



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 7, 2026 – 01:34 AM UTC

PDB ID : 3V6Q / pdb_00003v6q
Title : Crystal structure of the complex of bovine lactoperoxidase with Carbon monoxide at 2.0 Å resolution
Authors : Yamini, S.; Singh, A.K.; Pandey, N.; Sinha, M.; Kaur, P.; Sharma, S.; Singh, T.P.
Deposited on : 2011-12-20
Resolution : 2.00 Å(reported)

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

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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

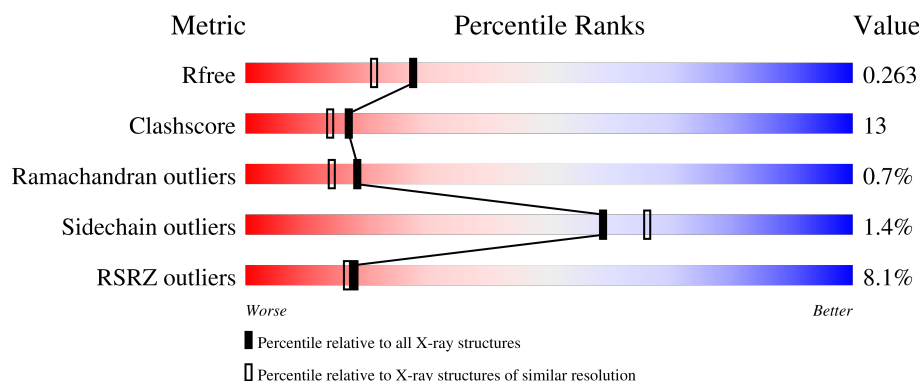
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

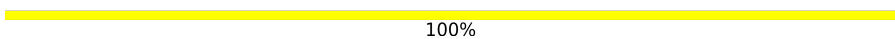
The reported resolution of this entry is 2.00 Å.

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	10052 (2.00-2.00)
Clashscore	190562	11152 (2.00-2.00)
Ramachandran outliers	187476	11031 (2.00-2.00)
Sidechain outliers	187428	11029 (2.00-2.00)
RSRZ outliers	180081	10067 (2.00-2.00)

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	595	
2	B	2	

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

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
11	CMO	A	621	-	-	X	-
6	IOD	A	615	-	-	X	-

2 Entry composition [i](#)

There are 12 unique types of molecules in this entry. The entry contains 5417 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 Lactoperoxidase.

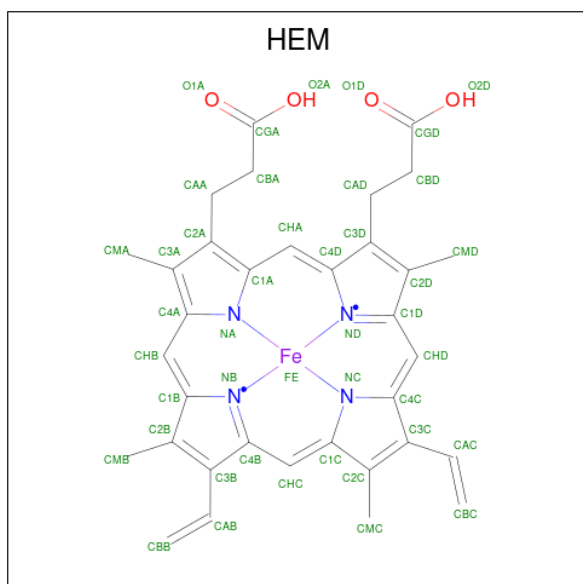
Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	595	Total	C	N	O	P	S	0	0	0
			4774	3037	847	863	1	26			

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



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

- Molecule 3 is PROTOPORPHYRIN IX CONTAINING FE (CCD ID: HEM) (formula: $C_{34}H_{32}FeN_4O_4$).

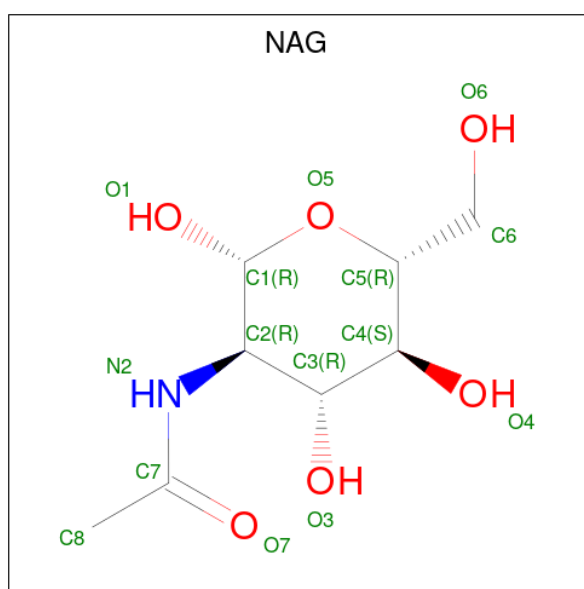


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	Fe	N	O	
			43	34	1	4	4	
								0
								0

- Molecule 4 is CALCIUM ION (CCD ID: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total	Ca		
			1	1	0	0

- Molecule 5 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: C₈H₁₅NO₆).



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

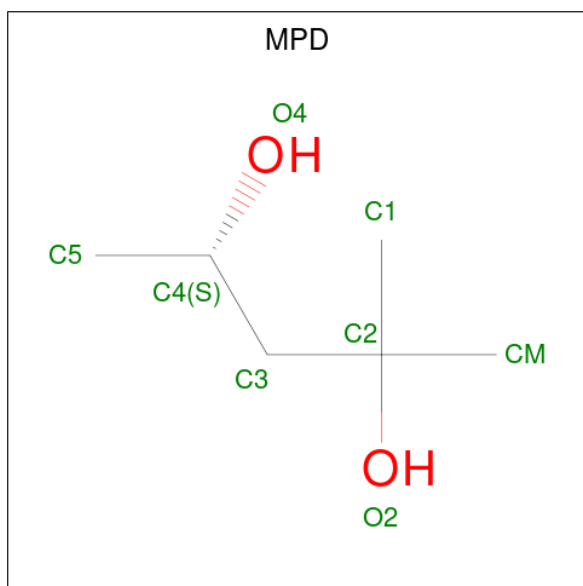
- Molecule 6 is IODIDE ION (CCD ID: IOD) (formula: I).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	9	Total	I		
			9	9	0	0

- Molecule 7 is BROMIDE ION (CCD ID: BR) (formula: Br).

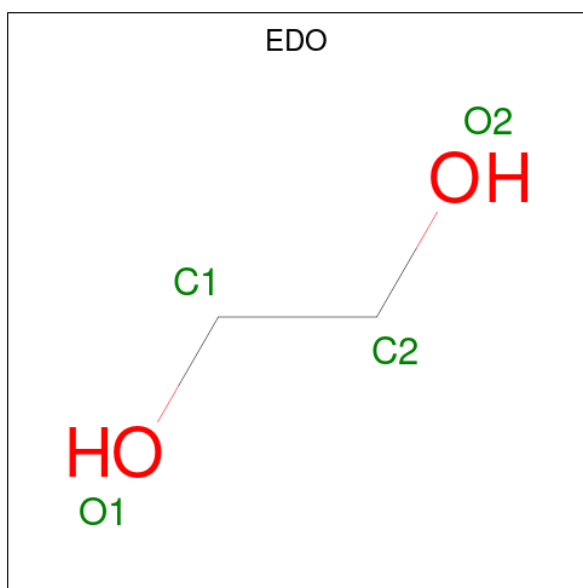
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	1	Total	Br	0	0
			1	1		

- Molecule 8 is (4S)-2-METHYL-2,4-PENTANEDIOL (CCD ID: MPD) (formula: $C_6H_{14}O_2$).



