O-C1Z	2.54	1.57	1.51
4	A	601[B]	HEM	CMB-C2B	2.50	1.55	1.50
4	A	601[B]	HEM	CMC-C2C	2.44	1.55	1.50
4	B	601[B]	HEM	CMC-C2C	2.37	1.55	1.50
5	B	602	D2V	C1S-C1U	2.36	1.55	1.51
4	B	601[B]	HEM	FE-NA	2.31	2.02	1.95
4	B	601[A]	HEM	FE-NC	2.29	2.02	1.95
5	B	602	D2V	C1K-C1J	2.29	1.39	1.32
4	A	601[A]	HEM	CMA-C3A	2.28	1.55	1.50
5	A	602	D2V	C2A-C1V	2.27	1.55	1.51
4	B	601[A]	HEM	CMB-C2B	2.20	1.55	1.50
4	A	601[B]	HEM	CMD-C2D	2.18	1.55	1.50
5	A	602	D2V	C1K-C1J	2.16	1.38	1.32
4	A	601[A]	HEM	CMC-C2C	2.15	1.55	1.50
4	B	601[B]	HEM	CMA-C3A	2.11	1.55	1.50
5	A	602	D2V	C1H-C1U	2.09	1.42	1.36
4	B	601[A]	HEM	C2A-C3A	-2.08	1.33	1.38
5	B	602	D2V	C1X-C1J	2.07	1.55	1.51
4	B	601[A]	HEM	CMD-C2D	2.07	1.55	1.50
4	B	601[B]	HEM	CMB-C2B	2.04	1.54	1.50

All (48) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	602	D2V	C2C-C2B-C1Y	-6.10	112.07	119.38
4	A	601[B]	HEM	C4D-ND-C1D	5.49	111.71	105.21
4	A	601[A]	HEM	C4D-ND-C1D	5.40	111.60	105.21
4	B	601[B]	HEM	C4D-ND-C1D	5.29	111.47	105.21
5	B	602	D2V	C1P-C2A-C1V	-5.26	114.59	120.68
4	B	601[A]	HEM	C4D-ND-C1D	5.19	111.35	105.21
5	B	602	D2V	C2C-C2B-C1Y	-4.47	114.03	119.38
5	A	602	D2V	C1P-C2A-C2C	-4.37	100.30	104.21
5	A	602	D2V	C1P-C2A-C1V	-3.80	116.28	120.68
4	A	601[A]	HEM	C1B-NB-C4B	3.44	109.28	105.21
5	A	602	D2V	C1L-C1R-C2C	-3.42	109.18	113.19
5	B	602	D2V	C1M-C1V-C1I	-3.33	121.67	125.27
5	A	602	D2V	C1M-C1V-C1I	-3.12	121.90	125.27
5	B	602	D2V	C1Z-C1S-C1U	-2.93	108.23	112.51
4	A	601[A]	HEM	CAA-CBA-CGA	-2.84	106.12	113.67
4	A	601[B]	HEM	C1B-NB-C4B	2.79	108.52	105.21
5	B	602	D2V	C1R-C1L-C1M	2.79	117.40	111.28
5	A	602	D2V	C1Q-C1P-C2A	-2.77	100.73	105.25
5	A	602	D2V	C1N-C1T-C1A	-2.72	118.09	122.42
5	B	602	D2V	C1R-C2C-C2B	-2.68	112.65	116.60
4	B	601[B]	HEM	C1B-NB-C4B	2.68	108.38	105.21
4	B	601[B]	HEM	CAD-C3D-C4D	2.67	129.34	124.70
4	A	601[B]	HEM	CAD-CBD-CGD	-2.63	106.69	113.67
5	B	602	D2V	C1Q-C1P-C2A	-2.61	101.00	105.25
5	B	602	D2V	C1S-C1U-C1T	2.61	116.83	113.86
5	A	602	D2V	C1S-C1U-C1T	2.54	116.75	113.86
5	B	602	D2V	C1N-C1T-C1U	2.49	118.07	114.51
4	A	601[B]	HEM	C2A-C1A-NA	-2.48	107.41	110.15
4	A	601[A]	HEM	C3B-C4B-NB	-2.47	107.69	109.47
5	B	602	D2V	C1H-C1I-C1V	-2.45	122.52	126.13
4	A	601[B]	HEM	CAD-C3D-C4D	2.44	128.95	124.70
4	B	601[B]	HEM	C3B-C4B-NB	-2.43	107.72	109.47
5	A	602	D2V	C1X-C1J-C1K	-2.41	118.51	125.81
5	A	602	D2V	C1A-C1T-C1U	2.32	125.64	123.05
4	A	601[A]	HEM	C2A-C1A-NA	-2.26	107.65	110.15
4	A	601[B]	HEM	CHD-C1D-ND	2.25	126.84	124.42
5	B	602	D2V	C1X-C1J-C1K	-2.15	119.31	125.81
4	A	601[A]	HEM	C3B-C2B-C1B	2.15	108.02	106.41
4	B	601[A]	HEM	C2A-C1A-NA	-2.12	107.80	110.15
4	A	601[A]	HEM	CHD-C4C-NC	2.08	126.72	124.45
4	B	601[A]	HEM	CHA-C4D-ND	2.07	126.93	124.37
4	B	601[A]	HEM	CAA-CBA-CGA	-2.05	108.24	113.67
4	A	601[B]	HEM	C3B-C4B-NB	-2.04	108.01	109.47

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	601[B]	HEM	CMD-C2D-C1D	2.03	128.21	125.03
4	A	601[A]	HEM	C2B-C1B-NB	-2.02	107.51	109.84
4	B	601[B]	HEM	CAD-CBD-CGD	-2.01	108.33	113.67
4	A	601[A]	HEM	CHD-C1D-ND	2.01	126.58	124.42
4	A	601[A]	HEM	O2A-CGA-CBA	2.00	120.33	114.00

There are no chirality outliers.

