



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 5, 2026 – 11:58 AM EDT

PDB ID : 3BXI / pdb_00003bxi
Title : Structure of the complex of bovine lactoperoxidase with its catalyzed product
hypo thiocyanate ion at 2.3Å resolution
Authors : Singh, A.K.; Singh, N.; Sharma, S.; Shin, K.; Takase, M.; Kaur, P.; Srinivasan,
A.; Singh, T.P.
Deposited on : 2008-01-14
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	:	9.0.015 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.50

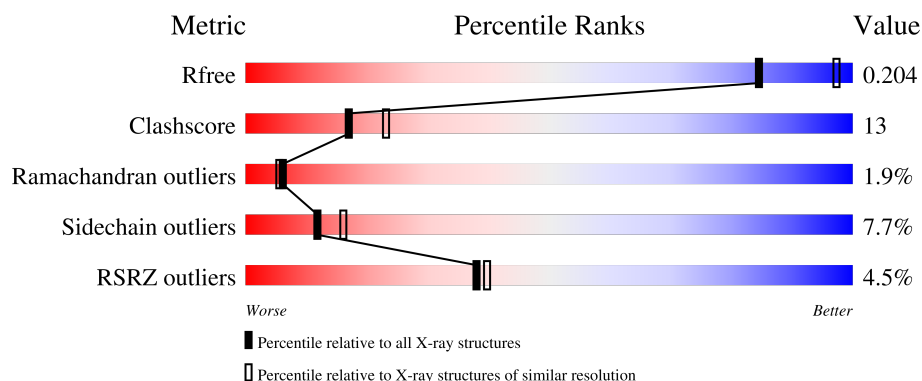
1 Overall quality at a glance

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	595	5% 69% 26% • •
2	B	3	67% 33%
2	D	3	33% 67%
3	C	2	50% 50%

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Mol	Chain	Length	Quality of chain
3	E	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
8	OSM	A	614	-	-	X	-

2 Entry composition [i](#)

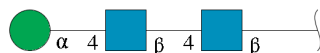
There are 9 unique types of molecules in this entry. The entry contains 5392 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 alpha-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



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

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

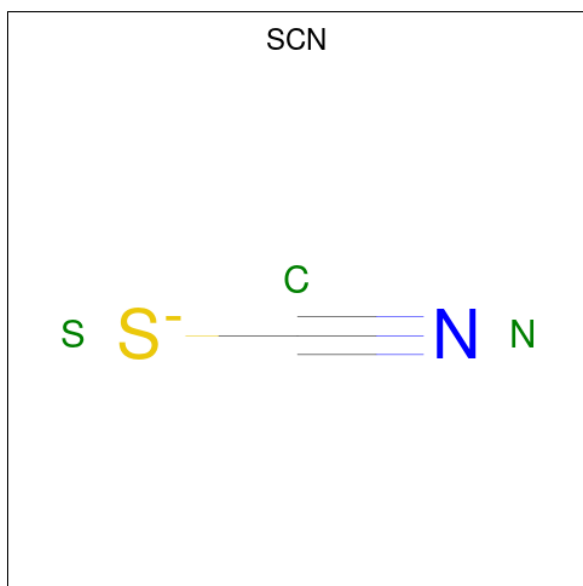


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	C	2	Total	C	N	O	0	0	0
			28	16	2	10			
3	E	2	Total	C	N	O	0	0	0
			28	16	2	10			

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

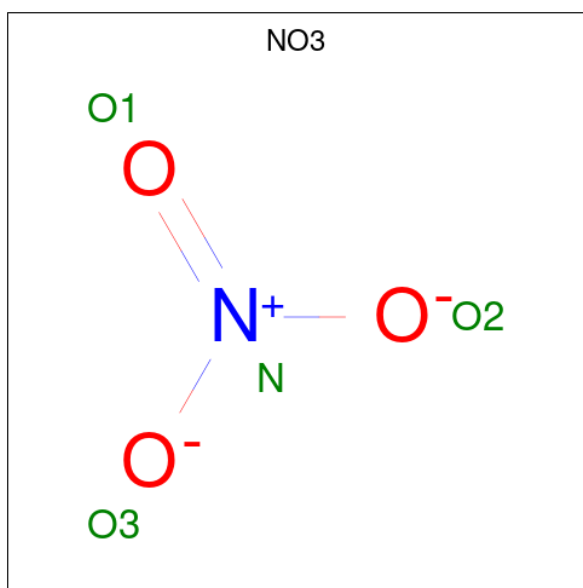
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total	Ca	0	0
			1	1		

- Molecule 5 is THIOCYANATE ION (CCD ID: SCN) (formula: CNS).



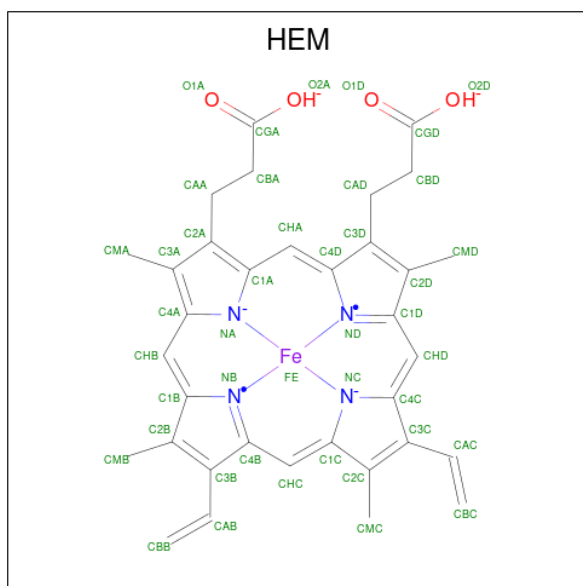
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	A	1	Total	C	N	S	0	0
			3	1	1	1		

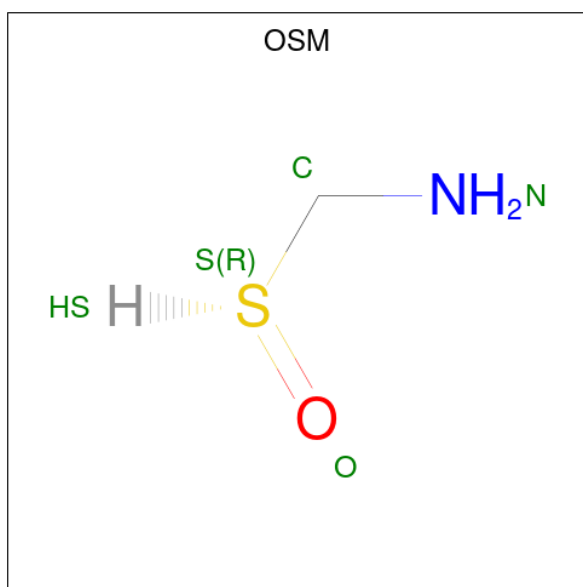
- Molecule 6 is NITRATE ION (CCD ID: NO3) (formula: NO₃).



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

- Molecule 7 is PROTOPORPHYRIN IX CONTAINING FE (CCD ID: HEM) (formula: $C_{34}H_{32}FeN_4O_4$).





Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
8	A	1	Total	C	N	O	S	0	0
			4	1	1	1	1		

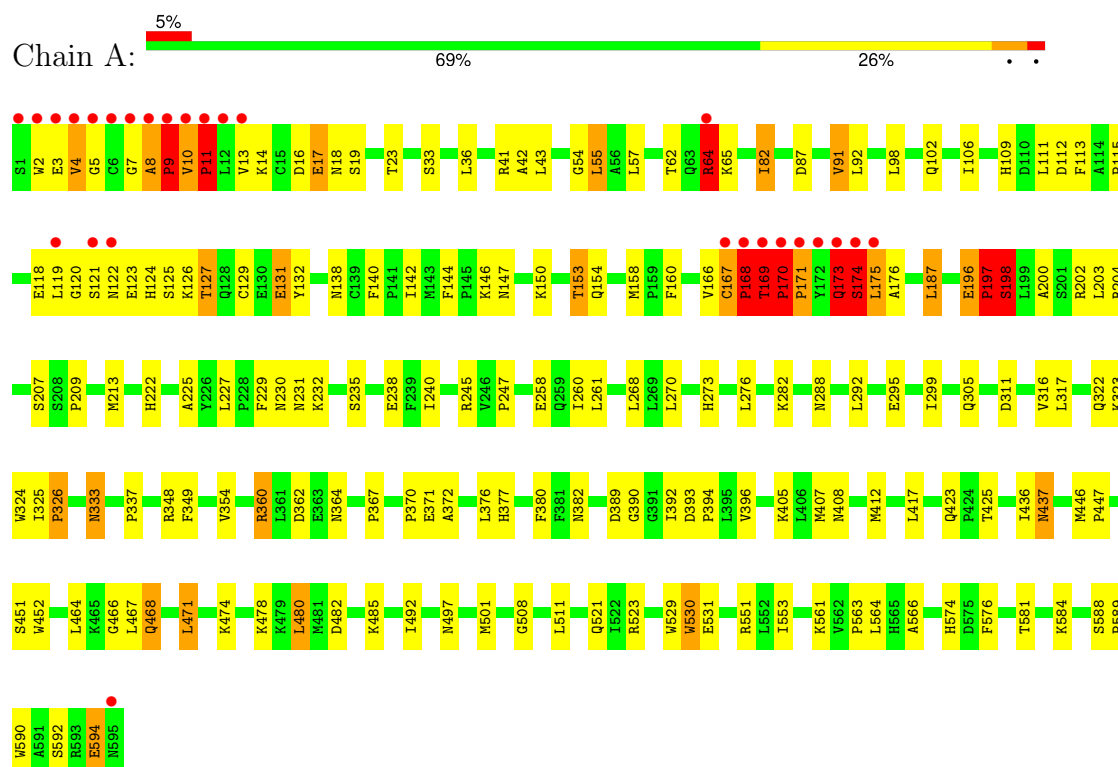
- Molecule 9 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
9	A	413	Total	O	0	0
			413	413		

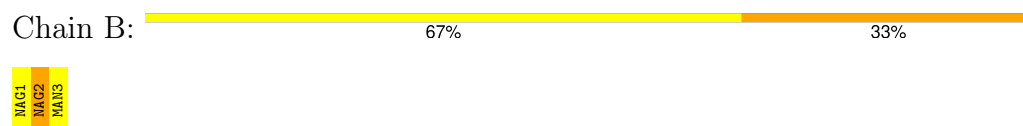
3 Residue-property plots [i](#)

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



• Molecule 2: alpha-D-mannopyranose-(1-4)-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

Chain C:  50% 50%

MAG1
MAG2

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

Chain E:  50% 50%

MAG1
MAG2

4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	54.63Å 80.67Å 77.68Å 90.00° 102.60° 90.00°	Depositor
Resolution (Å)	24.46 – 2.30 24.46 – 2.30	Depositor EDS
% Data completeness (in resolution range)	99.2 (24.46-2.30) 99.2 (24.46-2.30)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.41 (at 2.31Å)	Xtriage
Refinement program	REFMAC 5.0	Depositor
R, R_{free}	0.168 , 0.218 0.169 , 0.204	Depositor DCC
R_{free} test set	932 reflections (3.17%)	wwPDB-VP
Wilson B-factor (Å ²)	23.5	Xtriage
Anisotropy	0.105	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 48.0	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	5392	wwPDB-VP
Average B, all atoms (Å ²)	29.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.39% 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: SEP, OSM, CA, SCN, MAN, NO3, NAG, 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	0.68	10/4891 (0.2%)	1.28	66/6634 (1.0%)

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 (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	168	PRO	N-CA	9.73	1.59	1.47
1	A	10	VAL	CA-C	9.21	1.62	1.52
1	A	9	PRO	N-CA	7.47	1.56	1.47
1	A	196	GLU	CA-C	6.96	1.59	1.52
1	A	594	GLU	N-CA	6.49	1.54	1.46
1	A	258	GLU	N-CA	6.16	1.53	1.46
1	A	10	VAL	C-O	-5.69	1.17	1.24
1	A	167	CYS	CA-C	5.69	1.59	1.52
1	A	592	SER	N-CA	5.10	1.53	1.46
1	A	169	THR	C-N	5.09	1.39	1.33

All (66) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	10	VAL	CA-C-N	18.29	142.70	119.84
1	A	10	VAL	C-N-CA	18.29	142.70	119.84
1	A	10	VAL	N-CA-C	14.81	140.86	108.88

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	10	VAL	C-N-CD	-12.07	75.53	125.00
1	A	468	GLN	OE1-CD-NE2	-10.24	112.36	122.60
1	A	173	GLN	OE1-CD-NE2	-9.68	112.92	122.60
1	A	232	LYS	N-CA-C	9.63	124.20	110.23
1	A	316	VAL	N-CA-C	-9.41	102.73	111.67
1	A	2	TRP	N-CA-C	-8.93	97.09	109.95
1	A	9	PRO	CA-N-CD	-8.91	99.53	112.00
1	A	594	GLU	N-CA-C	8.87	122.69	111.24
1	A	174	SER	N-CA-C	8.62	129.17	110.80
1	A	168	PRO	CA-N-CD	-8.48	100.12	112.00
1	A	326	PRO	CA-N-CD	-7.98	100.83	112.00
1	A	167	CYS	CB-CA-C	7.45	124.84	110.17
1	A	260	ILE	N-CA-C	7.44	119.10	110.62
1	A	170	PRO	CA-N-CD	-7.38	101.67	112.00
1	A	508	GLY	CA-C-N	7.33	126.59	118.97
1	A	508	GLY	C-N-CA	7.33	126.59	118.97
1	A	437	ASN	CA-CB-CG	7.33	119.92	112.60
1	A	121	SER	N-CA-C	-7.28	95.21	107.80
1	A	173	GLN	CG-CD-NE2	6.98	126.86	116.40
1	A	225	ALA	N-CA-C	6.63	119.84	110.50
1	A	574	HIS	N-CA-C	6.34	118.72	111.11
1	A	468	GLN	CG-CD-NE2	6.34	125.91	116.40
1	A	592	SER	CA-C-N	6.26	129.72	120.95
1	A	592	SER	C-N-CA	6.26	129.72	120.95
1	A	566	ALA	N-CA-C	6.21	119.96	112.38
1	A	2	TRP	CB-CA-C	-6.21	99.62	111.48
1	A	167	CYS	O-C-N	-6.17	114.23	121.32
1	A	2	TRP	CA-C-N	-6.08	112.28	122.92
1	A	2	TRP	C-N-CA	-6.08	112.28	122.92
1	A	389	ASP	N-CA-C	6.07	117.72	110.44
1	A	7	GLY	N-CA-C	-6.00	98.97	113.18
1	A	390	GLY	N-CA-C	5.99	123.03	114.64
1	A	11	PRO	CA-N-CD	-5.98	103.63	112.00
1	A	197	PRO	CA-N-CD	-5.94	103.68	112.00
1	A	55	LEU	N-CA-C	5.92	120.80	112.93
1	A	33	SER	CA-C-N	5.86	126.27	119.47
1	A	33	SER	C-N-CA	5.86	126.27	119.47
1	A	447	PRO	N-CA-C	-5.85	102.63	111.41
1	A	111	LEU	N-CA-C	5.80	120.51	112.45
1	A	23	THR	N-CA-C	-5.79	103.06	110.53
1	A	167	CYS	CA-C-O	-5.73	112.31	120.16
1	A	530	TRP	N-CA-C	5.68	119.02	111.75