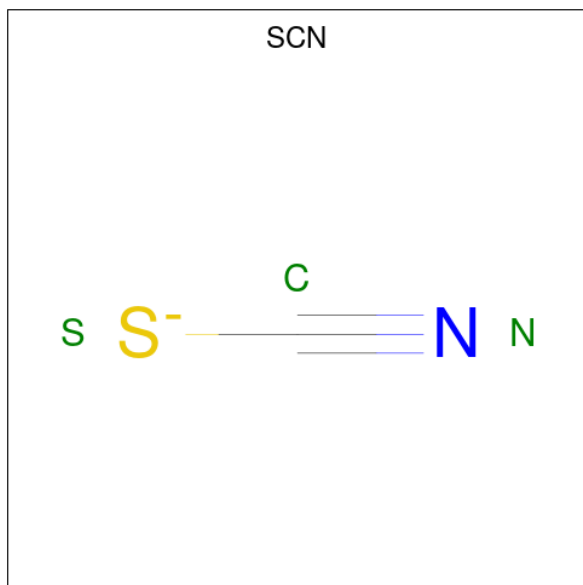
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
8	A	1	Total	C	O	0	0
			8	6	2		

- Molecule 9 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: $C_2H_6O_2$).



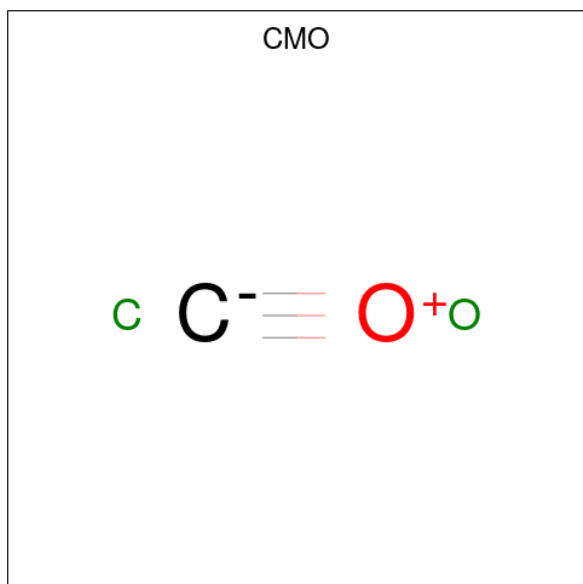
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
9	A	1	Total	C	O	0	0
			4	2	2		
9	A	1	Total	C	O	0	0
			4	2	2		

- Molecule 10 is THIOCYANATE ION (CCD ID: SCN) (formula: CNS).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
10	A	1	Total	C	N	S	0	0
			3	1	1	1		

- Molecule 11 is CARBON MONOXIDE (CCD ID: CMO) (formula: CO).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
11	A	1	Total	C	O	0	0
			2	1	1		

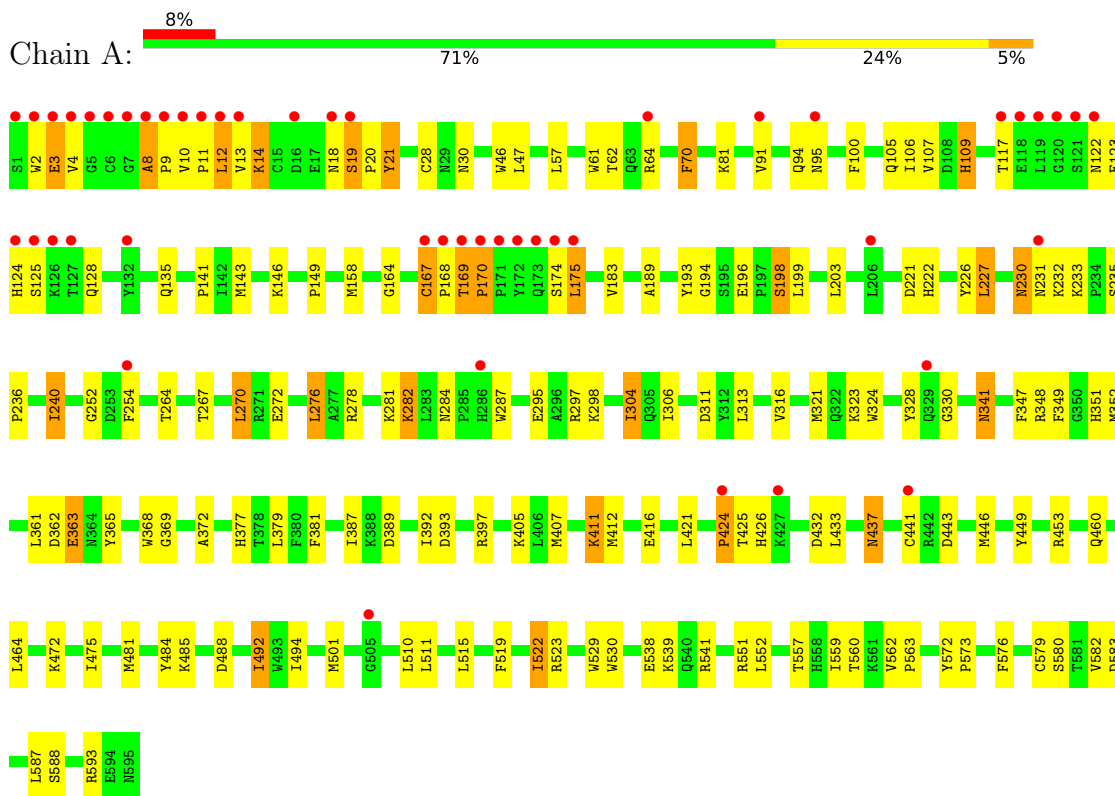
- Molecule 12 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
12	A	512	Total	O	0	0
			512	512		

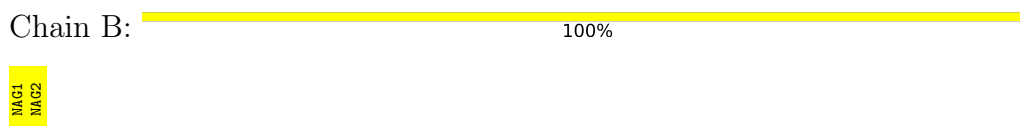
3 Residue-property plots [i](#)

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: Lactoperoxidase



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



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	54.30Å 80.00Å 76.55Å 90.00° 103.02° 90.00°	Depositor
Resolution (Å)	44.13 – 2.00 44.13 – 2.00	Depositor EDS
% Data completeness (in resolution range)	99.6 (44.13-2.00) 99.6 (44.13-2.00)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.05	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.71 (at 2.00Å)	Xtriage
Refinement program	REFMAC 5.5.0109	Depositor
R, R_{free}	0.191 , 0.248 0.224 , 0.263	Depositor DCC
R_{free} test set	2161 reflections (5.04%)	wwPDB-VP
Wilson B-factor (Å ²)	28.2	Xtriage
Anisotropy	0.620	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 44.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	5417	wwPDB-VP
Average B, all atoms (Å ²)	36.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.45% 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: MPD, CMO, CA, IOD, EDO, NAG, SCN, SEP, BR, HEM