All (28) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	B	601[A]	HEM	C1A-C2A-CAA-CBA
5	A	602	D2V	C1J-C1K-C1Y-C2B
4	B	601[A]	HEM	C3A-C2A-CAA-CBA
5	B	602	D2V	C1J-C1K-C1Y-C2B
4	A	601[A]	HEM	C1A-C2A-CAA-CBA
4	B	601[B]	HEM	C2D-C3D-CAD-CBD
4	B	601[B]	HEM	C4D-C3D-CAD-CBD
5	B	602	D2V	C1J-C1K-C1Y-C1E
4	A	601[A]	HEM	C3A-C2A-CAA-CBA
5	A	602	D2V	C1J-C1K-C1Y-C1E
4	B	601[B]	HEM	CAD-CBD-CGD-O2D
5	B	602	D2V	C1K-C1J-C1X-C1W
4	B	601[B]	HEM	CAD-CBD-CGD-O1D
4	A	601[B]	HEM	C2D-C3D-CAD-CBD
4	A	601[B]	HEM	CAD-CBD-CGD-O2D
4	B	601[A]	HEM	CAA-CBA-CGA-O2A
4	A	601[A]	HEM	CAA-CBA-CGA-O2A
5	B	602	D2V	C1K-C1J-C1X-C1D
4	A	601[A]	HEM	CAA-CBA-CGA-O1A
4	A	601[B]	HEM	CAD-CBD-CGD-O1D
4	B	601[A]	HEM	CAA-CBA-CGA-O1A
4	A	601[B]	HEM	C4D-C3D-CAD-CBD
5	A	602	D2V	C1K-C1J-C1X-C1D
4	A	601[A]	HEM	CAD-CBD-CGD-O2D
4	B	601[A]	HEM	CAD-CBD-CGD-O1D
4	A	601[B]	HEM	C2B-C3B-CAB-CBB
4	A	601[B]	HEM	C2C-C3C-CAC-CBC
4	B	601[A]	HEM	C2B-C3B-CAB-CBB

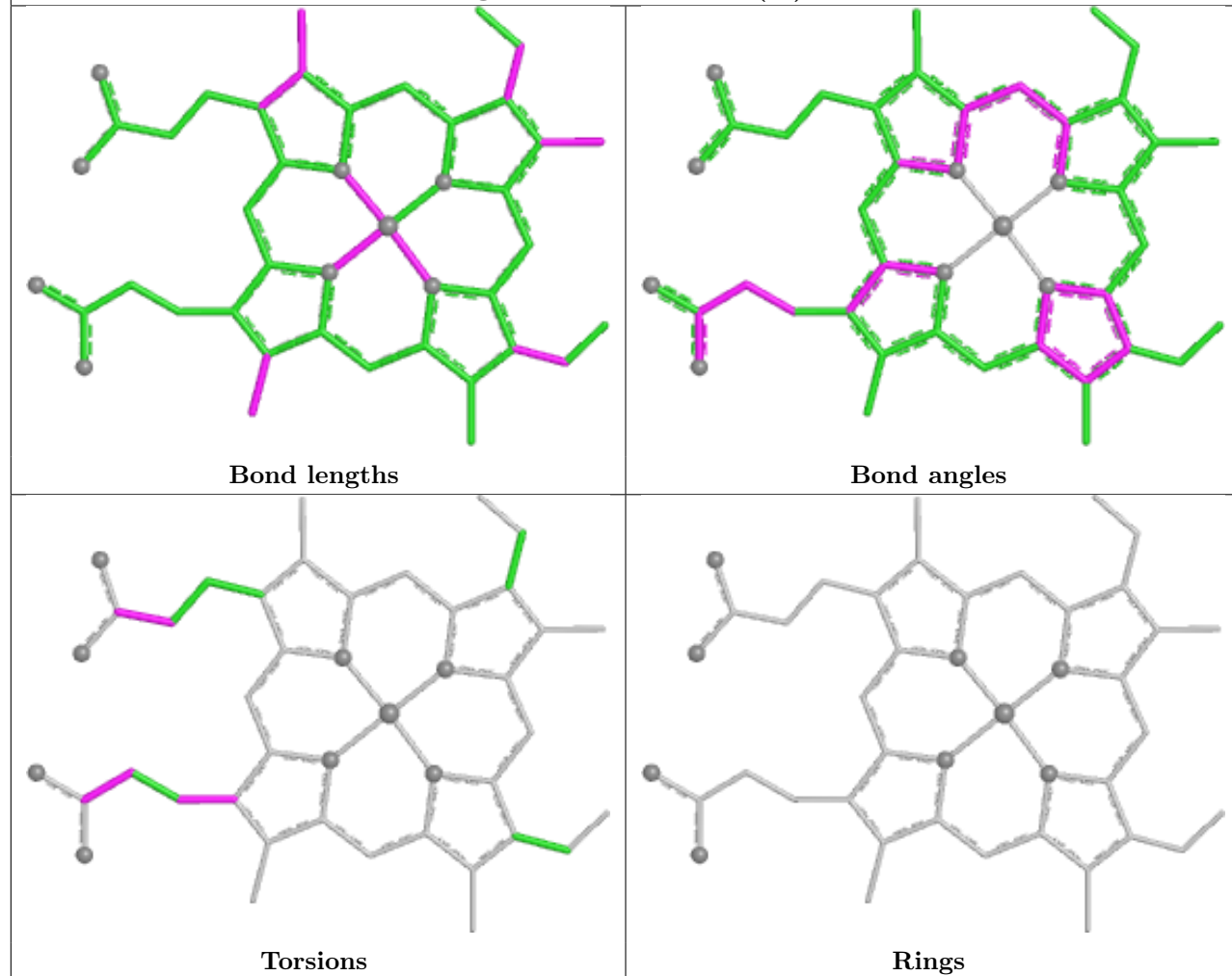
There are no ring outliers.

6 monomers are involved in 27 short contacts:

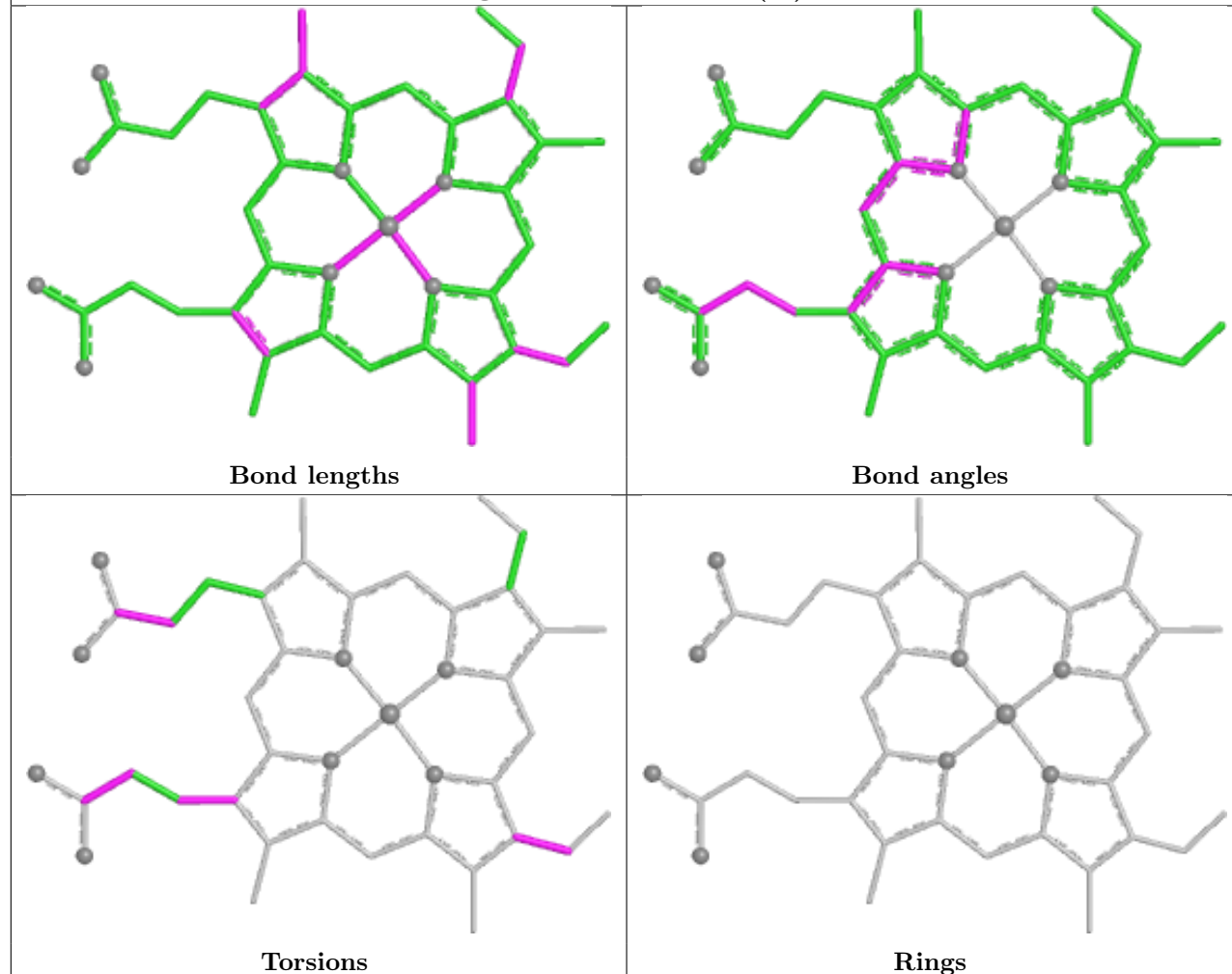
Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	601[A]	HEM	7	0
4	B	601[A]	HEM	8	0
5	A	602	D2V	1	0
4	A	601[B]	HEM	5	0
5	B	602	D2V	1	0
4	B	601[B]	HEM	5	0

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.

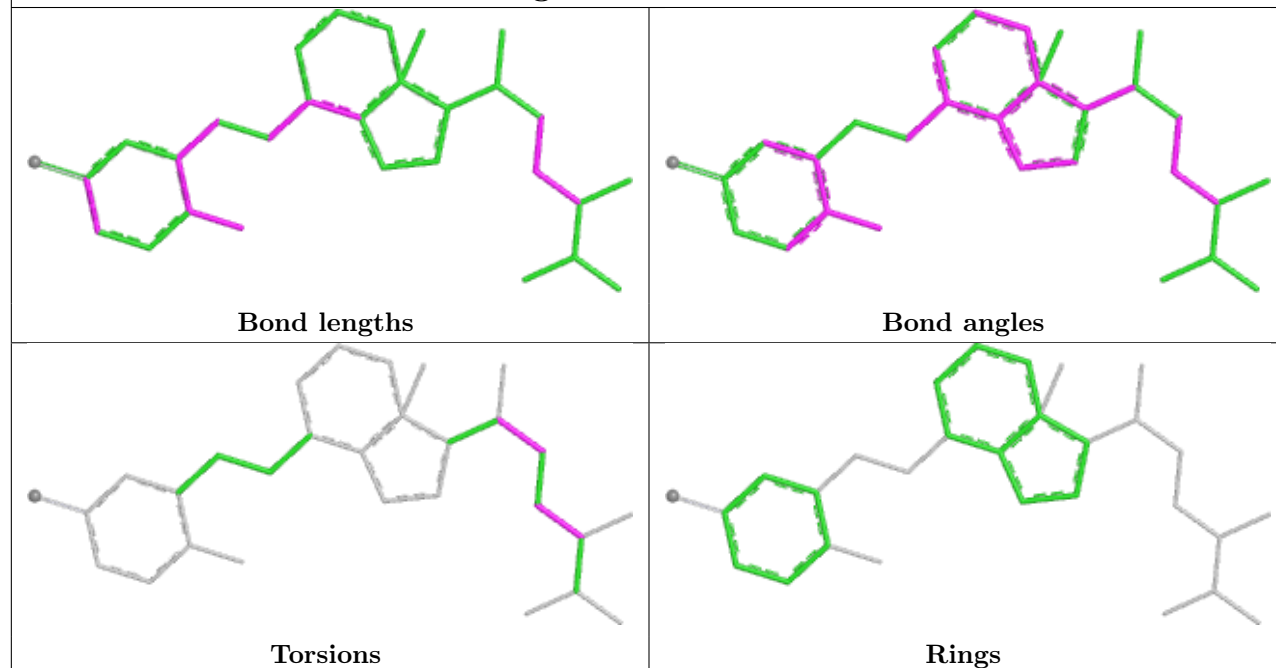
Ligand HEM A 601 (A)