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	222	HIS	CA-CB-CG	-5.62	108.18	113.80
1	A	364	ASN	N-CA-C	-5.62	106.32	112.72
1	A	466	GLY	CA-C-N	5.49	127.95	120.54
1	A	466	GLY	C-N-CA	5.49	127.95	120.54
1	A	153	THR	N-CA-CB	-5.49	103.71	111.00
1	A	140	PHE	CA-C-N	5.30	125.22	119.76
1	A	140	PHE	C-N-CA	5.30	125.22	119.76
1	A	412	MET	N-CA-C	5.27	119.63	111.56
1	A	521	GLN	N-CA-C	5.23	117.06	111.36
1	A	43	LEU	N-CA-C	-5.18	103.19	110.50
1	A	112	ASP	N-CA-C	5.17	117.10	108.99
1	A	92	LEU	N-CA-C	5.15	117.76	110.50
1	A	11	PRO	N-CA-C	-5.13	101.89	112.47
1	A	10	VAL	O-C-N	5.13	126.95	121.10
1	A	167	CYS	N-CA-C	-5.13	98.48	109.81
1	A	138	ASN	N-CA-C	5.10	119.26	113.19
1	A	207	SER	N-CA-C	-5.04	105.49	111.69
1	A	87	ASP	N-CA-C	5.03	116.52	108.41
1	A	10	VAL	CB-CA-C	-5.03	102.20	111.36
1	A	127	THR	N-CA-C	-5.03	107.21	113.55
1	A	311	ASP	N-CA-C	5.01	119.39	113.12

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	197	PRO	Peptide
1	A	198	SEP	Mainchain