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	1.56	45/4891 (0.9%)	1.34	57/6634 (0.9%)

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

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	2

All (45) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	313	LEU	C-O	-8.73	1.16	1.24
1	A	460	GLN	C-O	-7.78	1.17	1.24
1	A	47	LEU	C-O	-7.25	1.16	1.24
1	A	488	ASP	C-O	-6.72	1.15	1.24
1	A	10	VAL	CA-C	6.35	1.57	1.52
1	A	515	LEU	C-O	-6.16	1.17	1.24
1	A	272	GLU	C-O	-6.12	1.17	1.24
1	A	494	ILE	C-O	-6.10	1.16	1.24
1	A	106	ILE	C-O	-6.07	1.17	1.24
1	A	484	TYR	C-O	-6.03	1.16	1.23
1	A	464	LEU	C-O	-5.98	1.17	1.24
1	A	194	GLY	N-CA	5.92	1.51	1.45
1	A	316	VAL	C-O	-5.89	1.17	1.24
1	A	510	LEU	C-O	-5.83	1.17	1.24
1	A	562	VAL	C-O	-5.82	1.19	1.24
1	A	426	HIS	CG-ND1	-5.81	1.31	1.38
1	A	298	LYS	C-O	-5.78	1.17	1.24
1	A	412	MET	C-O	-5.78	1.18	1.23

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	492	ILE	C-O	-5.75	1.17	1.24
1	A	270	LEU	C-O	-5.70	1.17	1.24
1	A	306	ILE	C-O	-5.67	1.17	1.24
1	A	437	ASN	C-O	-5.62	1.17	1.24
1	A	107	VAL	C-O	-5.43	1.18	1.24
1	A	511	LEU	C-O	-5.42	1.17	1.24
1	A	443	ASP	C-O	-5.36	1.17	1.24
1	A	57	LEU	C-O	-5.36	1.17	1.24
1	A	421	LEU	C-O	-5.28	1.17	1.23
1	A	411	LYS	C-O	-5.27	1.16	1.23
1	A	341	ASN	C-O	-5.25	1.18	1.24
1	A	552	LEU	C-O	-5.23	1.18	1.24
1	A	109	HIS	C-O	-5.22	1.17	1.24
1	A	426	HIS	CD2-NE2	-5.21	1.32	1.37
1	A	183	VAL	N-CA	5.21	1.52	1.46
1	A	276	LEU	C-O	-5.16	1.18	1.24
1	A	321	MET	C-O	-5.16	1.18	1.24
1	A	433	LEU	C-O	-5.15	1.18	1.24
1	A	70	PHE	C-O	-5.12	1.17	1.23
1	A	8	ALA	CA-C	5.10	1.59	1.52
1	A	304	ILE	C-O	-5.09	1.18	1.24
1	A	453	ARG	C-O	-5.07	1.18	1.24
1	A	449	TYR	C-O	-5.05	1.18	1.24
1	A	379	LEU	C-O	-5.04	1.17	1.24
1	A	61	TRP	C-O	-5.04	1.18	1.24
1	A	143	MET	C-O	-5.03	1.17	1.23
1	A	100	PHE	C-O	-5.01	1.18	1.24

All (57) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	2	TRP	N-CA-C	16.80	129.59	111.28
1	A	481	MET	CB-CA-C	-16.42	84.69	111.13
1	A	14	LYS	N-CA-C	11.73	129.00	109.76
1	A	230	ASN	CA-C-N	11.73	138.41	120.82
1	A	230	ASN	C-N-CA	11.73	138.41	120.82
1	A	231	ASN	N-CA-C	11.11	129.71	111.37
1	A	240	ILE	CB-CA-C	11.07	126.14	111.97
1	A	593	ARG	N-CA-C	-10.11	90.85	108.23
1	A	174	SER	N-CA-C	9.81	125.40	113.23
1	A	231	ASN	CB-CA-C	-9.26	95.37	110.74
1	A	230	ASN	CB-CA-C	8.71	124.55	109.65

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	170	PRO	CB-CA-C	-8.34	100.74	110.92
1	A	230	ASN	N-CA-C	-7.73	98.84	110.28
1	A	562	VAL	N-CA-C	7.34	117.30	109.01
1	A	582	VAL	CB-CA-C	7.30	121.71	111.34
1	A	124	HIS	N-CA-C	-7.26	104.55	113.19
1	A	330	GLY	N-CA-C	7.23	118.82	111.95
1	A	254	PHE	N-CA-C	7.17	121.28	112.54
1	A	28	CYS	CB-CA-C	-7.07	100.43	111.39
1	A	81	LYS	CB-CA-C	6.98	122.52	110.79
1	A	174	SER	CB-CA-C	-6.97	95.40	109.55
1	A	484	TYR	CA-C-N	-6.94	109.16	122.06
1	A	484	TYR	C-N-CA	-6.94	109.16	122.06
1	A	328	TYR	N-CA-C	6.79	120.15	109.96
1	A	169	THR	CB-CA-C	-6.69	96.99	110.17
1	A	369	GLY	CA-C-N	6.58	126.16	119.19
1	A	369	GLY	C-N-CA	6.58	126.16	119.19
1	A	387	ILE	CB-CA-C	-6.54	103.20	112.22
1	A	588	SER	CA-C-N	-6.48	112.93	120.12
1	A	588	SER	C-N-CA	-6.48	112.93	120.12
1	A	240	ILE	N-CA-C	-6.24	104.43	110.42
1	A	282	LYS	N-CA-C	-6.23	104.54	111.71
1	A	235	SER	CA-C-N	-6.21	113.28	119.56
1	A	235	SER	C-N-CA	-6.21	113.28	119.56
1	A	46	TRP	N-CA-C	-6.20	105.20	112.89
1	A	284	ASN	CA-C-N	-6.09	114.00	120.94
1	A	284	ASN	C-N-CA	-6.09	114.00	120.94
1	A	484	TYR	O-C-N	-6.05	116.11	123.19
1	A	170	PRO	CA-C-N	5.96	125.97	119.90
1	A	170	PRO	C-N-CA	5.96	125.97	119.90
1	A	484	TYR	N-CA-C	-5.82	100.21	109.76
1	A	3	GLU	CB-CA-C	-5.71	100.56	110.45
1	A	146	LYS	N-CA-C	5.62	118.17	111.71
1	A	351	HIS	N-CA-C	5.51	118.05	111.71
1	A	170	PRO	N-CA-C	5.50	117.41	110.70
1	A	19	SER	CA-C-N	5.45	126.65	119.84
1	A	19	SER	C-N-CA	5.45	126.65	119.84
1	A	30	ASN	CB-CA-C	5.43	118.71	109.53
1	A	175	LEU	N-CA-C	-5.38	101.32	108.74
1	A	311	ASP	N-CA-C	5.38	120.49	113.88
1	A	472	LYS	CB-CA-C	5.37	121.21	111.06
1	A	30	ASN	N-CA-C	-5.26	101.89	110.20
1	A	426	HIS	N-CA-C	5.22	119.65	113.28

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	389	ASP	N-CA-C	5.14	117.46	110.88
1	A	21	TYR	CB-CA-C	5.08	119.51	109.66
1	A	12	LEU	CB-CA-C	-5.07	110.74	116.63
1	A	562	VAL	CB-CA-C	-5.04	104.45	109.89

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	198	SEP	Mainchain
1	A	230	ASN	Peptide

5.2 Too-close contacts

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	4774	0	4687	119	0
2	B	28	0	25	0	0
3	A	43	0	30	3	0
4	A	1	0	0	0	0
5	A	28	0	26	0	0
6	A	9	0	0	5	0
7	A	1	0	0	0	0
8	A	8	0	13	4	0
9	A	8	0	12	2	0
10	A	3	0	0	0	0
11	A	2	0	0	2	0
12	A	512	0	0	1	0
All	All	5417	0	4793	123	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 13.