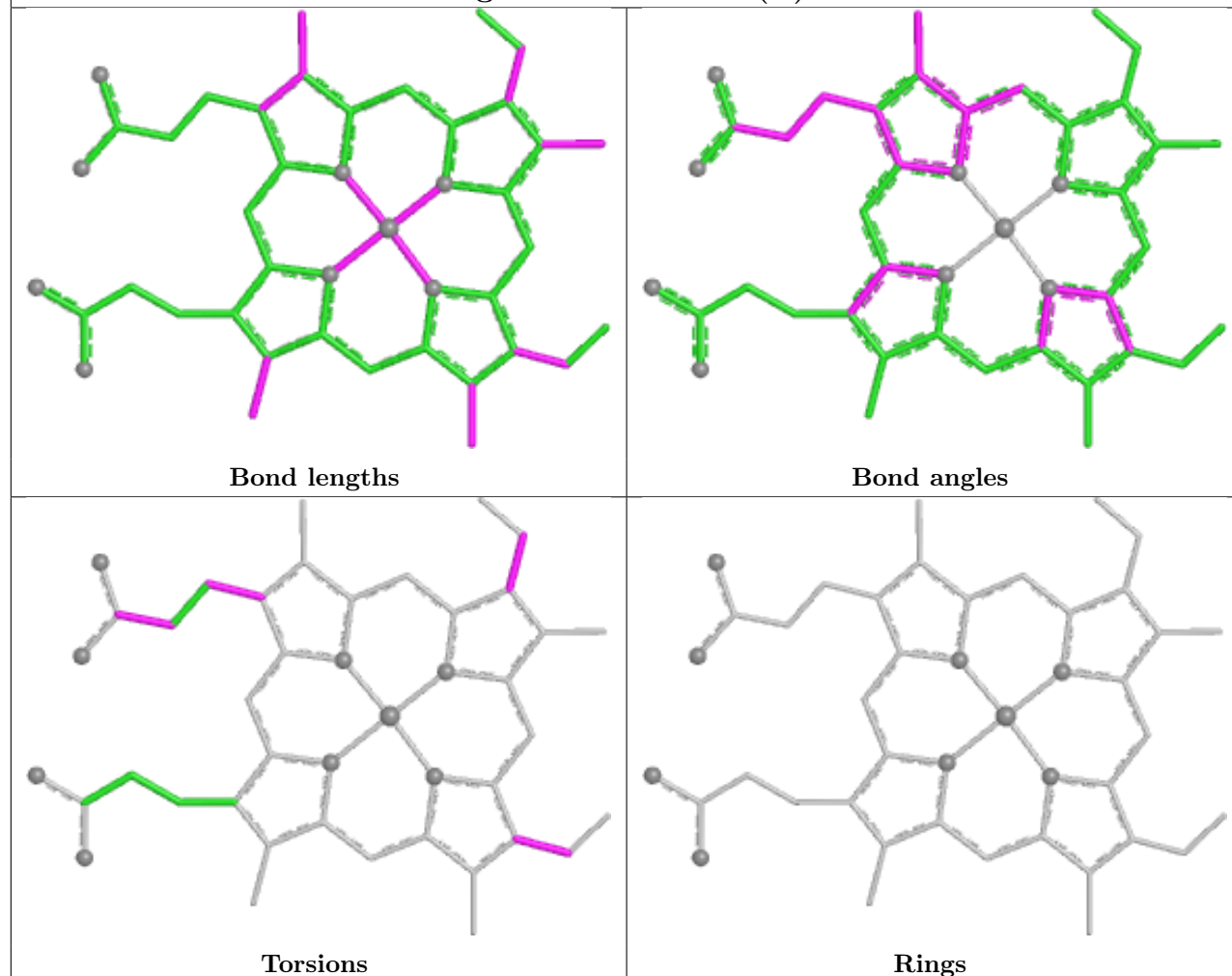
Ligand HEM B 601 (A)