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	4685	117	0
2	B	39	0	34	1	0
2	D	39	0	34	2	0
3	C	28	0	25	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	E	28	0	25	2	0
4	A	1	0	0	0	0
5	A	3	0	0	0	0
6	A	20	0	0	1	0
7	A	43	0	30	6	0
8	A	4	0	5	3	0
9	A	413	0	0	24	0
All	All	5392	0	4838	125	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 (125) 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:170:PRO:HD3	9:A:1010:HOH:O	1.56	1.02
1:A:55:LEU:HD23	9:A:1014:HOH:O	1.66	0.95
1:A:4:VAL:HG13	1:A:4:VAL:O	1.71	0.90
1:A:295:GLU:O	1:A:299:ILE:HD13	1.76	0.86
1:A:8:ALA:CB	1:A:9:PRO:HD2	2.08	0.83
1:A:360:ARG:NH1	1:A:372:ALA:HA	1.96	0.81
1:A:360:ARG:HH12	1:A:372:ALA:HA	1.45	0.80
1:A:8:ALA:HB3	1:A:9:PRO:HD2	1.64	0.80
1:A:240:ILE:HD11	1:A:382:ASN:HA	1.65	0.78
1:A:119:LEU:HD11	1:A:169:THR:CG2	2.12	0.78
1:A:227:LEU:HD23	1:A:270:LEU:HD22	1.64	0.78
9:A:966:HOH:O	3:E:2:NAG:H82	1.84	0.77
1:A:13:VAL:HA	9:A:1004:HOH:O	1.83	0.76
1:A:150:LYS:HZ2	1:A:154:GLN:HE22	1.31	0.75
1:A:13:VAL:HG12	1:A:14:LYS:H	1.49	0.75
9:A:699:HOH:O	2:D:3:MAN:H61	1.87	0.75
1:A:196:GLU:HB3	1:A:198:SEP:O2P	1.88	0.73
1:A:8:ALA:CB	1:A:9:PRO:CD	2.65	0.73
1:A:123:GLU:HB3	1:A:126:LYS:HE2	1.70	0.72
1:A:561:LYS:HG3	9:A:768:HOH:O	1.90	0.70
1:A:123:GLU:HG3	1:A:125:SER:H	1.57	0.70
1:A:82:ILE:HD12	1:A:480:LEU:HD13	1.71	0.69
1:A:13:VAL:HG12	1:A:14:LYS:N	2.09	0.68
1:A:235:SER:HB3	1:A:238:GLU:HG2	1.77	0.67
1:A:4:VAL:O	1:A:4:VAL:CG1	2.40	0.67
1:A:407:MET:HB3	1:A:501:MET:HE3	1.79	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:150:LYS:NZ	1:A:154:GLN:HE22	1.96	0.64
1:A:333:ASN:C	1:A:333:ASN:HD22	2.04	0.64
7:A:613:HEM:C1A	8:A:614:OSM:S	2.91	0.63
1:A:446:MET:HE1	6:A:608:NO3:O3	2.00	0.61
1:A:119:LEU:HD11	1:A:169:THR:HG22	1.82	0.60
1:A:530:TRP:CE2	1:A:531:GLU:HG3	2.37	0.59
1:A:19:SER:HB3	9:A:1026:HOH:O	2.01	0.59
1:A:119:LEU:HD11	1:A:169:THR:HG23	1.84	0.59
1:A:16:ASP:HB2	9:A:1002:HOH:O	2.02	0.58
1:A:102:GLN:O	1:A:106:ILE:HD13	2.03	0.58
1:A:64:ARG:HH12	1:A:65:LYS:HE2	1.69	0.58
1:A:323:LYS:HE3	1:A:324:TRP:CZ2	2.39	0.58
1:A:322:GLN:HG3	9:A:798:HOH:O	2.05	0.57
9:A:818:HOH:O	3:E:2:NAG:C1	2.53	0.57
1:A:408:ASN:HB2	9:A:784:HOH:O	2.04	0.57
1:A:169:THR:HG22	9:A:1010:HOH:O	2.03	0.57
1:A:467:LEU:HG	1:A:471:LEU:HD22	1.86	0.57
1:A:62:THR:O	1:A:64:ARG:N	2.36	0.56
1:A:8:ALA:HB1	1:A:9:PRO:CD	2.35	0.56
1:A:102:GLN:HE21	1:A:106:ILE:CD1	2.18	0.56
1:A:8:ALA:HB1	1:A:9:PRO:HD2	1.86	0.56
1:A:362:ASP:OD1	1:A:362:ASP:C	2.48	0.55
1:A:213:MET:HG2	1:A:273:HIS:CD2	2.41	0.55
1:A:396:VAL:HG11	1:A:553:ILE:HD12	1.88	0.54
7:A:613:HEM:NA	8:A:614:OSM:S	2.80	0.54
1:A:62:THR:HG22	1:A:64:ARG:HB3	1.91	0.53
1:A:197:PRO:HB2	1:A:198:SEP:O3P	2.09	0.53
1:A:142:ILE:CD1	1:A:436:ILE:HD13	2.40	0.52
1:A:425:THR:HB	9:A:1007:HOH:O	2.10	0.52
1:A:127:THR:O	1:A:131:GLU:HB2	2.10	0.51
1:A:17:GLU:HG3	9:A:1027:HOH:O	2.11	0.50
1:A:8:ALA:HB3	1:A:9:PRO:CD	2.37	0.50
1:A:142:ILE:HD11	1:A:436:ILE:HD13	1.93	0.50
1:A:588:SER:N	1:A:589:PRO:CD	2.75	0.50
1:A:19:SER:CB	9:A:1026:HOH:O	2.58	0.50
1:A:551:ARG:HD3	1:A:584:LYS:HA	1.94	0.49
1:A:118:GLU:HG2	1:A:120:GLY:H	1.77	0.49
1:A:42:ALA:HB2	1:A:166:VAL:HG11	1.94	0.49
1:A:407:MET:HB3	1:A:501:MET:CE	2.43	0.48
1:A:16:ASP:HB3	9:A:1026:HOH:O	2.14	0.48
1:A:150:LYS:NZ	1:A:154:GLN:NE2	2.60	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:54:GLY:HA2	9:A:797:HOH:O	2.13	0.48
1:A:10:VAL:HG11	1:A:41:ARG:CZ	2.43	0.48
1:A:160:PHE:HD1	1:A:436:ILE:CD1	2.26	0.48
1:A:118:GLU:C	1:A:120:GLY:N	2.71	0.47
1:A:144:PHE:HE2	1:A:158:MET:HE3	1.79	0.47
1:A:245:ARG:HD3	9:A:820:HOH:O	2.14	0.47
1:A:468:GLN:HG2	1:A:474:LYS:HA	1.96	0.47
1:A:417:LEU:HD21	7:A:613:HEM:HMB3	1.97	0.47
1:A:36:LEU:HG	1:A:337:PRO:HD2	1.97	0.47
1:A:3:GLU:C	1:A:5:GLY:H	2.22	0.47
1:A:392:ILE:C	1:A:392:ILE:HD12	2.40	0.46
1:A:348:ARG:HH11	1:A:437:ASN:ND2	2.13	0.46
1:A:333:ASN:C	1:A:333:ASN:ND2	2.74	0.46
1:A:282:LYS:HB2	1:A:282:LYS:HE3	1.56	0.46
7:A:613:HEM:HBC2	7:A:613:HEM:HMC2	1.98	0.45
1:A:131:GLU:HB3	1:A:132:TYR:CD1	2.51	0.45
9:A:994:HOH:O	2:D:2:NAG:H3	2.16	0.45
1:A:299:ILE:HD11	1:A:590:TRP:HE1	1.82	0.45
1:A:282:LYS:HD2	9:A:979:HOH:O	2.15	0.45
1:A:113:PHE:O	1:A:115:PRO:HD3	2.18	0.44
1:A:230:ASN:OD1	1:A:231:ASN:N	2.50	0.44
1:A:13:VAL:CG1	1:A:14:LYS:N	2.79	0.44
1:A:563:PRO:HD3	1:A:576:PHE:CE2	2.53	0.43
1:A:417:LEU:CD2	7:A:613:HEM:HMB3	2.48	0.43
1:A:478:LYS:HB2	1:A:478:LYS:HE2	1.41	0.43
1:A:62:THR:C	1:A:64:ARG:N	2.75	0.43
1:A:131:GLU:HB3	1:A:132:TYR:CE1	2.53	0.43
1:A:146:LYS:O	1:A:147:ASN:HB2	2.18	0.43
1:A:354:VAL:HG11	7:A:613:HEM:CBB	2.48	0.43
1:A:229:PHE:CD1	1:A:247:PRO:HG2	2.54	0.43
1:A:109:HIS:NE2	8:A:614:OSM:S	2.89	0.42
1:A:91:VAL:HG13	1:A:405:LYS:HG3	2.01	0.42
1:A:523:ARG:HG3	1:A:529:TRP:CE2	2.53	0.42
1:A:423:GLN:HE21	1:A:423:GLN:HA	1.84	0.42
1:A:377:HIS:HA	1:A:380:PHE:CE2	2.55	0.42
1:A:392:ILE:HD12	1:A:393:ASP:N	2.34	0.42
1:A:187:LEU:HD13	1:A:305:GLN:HA	2.01	0.42
1:A:299:ILE:HD11	1:A:590:TRP:NE1	2.35	0.42
1:A:425:THR:CB	9:A:1007:HOH:O	2.67	0.42
9:A:878:HOH:O	2:B:2:NAG:H62	2.20	0.42
1:A:229:PHE:CG	1:A:247:PRO:HG2	2.55	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:348:ARG:HH11	1:A:437:ASN:HD22	1.68	0.42
1:A:200:ALA:O	1:A:204:ARG:HG3	2.19	0.42
1:A:349:PHE:HA	1:A:497:ASN:HD21	1.83	0.41
1:A:288:ASN:O	1:A:292:LEU:HD23	2.21	0.41
1:A:17:GLU:N	9:A:1026:HOH:O	2.52	0.41
1:A:175:LEU:CD2	1:A:176:ALA:H	2.34	0.41
1:A:370:PRO:HG2	1:A:371:GLU:HG3	2.03	0.41
1:A:452:TRP:CD1	1:A:492:ILE:HD13	2.55	0.41
1:A:173:GLN:CG	1:A:174:SER:H	2.34	0.41
1:A:10:VAL:HG21	1:A:41:ARG:NH2	2.36	0.40
1:A:581:THR:O	1:A:581:THR:HG22	2.21	0.40
1:A:129:CYS:HB2	9:A:900:HOH:O	2.21	0.40
1:A:393:ASP:HB2	1:A:394:PRO:HD3	2.02	0.40
1:A:202:ARG:NH2	1:A:231:ASN:HB2	2.36	0.40
1:A:160:PHE:HD1	1:A:436:ILE:HD12	1.87	0.40
1:A:482:ASP:O	1:A:485:LYS:NZ	2.47	0.40
1:A:173:GLN:HG2	1:A:174:SER:H	1.87	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%)	557 (94%)	24 (4%)	11 (2%)	6 5

All (11) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	8	ALA
1	A	11	PRO
1	A	167	CYS
1	A	168	PRO

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Mol	Chain	Res	Type
1	A	169	THR
1	A	174	SER
1	A	122	ASN
1	A	170	PRO
1	A	173	GLN
1	A	64	ARG
1	A	171	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%)	477 (92%)	40 (8%)	12	16

All (40) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	4	VAL
1	A	9	PRO
1	A	11	PRO
1	A	17	GLU
1	A	18	ASN
1	A	57	LEU
1	A	64	ARG
1	A	82	ILE
1	A	91	VAL
1	A	98	LEU
1	A	124	HIS
1	A	131	GLU
1	A	153	THR
1	A	168	PRO
1	A	169	THR
1	A	170	PRO
1	A	171	PRO
1	A	174	SER
1	A	175	LEU

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Mol	Chain	Res	Type
1	A	187	LEU
1	A	197	PRO
1	A	203	LEU
1	A	209	PRO
1	A	261	LEU
1	A	268	LEU
1	A	276	LEU
1	A	317	LEU
1	A	325	ILE
1	A	326	PRO
1	A	333	ASN
1	A	360	ARG
1	A	367	PRO
1	A	376	LEU
1	A	451	SER
1	A	464	LEU
1	A	471	LEU
1	A	480	LEU
1	A	511	LEU
1	A	564	LEU
1	A	594	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	138	ASN
1	A	147	ASN
1	A	154	GLN
1	A	333	ASN
1	A	341	ASN
1	A	364	ASN
1	A	423	GLN
1	A	460	GLN
1	A	497	ASN
1	A	520	GLN
1	A	556	ASN
1	A	568	GLN
1	A	570	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	1.66	1 (12%)	7,12,14	2.11	2 (28%)