All (123) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:12:LEU:CB	1:A:13:VAL:HA	1.61	1.28
1:A:12:LEU:HB3	1:A:13:VAL:CA	1.71	1.20
1:A:12:LEU:CB	1:A:13:VAL:CA	2.30	1.06
1:A:227:LEU:HD22	1:A:270:LEU:CD2	1.87	1.03
1:A:167:CYS:CB	1:A:168:PRO:HD2	1.89	1.02
1:A:168:PRO:HB2	1:A:170:PRO:HD2	1.39	1.01
1:A:227:LEU:CD2	1:A:270:LEU:HD22	1.90	1.01
1:A:167:CYS:HB2	1:A:168:PRO:HD2	1.42	0.98
1:A:227:LEU:HD22	1:A:270:LEU:HD22	0.99	0.97
1:A:167:CYS:CB	1:A:168:PRO:CD	2.42	0.96
1:A:12:LEU:HB2	1:A:13:VAL:HB	1.53	0.90
1:A:167:CYS:HB3	1:A:168:PRO:CD	2.01	0.89
1:A:3:GLU:CD	1:A:4:VAL:H	1.81	0.88
1:A:9:PRO:HD3	1:A:167:CYS:O	1.74	0.87
1:A:3:GLU:HB2	1:A:175:LEU:HD22	1.57	0.84
1:A:551:ARG:HD3	1:A:583:ASP:O	1.80	0.82
1:A:12:LEU:HB2	1:A:13:VAL:CB	2.09	0.80
1:A:522:ILE:HD12	1:A:522:ILE:C	2.09	0.77
1:A:91:VAL:HG12	1:A:405:LYS:HG3	1.66	0.77
1:A:167:CYS:HB3	1:A:168:PRO:HD3	1.65	0.77
1:A:12:LEU:HB3	1:A:13:VAL:HA	0.79	0.75
1:A:123:GLU:C	1:A:125:SER:H	1.95	0.72
1:A:94:GLN:C	1:A:95:ASN:HD22	1.98	0.72
1:A:3:GLU:HB2	1:A:175:LEU:CD2	2.20	0.71
1:A:19:SER:O	1:A:21:TYR:N	2.22	0.70
1:A:236:PRO:HG3	1:A:424:PRO:HB3	1.74	0.70
1:A:362:ASP:OD1	1:A:362:ASP:C	2.35	0.69
1:A:227:LEU:HD23	1:A:227:LEU:N	2.07	0.69
1:A:117:THR:HG22	1:A:164:GLY:HA2	1.75	0.68
1:A:12:LEU:HB2	1:A:13:VAL:CA	2.24	0.66
3:A:601:HEM:HBC2	3:A:601:HEM:HMC1	1.76	0.65
1:A:381:PHE:CZ	1:A:424:PRO:HG3	2.33	0.64
1:A:199:LEU:HA	8:A:617:MPD:HM2	1.81	0.63
1:A:3:GLU:CD	1:A:4:VAL:N	2.56	0.63
1:A:227:LEU:CD2	1:A:270:LEU:CD2	2.64	0.63
1:A:169:THR:N	1:A:170:PRO:CD	2.62	0.62
1:A:522:ILE:HD12	1:A:523:ARG:N	2.13	0.62
1:A:278:ARG:O	1:A:282:LYS:HG3	1.99	0.62
1:A:169:THR:OG1	1:A:170:PRO:HD3	2.00	0.62
1:A:122:ASN:CG	1:A:122:ASN:O	2.42	0.61
1:A:12:LEU:CB	1:A:13:VAL:CB	2.72	0.61
1:A:123:GLU:C	1:A:125:SER:N	2.57	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:580:SER:CB	6:A:615:IOD:I	3.19	0.61
1:A:3:GLU:CG	1:A:4:VAL:N	2.64	0.60
1:A:323:LYS:HE3	1:A:324:TRP:CZ2	2.38	0.59
1:A:341:ASN:HB3	1:A:446:MET:HE1	1.84	0.59
1:A:368:TRP:O	1:A:372:ALA:HB2	2.03	0.58
1:A:3:GLU:HA	1:A:3:GLU:OE1	2.05	0.57
1:A:377:HIS:CD2	9:A:618:EDO:H21	2.41	0.56
1:A:530:TRP:CZ3	1:A:541:ARG:HG2	2.41	0.55
1:A:18:ASN:O	1:A:19:SER:C	2.49	0.55
1:A:563:PRO:HD3	1:A:576:PHE:CE2	2.42	0.55
1:A:377:HIS:ND1	1:A:416:GLU:OE2	2.35	0.54
1:A:425:THR:HG22	1:A:425:THR:O	2.08	0.54
1:A:167:CYS:HB3	1:A:168:PRO:HD2	1.71	0.53
1:A:19:SER:C	1:A:21:TYR:H	2.16	0.53
1:A:149:PRO:HB2	6:A:614:IOD:I	2.79	0.53
1:A:12:LEU:CD1	1:A:13:VAL:HG23	2.38	0.53
1:A:221:ASP:HB2	1:A:226:TYR:CZ	2.44	0.52
1:A:377:HIS:HD2	9:A:618:EDO:H21	1.73	0.51
1:A:323:LYS:HE3	1:A:324:TRP:CE2	2.46	0.51
1:A:91:VAL:CG2	1:A:411:LYS:HD3	2.41	0.51
1:A:199:LEU:CA	8:A:617:MPD:HM2	2.42	0.50
1:A:199:LEU:HB2	8:A:617:MPD:HM2	1.93	0.50
1:A:407:MET:HE3	1:A:407:MET:C	2.36	0.50
1:A:352:MET:CB	1:A:407:MET:HG2	2.42	0.50
1:A:393:ASP:OD1	1:A:557:THR:HB	2.11	0.50
1:A:193:TYR:OH	1:A:297:ARG:HA	2.11	0.50
1:A:441:CYS:SG	1:A:492:ILE:HG22	2.52	0.50
1:A:580:SER:HB3	6:A:615:IOD:I	2.82	0.50
1:A:122:ASN:O	1:A:122:ASN:ND2	2.44	0.49
1:A:407:MET:HB3	1:A:501:MET:CE	2.43	0.49
1:A:362:ASP:OD1	1:A:363:GLU:N	2.46	0.48
1:A:580:SER:HB2	6:A:615:IOD:I	2.83	0.48
1:A:91:VAL:HG22	1:A:411:LYS:CD	2.44	0.47
1:A:352:MET:HB2	1:A:407:MET:HG2	1.95	0.47
1:A:348:ARG:HH11	1:A:437:ASN:ND2	2.12	0.47
1:A:3:GLU:CG	1:A:4:VAL:H	2.23	0.47
1:A:3:GLU:HG3	1:A:4:VAL:N	2.29	0.47
3:A:601:HEM:CMB	3:A:601:HEM:HBB2	2.44	0.47
1:A:189:ALA:HB2	1:A:304:ILE:HD12	1.97	0.47
1:A:91:VAL:HG22	1:A:411:LYS:HD3	1.97	0.46
1:A:62:THR:HG22	1:A:64:ARG:HG2	1.97	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:559:ILE:HG23	6:A:611:IOD:I	2.85	0.46
1:A:8:ALA:HB3	1:A:167:CYS:HA	1.98	0.46
1:A:523:ARG:HG3	1:A:529:TRP:CE2	2.51	0.46
1:A:168:PRO:CB	1:A:170:PRO:HD2	2.28	0.45
1:A:538:GLU:HG3	1:A:539:LYS:N	2.32	0.44
1:A:196:GLU:HB3	1:A:198:SEP:O3P	2.18	0.44
1:A:362:ASP:O	1:A:365:TYR:N	2.44	0.44
1:A:3:GLU:CB	1:A:175:LEU:CD2	2.93	0.44
1:A:227:LEU:CD2	1:A:227:LEU:N	2.79	0.44
1:A:276:LEU:HD23	1:A:587:LEU:HD11	2.00	0.44
1:A:158:MET:HE1	1:A:432:ASP:H	1.83	0.43
1:A:94:GLN:C	1:A:95:ASN:ND2	2.72	0.43
1:A:227:LEU:HD21	1:A:267:THR:HA	2.00	0.43
1:A:70:PHE:CG	1:A:485:LYS:HB2	2.54	0.43
1:A:109:HIS:NE2	11:A:621:CMO:O	2.49	0.43
1:A:560:THR:HA	1:A:579:CYS:SG	2.59	0.43
1:A:13:VAL:HG22	1:A:14:LYS:N	2.33	0.43
1:A:281:LYS:O	1:A:282:LYS:C	2.62	0.43
1:A:9:PRO:CD	1:A:167:CYS:O	2.57	0.43
1:A:441:CYS:SG	1:A:446:MET:HG3	2.58	0.42
1:A:105:GLN:HE21	11:A:621:CMO:C	2.33	0.42
1:A:12:LEU:HD12	1:A:13:VAL:HG23	2.01	0.42
1:A:349:PHE:CD1	1:A:349:PHE:C	2.97	0.42
1:A:95:ASN:HD22	1:A:95:ASN:N	2.16	0.42
1:A:135:GLN:HB2	1:A:141:PRO:HD2	2.02	0.41
1:A:232:LYS:HG3	1:A:233:LYS:N	2.35	0.41
1:A:361:LEU:O	1:A:397:ARG:HD2	2.20	0.41
1:A:287:TRP:CZ3	1:A:295:GLU:HG2	2.55	0.41
1:A:519:PHE:HA	1:A:522:ILE:HG13	2.02	0.41
1:A:9:PRO:HG3	12:A:704:HOH:O	2.20	0.41
1:A:563:PRO:HD3	1:A:576:PHE:CD2	2.55	0.41
1:A:221:ASP:O	1:A:222:HIS:C	2.64	0.41
1:A:264:THR:HG23	1:A:392:ILE:HB	2.02	0.41
3:A:601:HEM:HBC2	3:A:601:HEM:CMC	2.46	0.41
1:A:341:ASN:HB3	1:A:446:MET:CE	2.51	0.41
1:A:175:LEU:HD12	1:A:175:LEU:HA	1.75	0.40
1:A:203:LEU:HD11	1:A:252:GLY:HA2	2.04	0.40
1:A:572:TYR:CG	1:A:573:PRO:HA	2.57	0.40
1:A:3:GLU:OE1	1:A:3:GLU:CA	2.65	0.40
8:A:617:MPD:H4	8:A:617:MPD:HM1	1.90	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	592/595 (100%)	568 (96%)	20 (3%)	4 (1%)	18 14