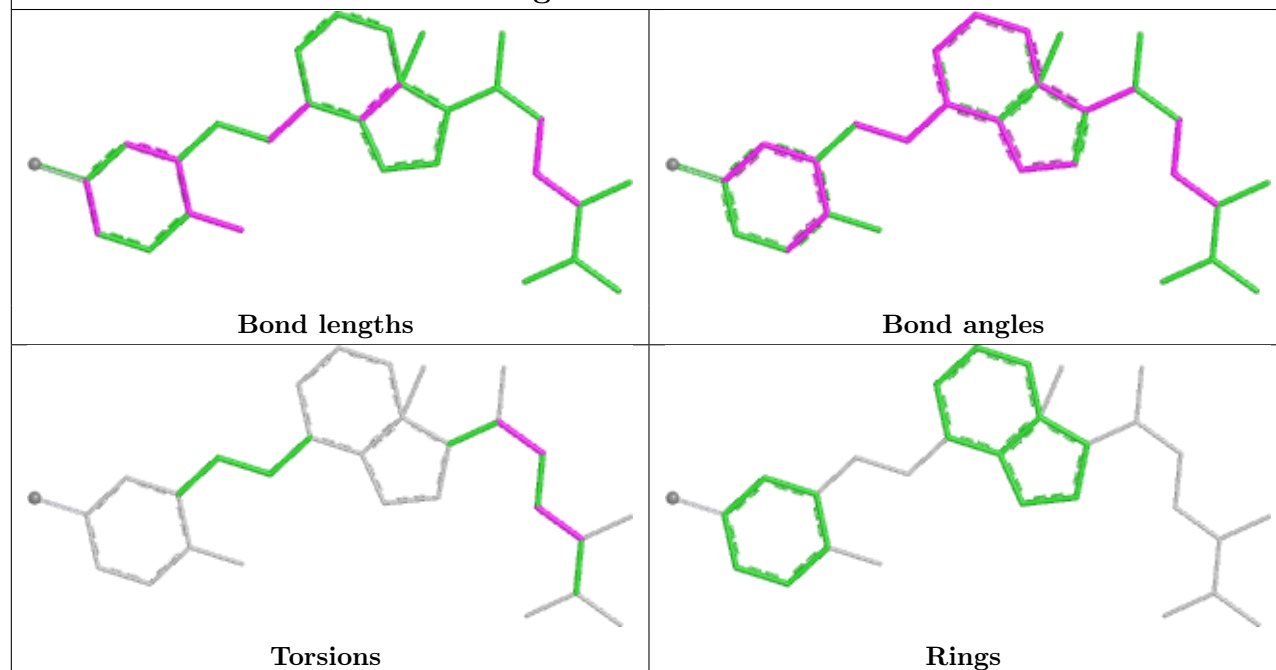
Ligand D2V A 602

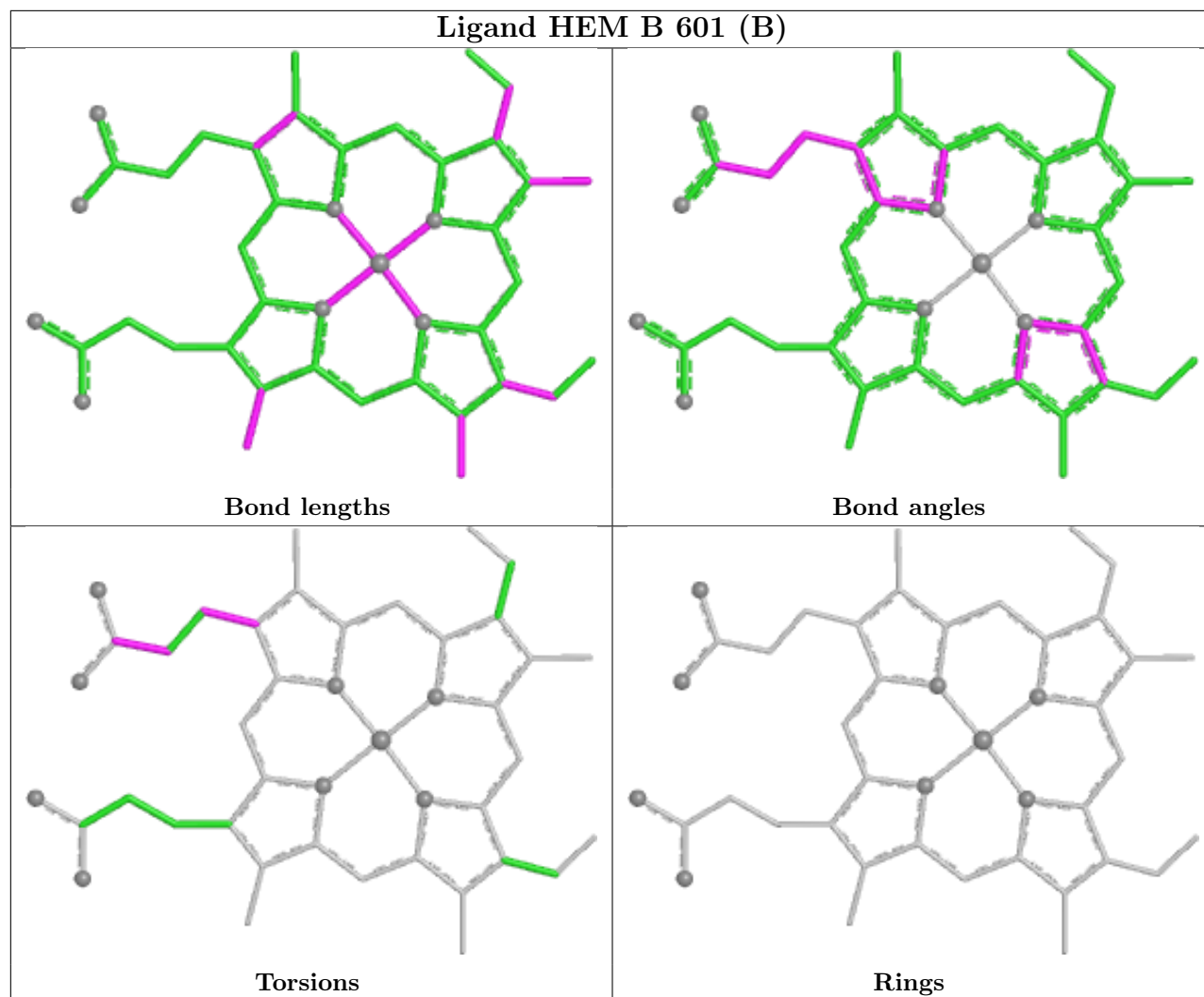


Ligand HEM A 601 (B)



Ligand D2V B 602





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	465/481 (96%)	0.35	24 (5%) 33 34	25, 40, 59, 80	0
1	B	467/481 (97%)	0.61	37 (7%) 18 20	30, 46, 68, 86	0
All	All	932/962 (96%)	0.48	61 (6%) 25 27	25, 43, 66, 86	0

All (61) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	477	LEU	5.9
1	A	502	HIS	5.5
1	B	429	SER	4.8
1	A	291	PRO	4.6
1	A	48	PHE	4.6
1	A	476	GLU	4.3
1	A	423	GLU	4.2
1	B	431	TYR	4.2
1	B	291	PRO	3.9
1	B	428	SER	3.9
1	A	478	VAL	3.8
1	A	475	HIS	3.8
1	B	503	HIS	3.7
1	A	346	PRO	3.6
1	B	101	HIS	3.4
1	B	75	GLU	3.3
1	A	271	GLN	3.3
1	A	286	GLN	3.3
1	B	391	GLY	3.2
1	B	266	VAL	3.2
1	B	505	HIS	3.2
1	A	46	LEU	3.1
1	B	504	HIS	3.0
1	A	212	GLU	3.0

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Mol	Chain	Res	Type	RSRZ
1	B	289	ASN	2.9
1	B	39	PHE	2.9
1	B	385	GLU	2.8
1	B	293	SER	2.8
1	A	116	MET	2.6
1	B	272	LEU	2.6
1	B	203	GLU	2.5
1	A	347	ASN	2.5
1	B	237	ILE	2.5
1	A	240	PHE	2.5
1	B	58	SER	2.5
1	B	175	ARG	2.5
1	B	386	ASP	2.5
1	A	266	VAL	2.4
1	B	313	GLU	2.4
1	B	430	GLY	2.4
1	A	237	ILE	2.4
1	B	240	PHE	2.4
1	B	287	GLY	2.3
1	B	57	ALA	2.3
1	A	288	LYS	2.3
1	B	415	ARG	2.2
1	A	38	GLY	2.2
1	B	416	ASP	2.2
1	A	353	ASP	2.2
1	B	239	PRO	2.2
1	A	259	ARG	2.1
1	A	45	GLY	2.1
1	B	477	LEU	2.1
1	B	212	GLU	2.1
1	B	288	LYS	2.1
1	B	238	LEU	2.1
1	B	271	GLN	2.1
1	A	232	PHE	2.1
1	B	347	ASN	2.0
1	B	259	ARG	2.0
1	B	436	GLU	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