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	-

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	198	SEP	P-O1P	3.52	1.61	1.50

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	198	SEP	OG-P-O1P	4.43	118.41	106.44
1	A	198	SEP	O2P-P-OG	-3.25	98.19	106.67

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 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	A	198	SEP	2	0

5.5 Carbohydrates [i](#)

10 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	0.83	0	17,19,21	1.22	2 (11%)
2	NAG	B	2	2	14,14,15	1.04	1 (7%)	17,19,21	1.05	1 (5%)
2	MAN	B	3	2	11,11,12	0.70	0	15,15,17	1.14	1 (6%)
3	NAG	C	1	1,3	14,14,15	0.70	0	17,19,21	0.72	0
3	NAG	C	2	3	14,14,15	0.68	0	17,19,21	1.49	3 (17%)
2	NAG	D	1	2,1	14,14,15	0.63	0	17,19,21	0.91	0
2	NAG	D	2	2	14,14,15	1.16	2 (14%)	17,19,21	1.56	2 (11%)
2	MAN	D	3	2	11,11,12	0.88	1 (9%)	15,15,17	0.97	1 (6%)
3	NAG	E	1	1,3	14,14,15	1.10	2 (14%)	17,19,21	1.34	1 (5%)
3	NAG	E	2	3	14,14,15	0.74	1 (7%)	17,19,21	1.75	2 (11%)

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	-	0/6/23/26	0/1/1/1
2	NAG	B	2	2	-	2/6/23/26	0/1/1/1
2	MAN	B	3	2	-	2/2/19/22	0/1/1/1
3	NAG	C	1	1,3	-	0/6/23/26	0/1/1/1
3	NAG	C	2	3	-	1/6/23/26	0/1/1/1
2	NAG	D	1	2,1	-	0/6/23/26	0/1/1/1
2	NAG	D	2	2	-	0/6/23/26	0/1/1/1
2	MAN	D	3	2	-	0/2/19/22	1/1/1/1
3	NAG	E	1	1,3	-	0/6/23/26	0/1/1/1
3	NAG	E	2	3	-	2/6/23/26	0/1/1/1

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	2	NAG	O5-C5	3.00	1.49	1.43
2	B	2	NAG	C1-C2	2.67	1.56	1.52
3	E	1	NAG	C3-C2	2.41	1.57	1.52
3	E	2	NAG	C1-C2	2.27	1.55	1.52
3	E	1	NAG	C1-C2	2.22	1.55	1.52
2	D	3	MAN	O5-C5	2.15	1.47	1.43
2	D	2	NAG	O5-C1	2.01	1.47	1.43

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	E	2	NAG	C4-C3-C2	5.80	119.51	111.02
2	D	2	NAG	C1-O5-C5	4.90	118.75	112.19
3	E	1	NAG	C4-C3-C2	4.67	117.86	111.02
3	C	2	NAG	C4-C3-C2	4.20	117.18	111.02
2	B	2	NAG	C4-C3-C2	3.34	115.91	111.02
2	B	3	MAN	C3-C4-C5	3.09	115.83	110.23
3	E	2	NAG	C3-C4-C5	2.86	115.41	110.23
2	D	3	MAN	C1-O5-C5	2.70	115.81	112.19
2	B	1	NAG	C2-N2-C7	-2.34	119.76	122.90
3	C	2	NAG	C3-C4-C5	2.28	114.36	110.23
2	D	2	NAG	C3-C4-C5	2.22	114.27	110.23
2	B	1	NAG	C1-O5-C5	2.10	115.00	112.19
3	C	2	NAG	C2-N2-C7	-2.02	120.19	122.90

There are no chirality outliers.

All (7) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	2	NAG	O5-C5-C6-O6
3	E	2	NAG	C4-C5-C6-O6
2	B	3	MAN	C4-C5-C6-O6
3	E	2	NAG	O5-C5-C6-O6
2	B	2	NAG	C4-C5-C6-O6
2	B	3	MAN	O5-C5-C6-O6
3	C	2	NAG	O5-C5-C6-O6

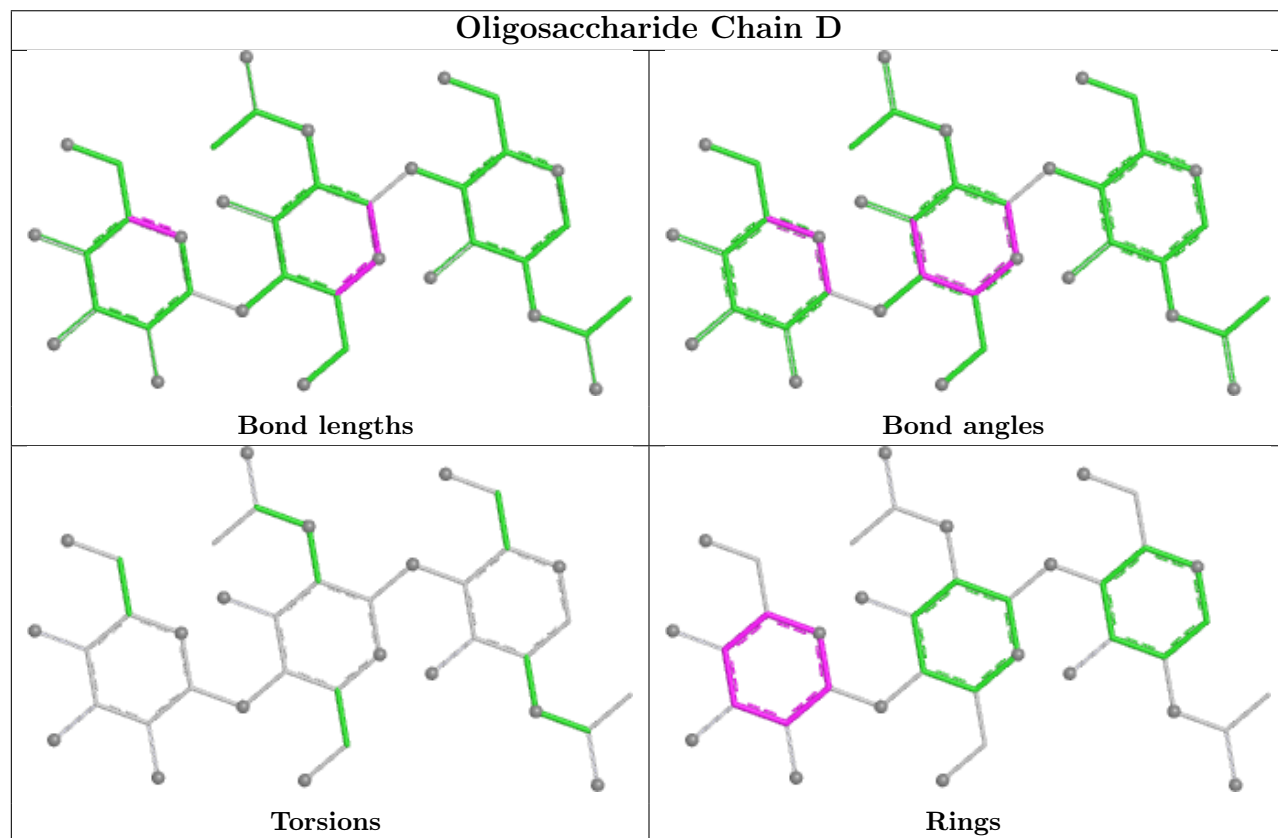
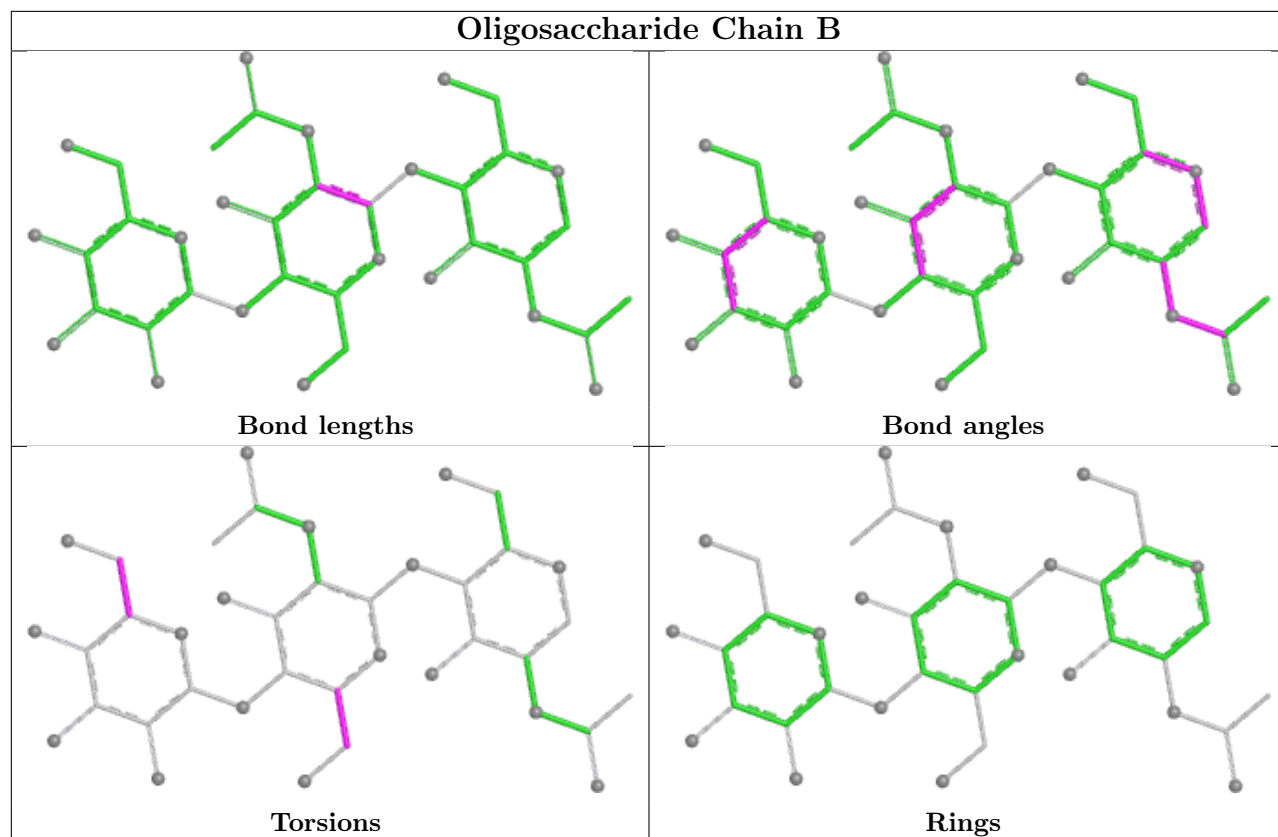
All (1) ring outliers are listed below:

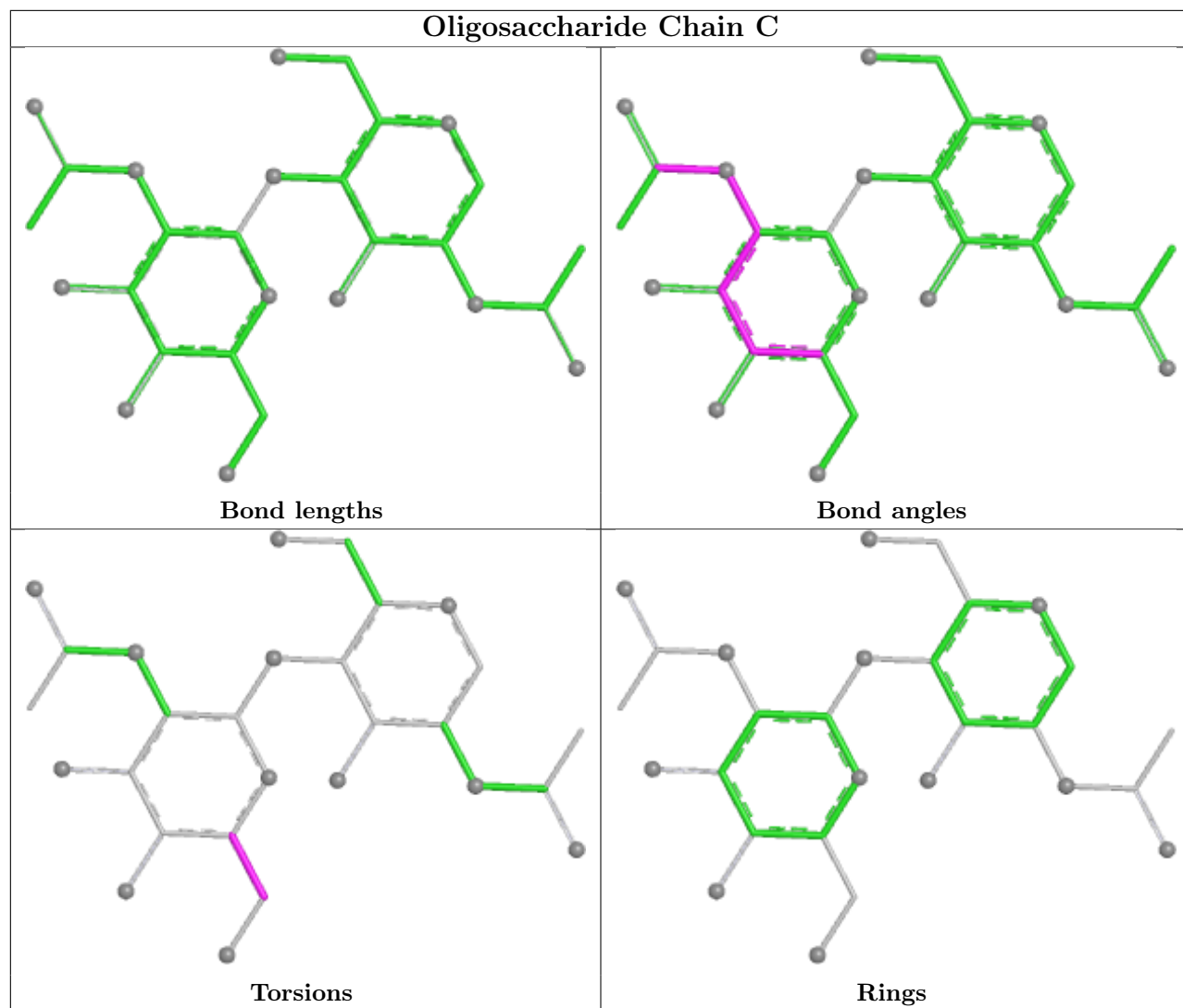
Mol	Chain	Res	Type	Atoms
2	D	3	MAN	C1-C2-C3-C4-C5-O5