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	20	PRO
1	A	167	CYS
1	A	11	PRO
1	A	424	PRO

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	517/517 (100%)	510 (99%)	7 (1%)	59 66

All (7) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	128	GLN
1	A	227	LEU
1	A	240	ILE
1	A	347	PHE
1	A	363	GLU
1	A	475	ILE
1	A	522	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	95	ASN
1	A	128	GLN
1	A	138	ASN
1	A	147	ASN
1	A	426	HIS
1	A	468	GLN
1	A	497	ASN
1	A	520	GLN
1	A	556	ASN
1	A	595	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

1 non-standard protein/DNA/RNA residue is 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	$\# Z > 2$
1	SEP	A	198	1	8,9,10	0.78	0	7,12,14	1.16	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
1	SEP	A	198	1	-	4/6/8/10	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	198	SEP	CB-OG-P-O1P
1	A	198	SEP	CB-OG-P-O2P
1	A	198	SEP	CB-OG-P-O3P
1	A	198	SEP	CA-CB-OG-P

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	A	198	SEP	1	0

5.5 Carbohydrates [i](#)

2 monosaccharides are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	NAG	B	1	2,1	14,14,15	0.72	0	17,19,21	1.66	4 (23%)
2	NAG	B	2	2	14,14,15	0.98	0	17,19,21	2.59	6 (35%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	NAG	B	1	2,1	-	2/6/23/26	0/1/1/1
2	NAG	B	2	2	-	0/6/23/26	0/1/1/1

There are no bond length outliers.

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	2	NAG	C1-O5-C5	8.23	123.21	112.19
2	B	1	NAG	C2-N2-C7	3.43	127.50	122.90
2	B	2	NAG	O5-C1-C2	-3.42	105.99	111.29
2	B	1	NAG	O4-C4-C3	-3.13	103.01	110.38
2	B	2	NAG	C4-C3-C2	-3.00	106.62	111.02
2	B	2	NAG	O5-C5-C4	2.43	116.73	110.83
2	B	1	NAG	O3-C3-C2	-2.34	104.54	109.40
2	B	1	NAG	O6-C6-C5	-2.28	103.56	111.33
2	B	2	NAG	C1-C2-N2	2.09	113.73	110.43
2	B	2	NAG	O3-C3-C4	-2.02	105.62	110.38

There are no chirality outliers.