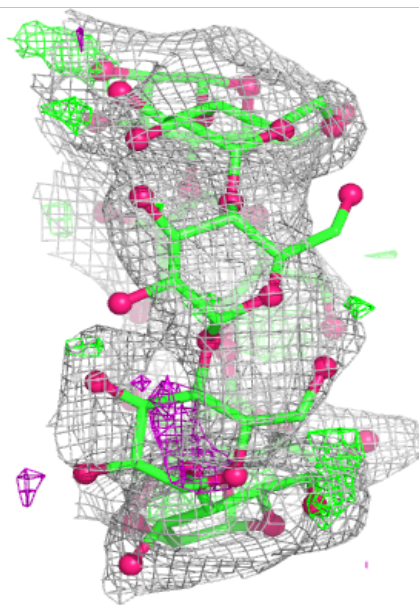
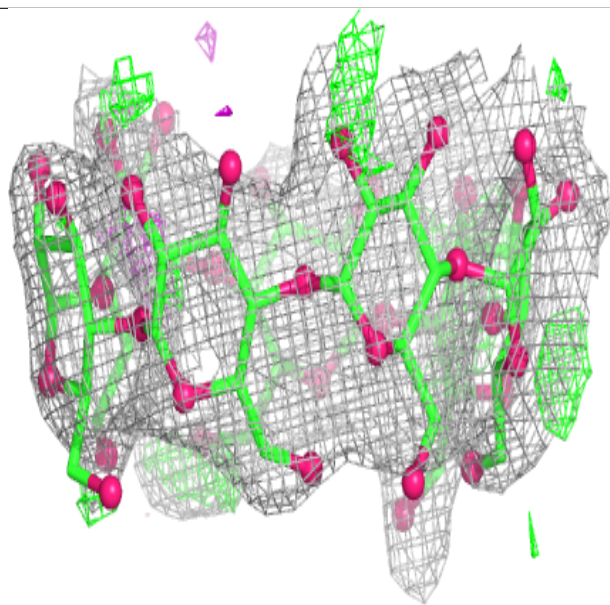
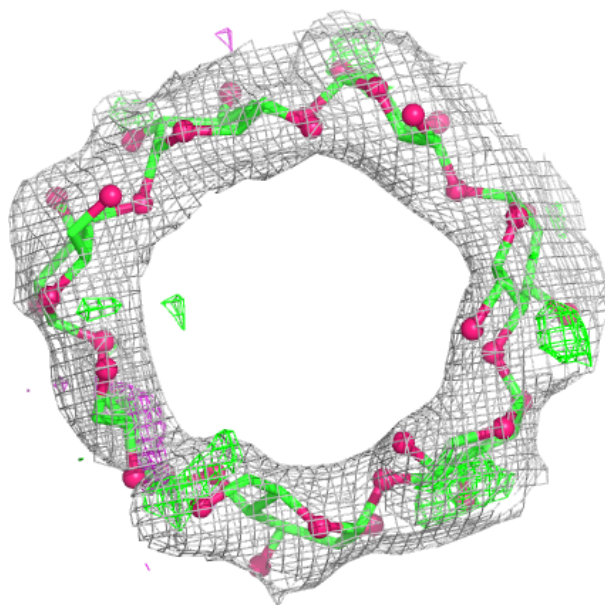
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

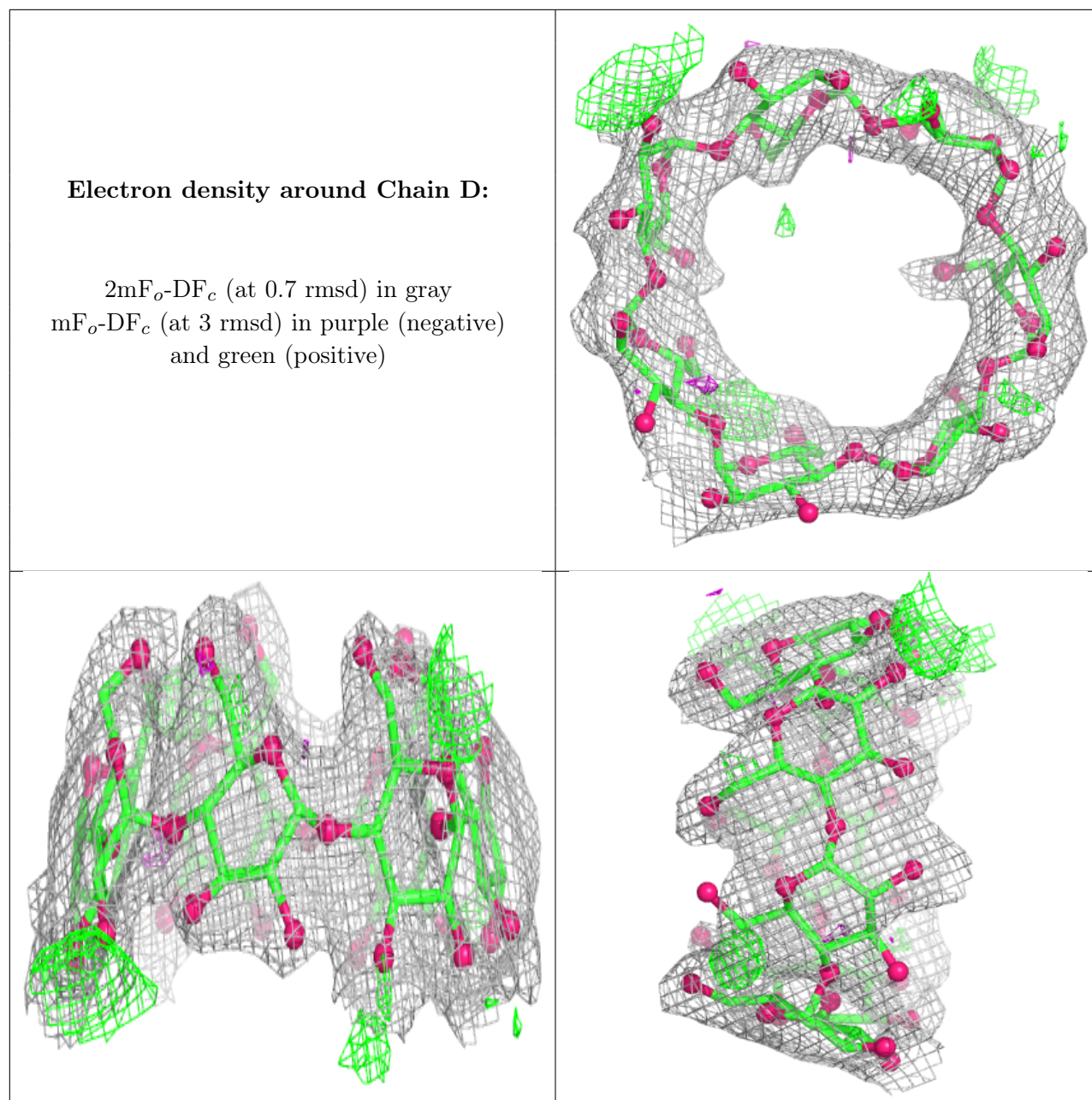
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	GLC	C	6	11/12	0.62	0.31	101,102,102,103	0
2	GLC	C	1	11/12	0.74	0.18	98,98,98,99	0
2	GLC	D	6	11/12	0.74	0.19	75,78,79,80	0
2	GLC	C	3	11/12	0.75	0.20	99,99,99,99	0
2	GLC	D	5	11/12	0.76	0.16	71,72,72,74	0
2	GLC	C	5	11/12	0.76	0.19	99,100,100,101	0
2	GLC	C	7	11/12	0.78	0.22	99,101,101,102	0
2	GLC	D	3	11/12	0.78	0.15	68,70,72,72	0
2	GLC	D	2	11/12	0.79	0.14	74,75,76,76	0
2	GLC	D	7	11/12	0.80	0.17	79,80,80,81	0
2	GLC	D	1	11/12	0.82	0.13	77,79,79,79	0
2	GLC	C	2	11/12	0.82	0.13	97,99,99,99	0
2	GLC	D	4	11/12	0.84	0.15	66,68,69,69	0
2	GLC	C	4	11/12	0.87	0.21	97,98,99,99	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)