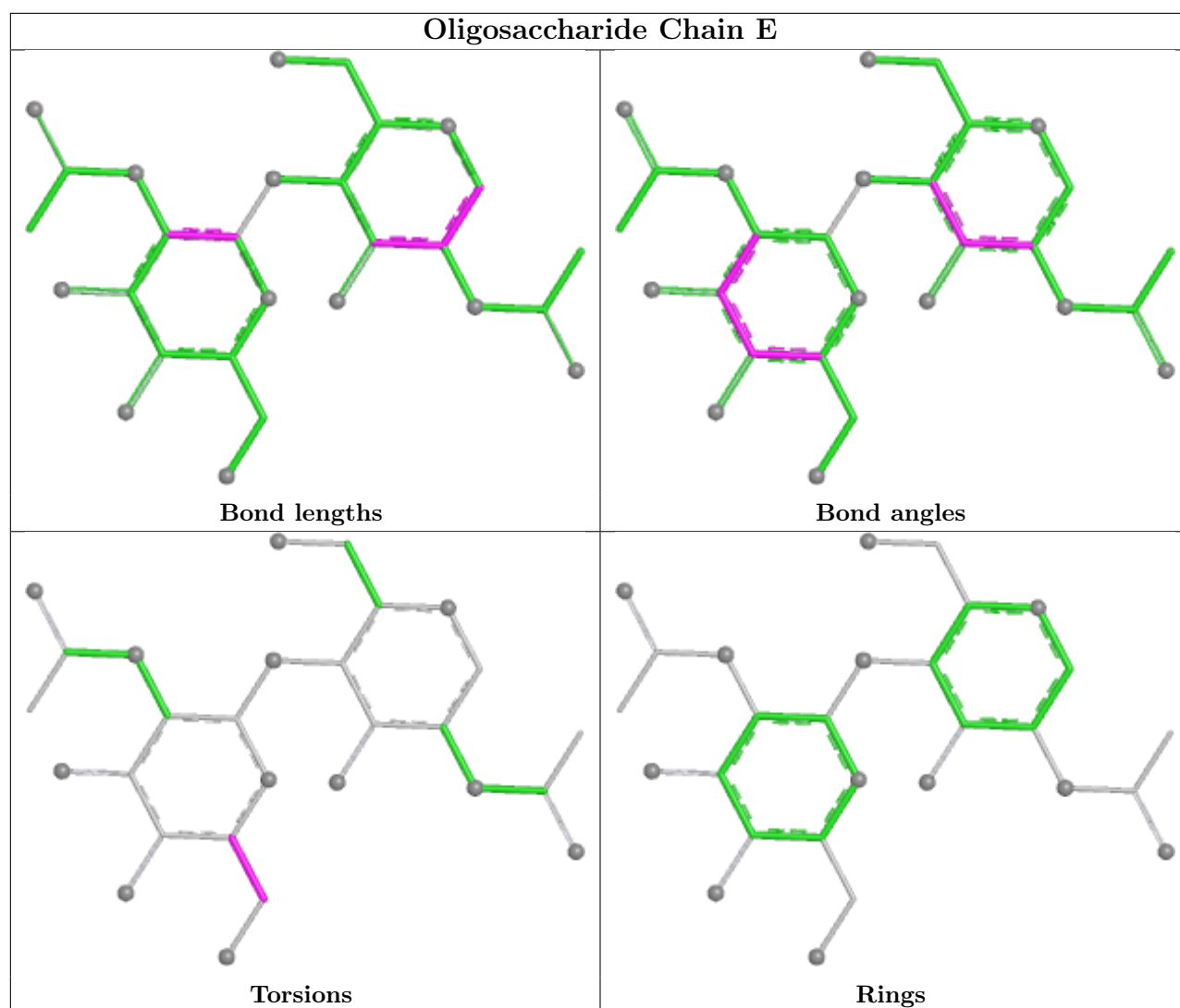
4 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	E	2	NAG	2	0
2	B	2	NAG	1	0
2	D	3	MAN	1	0
2	D	2	NAG	1	0

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







5.6 Ligand geometry [i](#)

Of 9 ligands modelled in this entry, 1 is 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	$\# Z > 2$
6	NO3	A	612	-	1,3,3	3.26	1 (100%)	0,3,3	-	-
5	SCN	A	607	-	1,2,2	1.72	0	0,1,1	-	-
6	NO3	A	609	-	1,3,3	3.21	1 (100%)	0,3,3	-	-

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
6	NO3	A	611	-	1,3,3	3.13	1 (100%)	0,3,3	-	-
8	OSM	A	614	7	1,3,3	0.01	0	0,2,2	-	-
6	NO3	A	608	-	1,3,3	3.02	1 (100%)	0,3,3	-	-
6	NO3	A	610	-	1,3,3	3.12	1 (100%)	0,3,3	-	-
7	HEM	A	613	1,8	50,50,50	2.23	20 (40%)	67,82,82	1.87	12 (17%)

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
7	HEM	A	613	1,8	-	4/14/54/54	-
8	OSM	A	614	7	-	0/0/1/1	-

All (25) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
7	A	613	HEM	CMC-C2C	4.76	1.60	1.50
7	A	613	HEM	CAC-C3C	4.55	1.59	1.47
7	A	613	HEM	FE-NB	4.38	2.08	1.94
7	A	613	HEM	CAB-C3B	4.32	1.58	1.47
7	A	613	HEM	C1B-NB	-4.06	1.33	1.40
7	A	613	HEM	C1A-NA	-3.55	1.32	1.39
7	A	613	HEM	CBD-CAD	3.29	1.63	1.51
6	A	612	NO3	O1-N	3.26	1.40	1.24
6	A	609	NO3	O1-N	3.21	1.40	1.24
6	A	611	NO3	O1-N	3.13	1.39	1.24
7	A	613	HEM	CHA-C1A	-3.12	1.32	1.39
6	A	610	NO3	O1-N	3.12	1.39	1.24
7	A	613	HEM	CHC-C4B	-3.09	1.32	1.39
7	A	613	HEM	CAD-C3D	3.06	1.59	1.51
7	A	613	HEM	C4D-ND	-3.04	1.35	1.40
6	A	608	NO3	O1-N	3.02	1.39	1.24
7	A	613	HEM	C3D-C2D	2.95	1.43	1.36
7	A	613	HEM	C4A-NA	-2.73	1.34	1.39
7	A	613	HEM	C1C-C2C	2.66	1.50	1.45
7	A	613	HEM	C4B-NB	-2.62	1.33	1.38
7	A	613	HEM	O2D-CGD	-2.55	1.22	1.30
7	A	613	HEM	FE-NC	2.53	2.03	1.95
7	A	613	HEM	FE-NA	2.49	2.03	1.95

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
7	A	613	HEM	C2A-C3A	-2.46	1.32	1.38
7	A	613	HEM	O1A-CGA	2.28	1.29	1.22

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	A	613	HEM	CAD-C3D-C4D	7.61	137.95	124.70
7	A	613	HEM	CMD-C2D-C1D	6.51	135.21	125.03
7	A	613	HEM	CAD-C3D-C2D	-4.66	119.14	127.87
7	A	613	HEM	CMD-C2D-C3D	-3.37	117.04	126.15
7	A	613	HEM	C3B-C4B-NB	3.36	111.88	109.47
7	A	613	HEM	O1A-CGA-CBA	-2.90	113.91	123.09
7	A	613	HEM	C4D-C3D-C2D	-2.75	102.89	106.89
7	A	613	HEM	CMC-C2C-C1C	2.46	129.06	124.73
7	A	613	HEM	C2A-C1A-NA	2.42	112.84	110.15
7	A	613	HEM	CHD-C1D-ND	2.34	126.95	124.42
7	A	613	HEM	C4D-ND-C1D	2.25	107.87	105.21
7	A	613	HEM	O2A-CGA-CBA	2.09	120.60	114.00