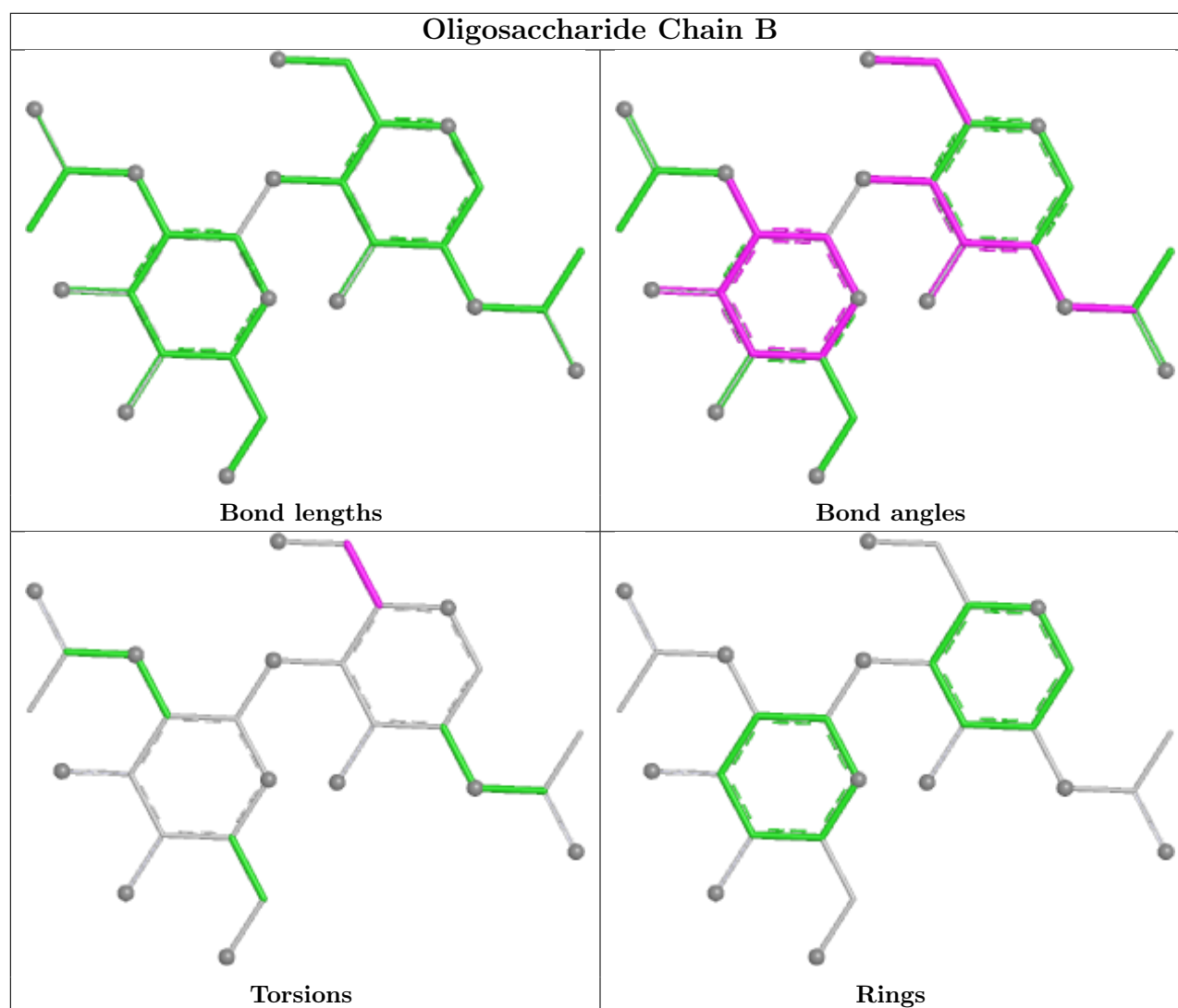
All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	1	NAG	O5-C5-C6-O6
2	B	1	NAG	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.



5.6 Ligand geometry [i](#)

Of 19 ligands modelled in this entry, 11 are monoatomic - leaving 8 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	$\# Z > 2$
3	HEM	A	601	11,1	50,50,50	2.22	10 (20%)	67,82,82	2.22	19 (28%)
5	NAG	A	606	1	14,14,15	0.70	0	17,19,21	1.29	2 (11%)
5	NAG	A	603	1	14,14,15	1.00	1 (7%)	17,19,21	1.70	6 (35%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
11	CMO	A	621	3	0,1,1	-	-	-		
9	EDO	A	618	-	3,3,3	1.06	0	2,2,2	1.26	0
8	MPD	A	617	-	7,7,7	4.85	3 (42%)	9,10,10	1.32	1 (11%)
10	SCN	A	620	-	1,2,2	1.07	0	0,1,1	-	-
9	EDO	A	619	-	3,3,3	0.64	0	2,2,2	1.17	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	HEM	A	601	11,1	-	4/14/54/54	-
5	NAG	A	606	1	-	2/6/23/26	0/1/1/1
5	NAG	A	603	1	-	2/6/23/26	0/1/1/1
9	EDO	A	618	-	-	0/1/1/1	-
8	MPD	A	617	-	-	2/5/5/5	-
9	EDO	A	619	-	-	1/1/1/1	-

All (14) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
8	A	617	MPD	C1-C2	-8.22	1.28	1.52
3	A	601	HEM	FE-NA	7.59	2.20	1.95
8	A	617	MPD	O2-C2	-6.94	1.27	1.44
8	A	617	MPD	CM-C2	-6.62	1.32	1.52
3	A	601	HEM	FE-NB	6.40	2.14	1.94
3	A	601	HEM	FE-ND	5.58	2.12	1.94
3	A	601	HEM	C3D-C2D	5.53	1.48	1.36
3	A	601	HEM	FE-NC	3.92	2.08	1.95
3	A	601	HEM	CAC-C3C	3.34	1.56	1.47
3	A	601	HEM	CMC-C2C	3.17	1.57	1.50
3	A	601	HEM	CAB-C3B	2.56	1.54	1.47
5	A	603	NAG	O7-C7	2.19	1.28	1.23
3	A	601	HEM	O1A-CGA	2.07	1.28	1.22
3	A	601	HEM	O1D-CGD	2.02	1.28	1.22

All (28) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	601	HEM	C4D-ND-C1D	6.86	113.33	105.21

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	601	HEM	C1B-NB-C4B	6.62	113.05	105.21
3	A	601	HEM	C3B-C4B-NB	-6.39	104.88	109.47
3	A	601	HEM	C2B-C1B-NB	-5.07	104.02	109.84
3	A	601	HEM	C3B-C2B-C1B	3.93	109.36	106.41
3	A	601	HEM	CHB-C1B-NB	3.77	129.03	124.37
5	A	603	NAG	O5-C5-C6	3.75	114.96	107.66
3	A	601	HEM	C4A-NA-C1A	3.18	111.00	105.82
5	A	606	NAG	C3-C4-C5	-3.14	104.54	110.23
3	A	601	HEM	CBA-CAA-C2A	-3.07	104.03	112.53
3	A	601	HEM	CHA-C4D-ND	2.96	128.02	124.37
3	A	601	HEM	C2A-C1A-NA	-2.93	106.90	110.15
3	A	601	HEM	CAD-CBD-CGD	-2.90	105.97	113.67
3	A	601	HEM	C2D-C1D-ND	-2.90	106.56	109.90
5	A	603	NAG	C8-C7-N2	-2.87	111.36	116.12
3	A	601	HEM	C3A-C4A-NA	-2.73	105.76	110.14
3	A	601	HEM	CMD-C2D-C1D	2.66	129.19	125.03
3	A	601	HEM	CHC-C4B-NB	2.64	127.27	124.42
3	A	601	HEM	C3D-C4D-ND	-2.56	107.37	110.17
5	A	606	NAG	O5-C5-C6	2.50	112.53	107.66
5	A	603	NAG	O3-C3-C4	-2.41	104.69	110.38
3	A	601	HEM	CBD-CAD-C3D	-2.40	105.89	112.53
8	A	617	MPD	O2-C2-C1	-2.36	100.63	107.99
5	A	603	NAG	C3-C4-C5	-2.35	105.97	110.23
3	A	601	HEM	C4C-NC-C1C	2.28	109.54	105.82
5	A	603	NAG	O7-C7-N2	2.20	125.87	121.98
3	A	601	HEM	C4C-C3C-C2C	2.02	108.57	106.81
5	A	603	NAG	C6-C5-C4	-2.02	108.07	113.02

There are no chirality outliers.