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
3	UNX	B	2	1/1	-0.57	1.92	2,2,2,2	1
3	UNX	B	10	1/1	-0.47	3.18	2,2,2,2	1
3	UNX	A	4	1/1	-0.36	1.67	2,2,2,2	1

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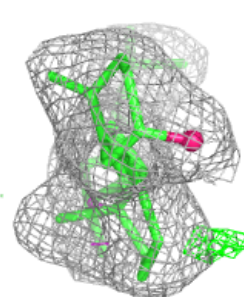
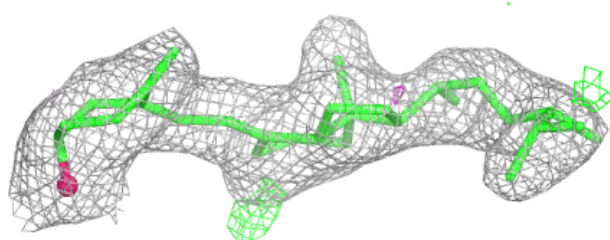
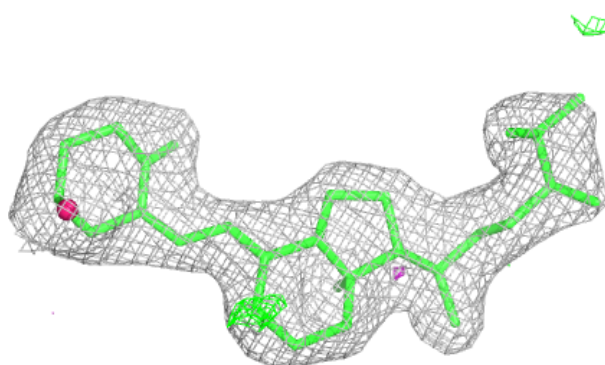
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	UNX	B	11	1/1	-0.22	1.56	2,2,2,2	1
3	UNX	A	1	1/1	-0.19	1.29	2,2,2,2	1
3	UNX	A	19	1/1	-0.16	1.60	2,2,2,2	1
3	UNX	A	6	1/1	-0.03	2.98	2,2,2,2	1
3	UNX	B	13	1/1	0.06	2.81	2,2,2,2	1
3	UNX	A	7	1/1	0.08	1.54	2,2,2,2	1
3	UNX	A	3	1/1	0.25	1.68	2,2,2,2	1
3	UNX	B	12	1/1	0.33	1.39	2,2,2,2	1
3	UNX	B	18	1/1	0.38	2.65	2,2,2,2	1
3	UNX	B	20	1/1	0.44	2.38	2,2,2,2	1
3	UNX	A	17	1/1	0.50	3.55	2,2,2,2	1
3	UNX	A	5	1/1	0.53	2.48	2,2,2,2	1
3	UNX	B	9	1/1	0.53	2.61	2,2,2,2	1
3	UNX	B	14	1/1	0.54	2.49	2,2,2,2	1
3	UNX	B	21	1/1	0.58	2.25	2,2,2,2	1
3	UNX	A	16	1/1	0.63	2.58	2,2,2,2	1
3	UNX	A	8	1/1	0.71	2.09	2,2,2,2	1
3	UNX	B	15	1/1	0.77	2.21	2,2,2,2	1
5	D2V	A	602	29/29	0.92	0.10	34,37,43,45	0
5	D2V	B	602	29/29	0.95	0.10	35,37,46,47	0
4	HEM	B	601[A]	43/43	0.98	0.08	32,34,36,36	43
4	HEM	B	601[B]	43/43	0.98	0.08	32,34,37,38	43
4	HEM	A	601[A]	43/43	0.98	0.06	25,27,28,29	43
4	HEM	A	601[B]	43/43	0.98	0.06	29,31,33,35	43

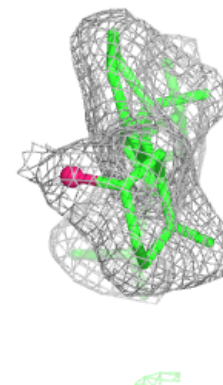
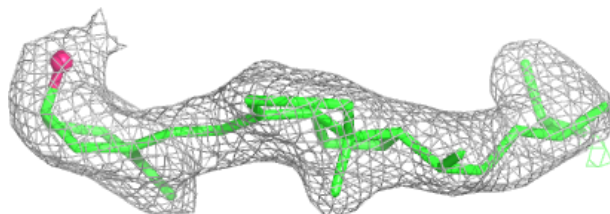
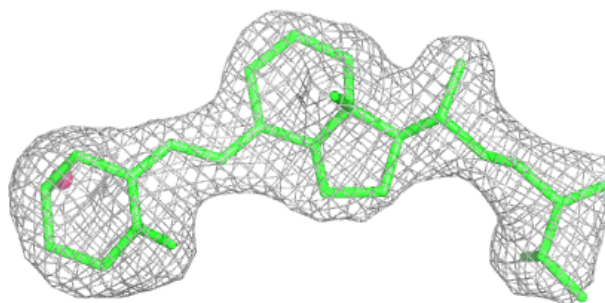
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around D2V A 602:

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

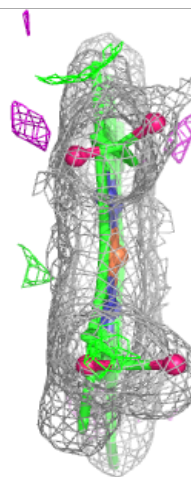
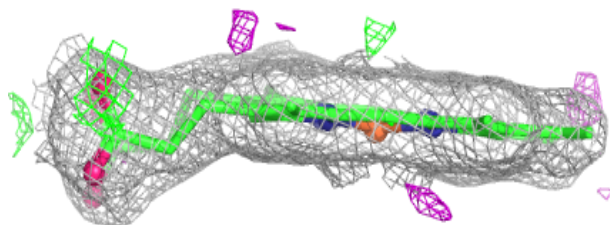
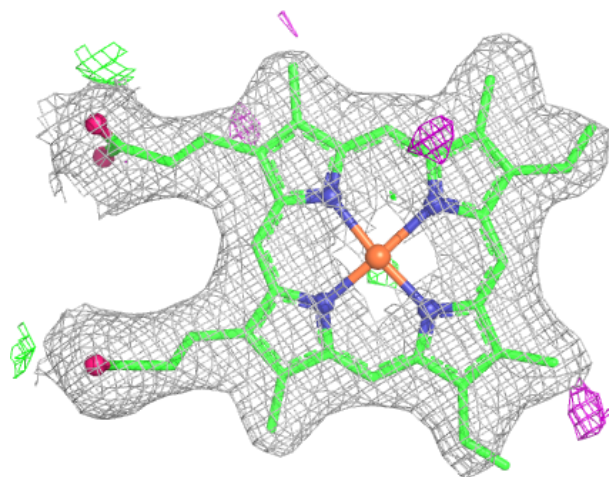
**Electron density around D2V B 602:**

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



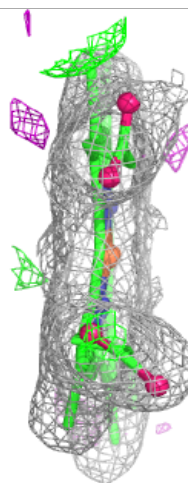
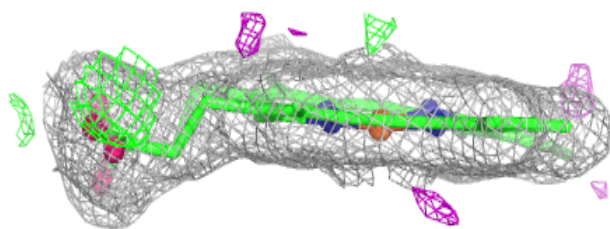
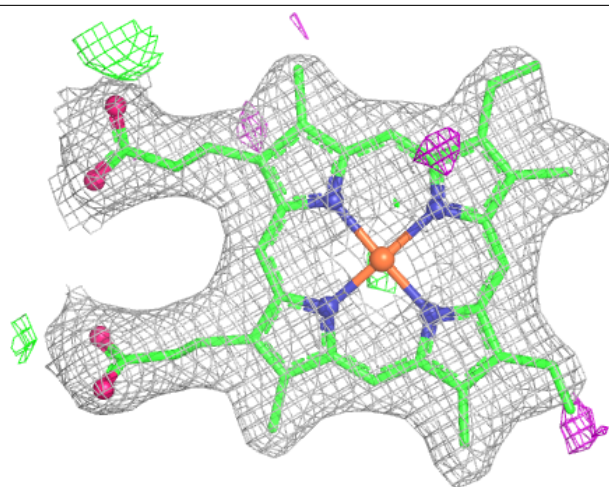
Electron density around HEM B 601 (A):

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



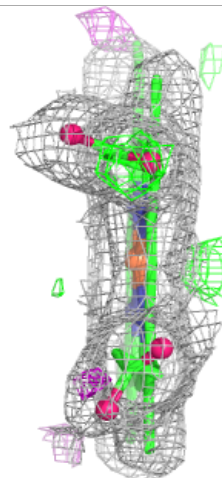
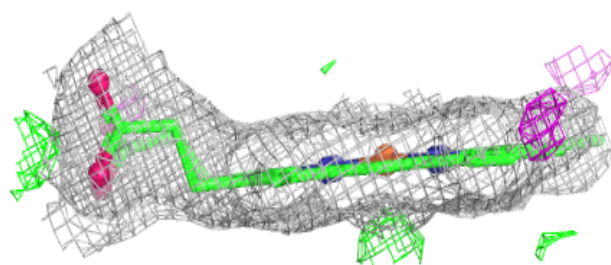
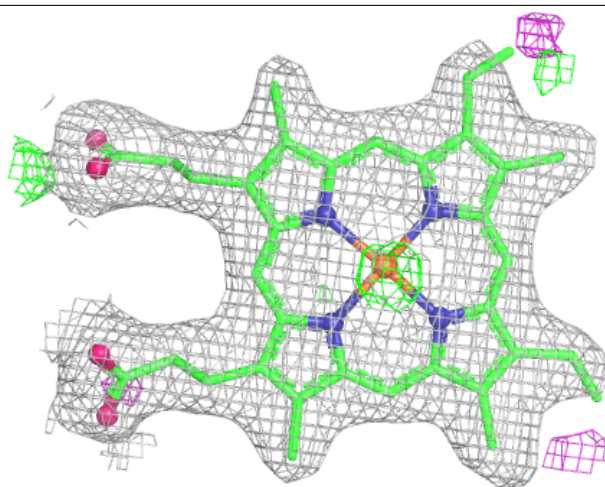
Electron density around HEM B 601 (B):

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



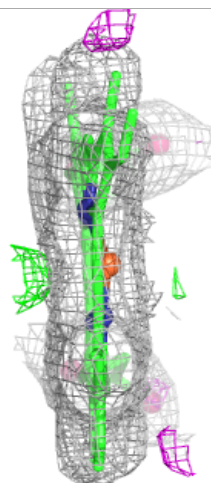
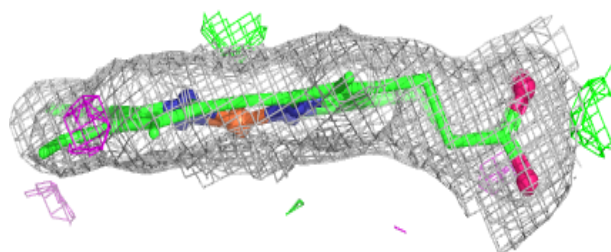
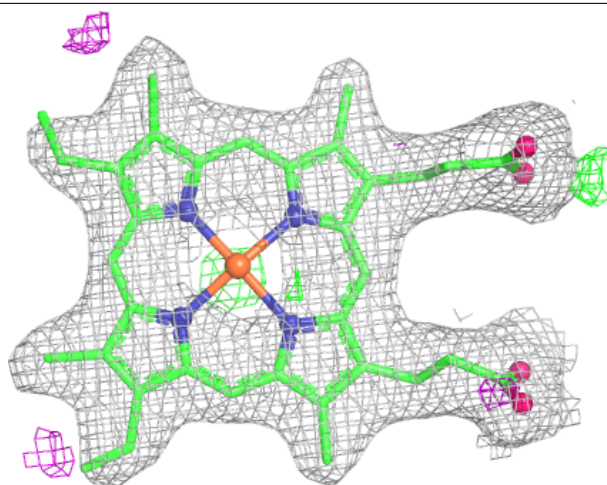
Electron density around HEM A 601 (A):

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



Electron density around HEM A 601 (B):

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



6.5 Other polymers ⓘ

There are no such residues in this entry.