All (11) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
8	A	617	MPD	C2-C3-C4-O4
8	A	617	MPD	C2-C3-C4-C5
5	A	603	NAG	O5-C5-C6-O6
5	A	603	NAG	C4-C5-C6-O6
5	A	606	NAG	C4-C5-C6-O6
9	A	619	EDO	O1-C1-C2-O2
5	A	606	NAG	O5-C5-C6-O6
3	A	601	HEM	CAD-CBD-CGD-O2D
3	A	601	HEM	CAA-CBA-CGA-O2A
3	A	601	HEM	CAA-CBA-CGA-O1A

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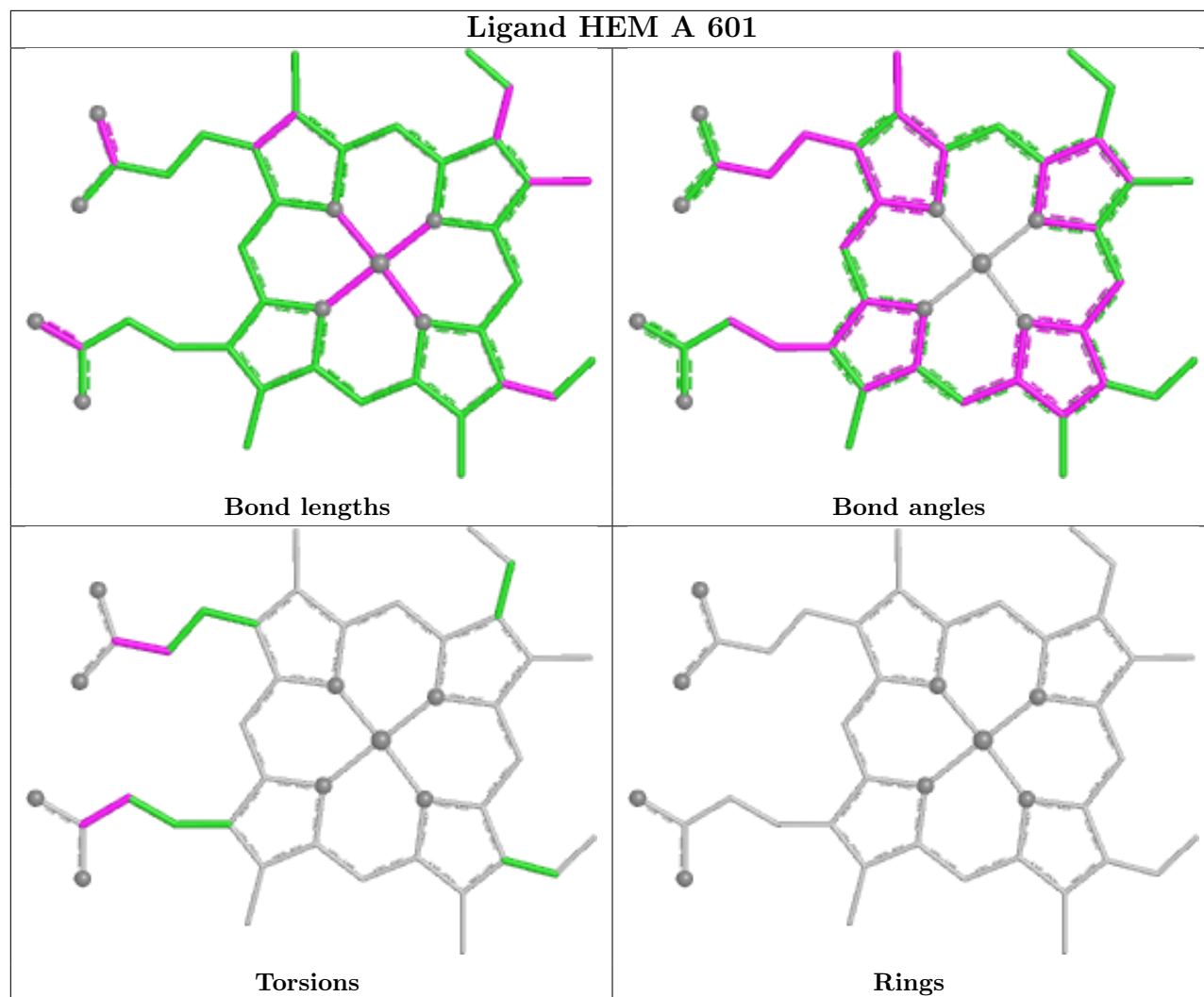
Mol	Chain	Res	Type	Atoms
3	A	601	HEM	CAD-CBD-CGD-O1D

There are no ring outliers.

4 monomers are involved in 11 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	601	HEM	3	0
11	A	621	CMO	2	0
9	A	618	EDO	2	0
8	A	617	MPD	4	0

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



5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	594/595 (99%)	0.39	48 (8%) 18 17	20, 32, 66, 99	0

All (48) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	13	VAL	7.2
1	A	10	VAL	7.1
1	A	8	ALA	6.8
1	A	12	LEU	6.8
1	A	11	PRO	6.7
1	A	2	TRP	6.4
1	A	172	TYR	6.0
1	A	168	PRO	5.6
1	A	4	VAL	5.6
1	A	9	PRO	5.5
1	A	174	SER	4.8
1	A	18	ASN	4.6
1	A	170	PRO	4.5
1	A	169	THR	4.4
1	A	119	LEU	4.2
1	A	175	LEU	4.1
1	A	120	GLY	4.0
1	A	5	GLY	3.9
1	A	1	SER	3.7
1	A	254	PHE	3.5
1	A	121	SER	3.4
1	A	441	CYS	3.4
1	A	117	THR	3.3
1	A	231	ASN	3.3
1	A	7	GLY	3.1
1	A	95	ASN	3.1
1	A	132	TYR	3.0

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Mol	Chain	Res	Type	RSRZ
1	A	173	GLN	2.9
1	A	427	LYS	2.9
1	A	16	ASP	2.9
1	A	126	LYS	2.8
1	A	171	PRO	2.8
1	A	125	SER	2.8
1	A	424	PRO	2.8
1	A	124	HIS	2.7
1	A	3	GLU	2.7
1	A	505	GLY	2.7
1	A	167	CYS	2.4
1	A	19	SER	2.4
1	A	6	CYS	2.3
1	A	329	GLN	2.2
1	A	127	THR	2.2
1	A	118	GLU	2.1
1	A	91	VAL	2.1
1	A	286	HIS	2.1
1	A	206	LEU	2.0
1	A	64	ARG	2.0
1	A	122	ASN	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [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
1	SEP	A	198	10/11	0.73	0.34	51,52,53,54	0

6.3 Carbohydrates [i](#)

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

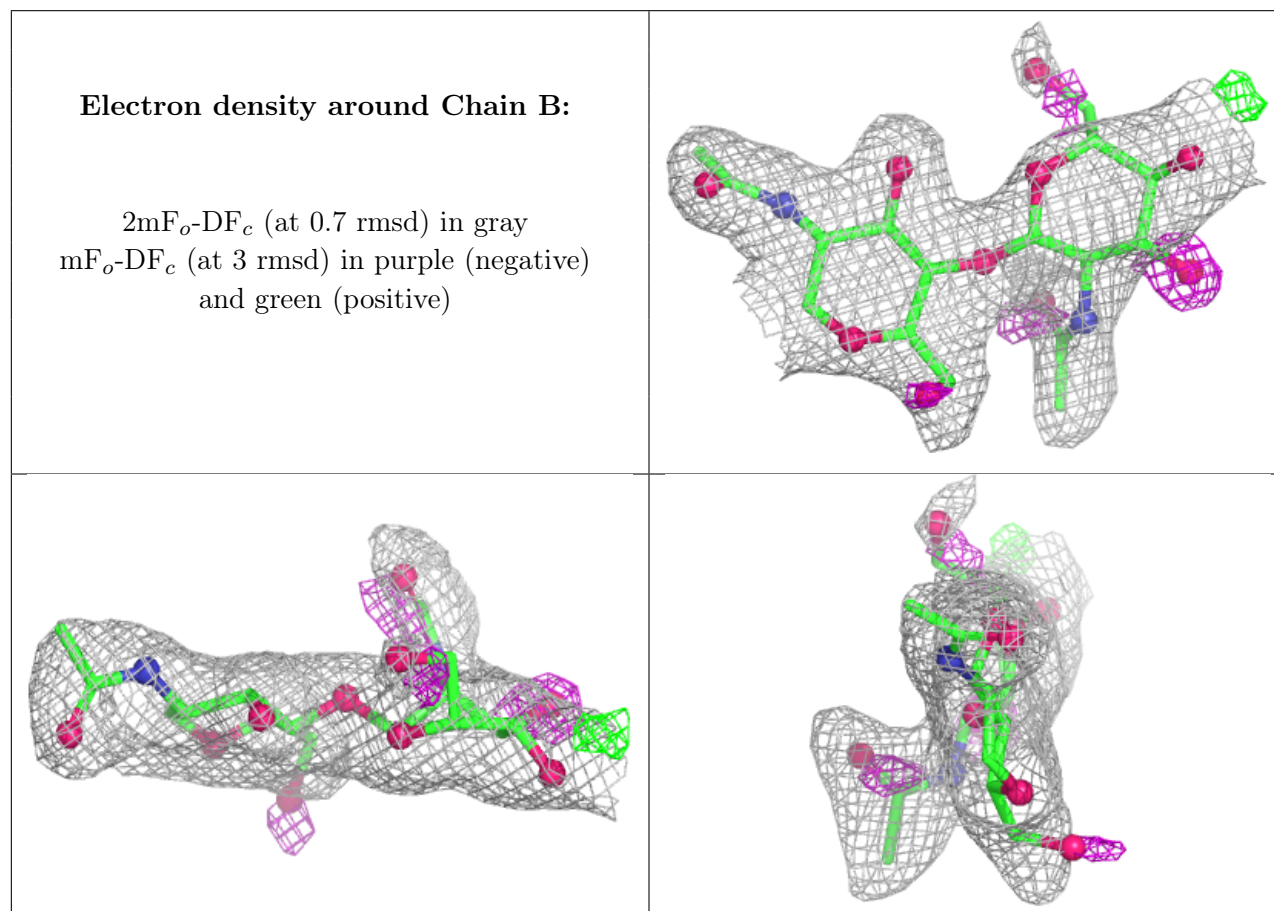
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	NAG	B	2	14/15	0.61	0.14	34,37,40,46	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	NAG	B	1	14/15	0.88	0.09	29,31,34,38	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.



6.4 Ligands [i](#)

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
9	EDO	A	618	4/4	0.66	0.26	26,26,27,28	0
9	EDO	A	619	4/4	0.69	0.19	26,28,33,34	0
5	NAG	A	606	14/15	0.76	0.11	31,34,38,41	0
8	MPD	A	617	8/8	0.76	0.26	18,26,33,37	0
5	NAG	A	603	14/15	0.83	0.09	25,32,34,37	0

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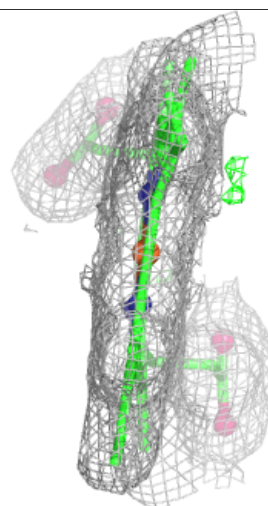
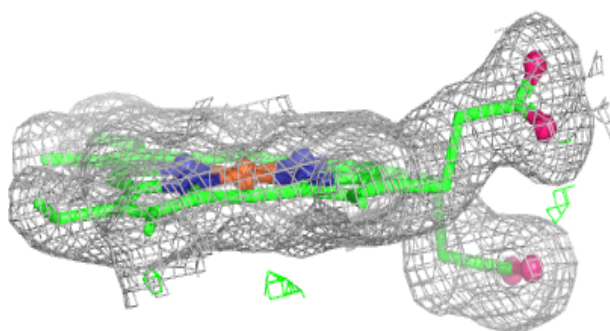
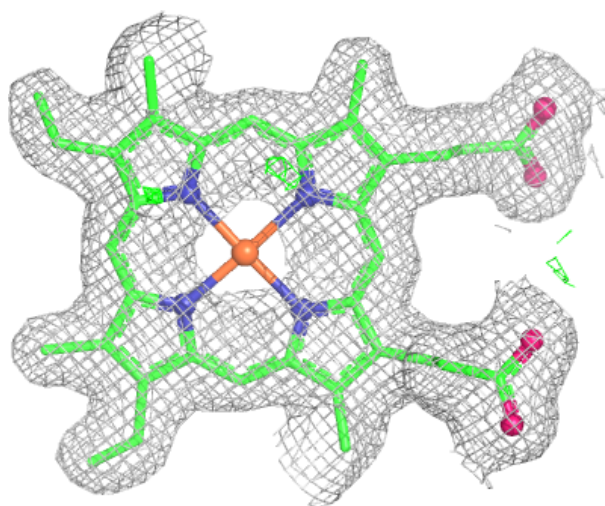
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
10	SCN	A	620	3/3	0.89	0.13	34,34,34,34	0
6	IOD	A	615	1/1	0.93	0.31	39,39,39,39	0
11	CMO	A	621	2/2	0.94	0.07	33,33,33,36	0
6	IOD	A	611	1/1	0.95	0.17	37,37,37,37	0
7	BR	A	616	1/1	0.96	0.04	29,29,29,29	0
6	IOD	A	607	1/1	0.97	0.04	23,23,23,23	0
6	IOD	A	612	1/1	0.98	0.14	35,35,35,35	0
6	IOD	A	614	1/1	0.98	0.34	39,39,39,39	0
3	HEM	A	601	43/43	0.98	0.05	16,22,26,27	0
6	IOD	A	609	1/1	0.99	0.19	36,36,36,36	0
6	IOD	A	610	1/1	0.99	0.04	31,31,31,31	0
4	CA	A	602	1/1	0.99	0.02	24,24,24,24	0
6	IOD	A	608	1/1	0.99	0.13	34,34,34,34	0
6	IOD	A	613	1/1	1.00	0.06	29,29,29,29	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 HEM A 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.