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
7	A	613	HEM	CAD-CBD-CGD-O1D
7	A	613	HEM	CAD-CBD-CGD-O2D
7	A	613	HEM	CAA-CBA-CGA-O1A
7	A	613	HEM	CAA-CBA-CGA-O2A

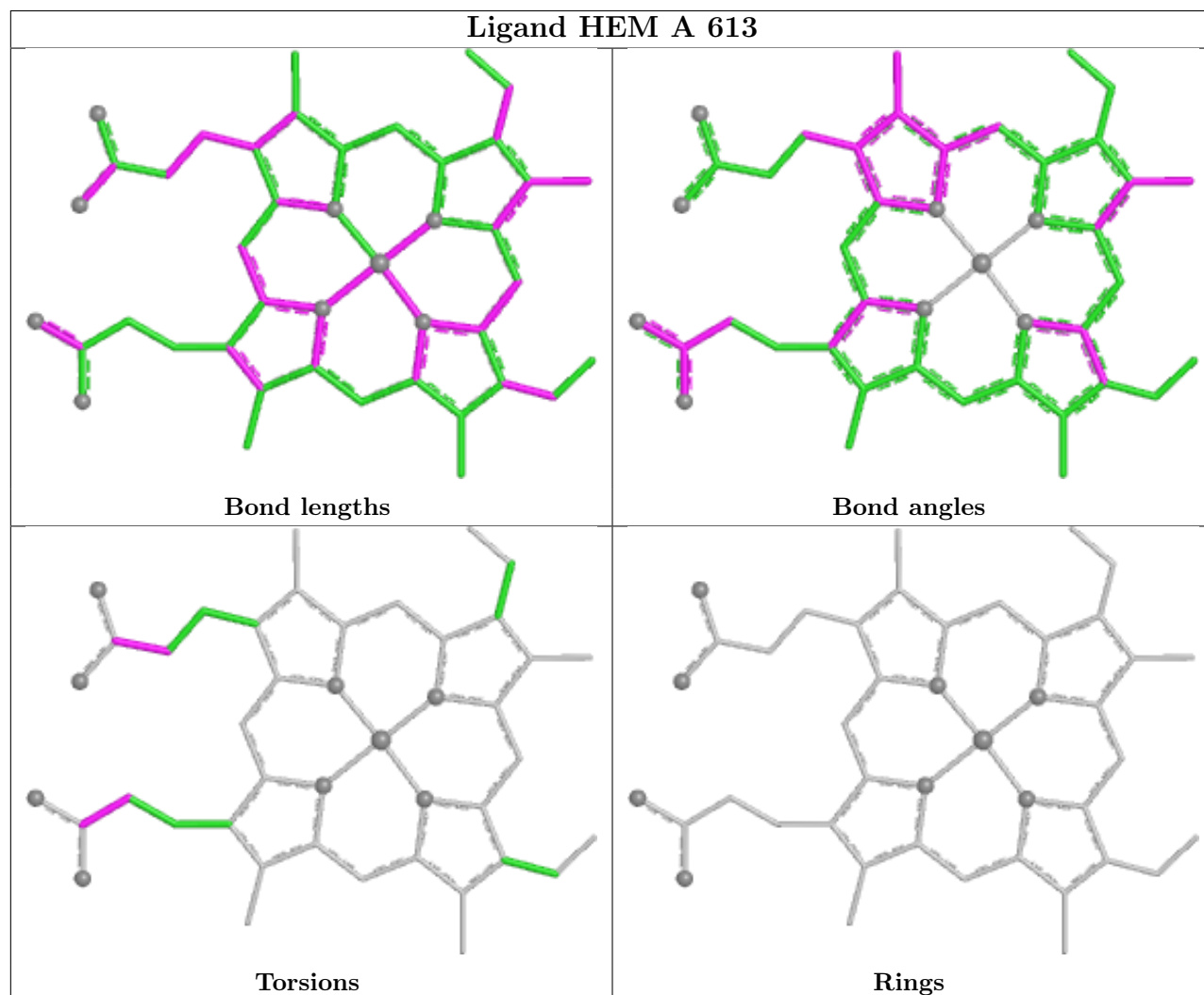
There are no ring outliers.

3 monomers are involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
8	A	614	OSM	3	0
6	A	608	NO3	1	0
7	A	613	HEM	6	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%)	-0.21	27 (4%) 38 40	9, 24, 59, 82	0

All (27) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	7	GLY	6.5
1	A	10	VAL	6.3
1	A	4	VAL	5.5
1	A	2	TRP	5.2
1	A	11	PRO	4.5
1	A	12	LEU	4.3
1	A	167	CYS	4.3
1	A	9	PRO	4.2
1	A	172	TYR	4.2
1	A	174	SER	4.0
1	A	168	PRO	3.9
1	A	8	ALA	3.8
1	A	3	GLU	3.8
1	A	5	GLY	3.8
1	A	121	SER	3.8
1	A	6	CYS	3.6
1	A	170	PRO	3.4
1	A	1	SER	3.3
1	A	13	VAL	3.3
1	A	175	LEU	3.3
1	A	169	THR	3.0
1	A	173	GLN	2.8
1	A	595	ASN	2.7
1	A	171	PRO	2.6
1	A	119	LEU	2.6
1	A	122	ASN	2.1
1	A	64	ARG	2.1

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	0.72	0.17	27,30,51,51	0

6.3 Carbohydrates [i](#)

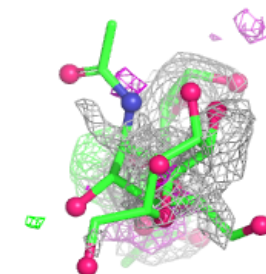
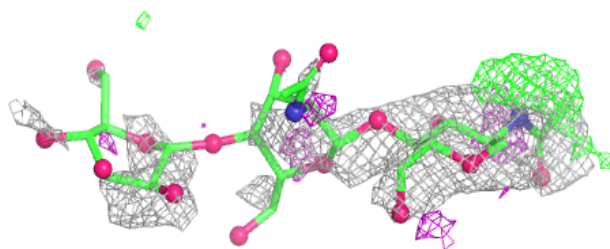
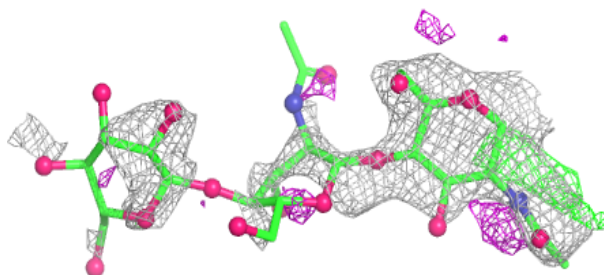
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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	MAN	B	3	11/12	0.21	0.16	92,94,94,95	0
2	NAG	B	2	14/15	0.58	0.17	77,82,84,88	0
2	MAN	D	3	11/12	0.58	0.12	79,82,82,83	0
2	NAG	B	1	14/15	0.63	0.16	51,56,61,69	0
2	NAG	D	2	14/15	0.66	0.14	63,68,71,76	0
3	NAG	E	1	14/15	0.66	0.13	55,62,64,70	0
3	NAG	C	2	14/15	0.71	0.12	60,63,66,68	0
3	NAG	E	2	14/15	0.72	0.13	74,77,78,78	0
3	NAG	C	1	14/15	0.90	0.08	33,44,49,54	0
2	NAG	D	1	14/15	0.90	0.11	42,44,51,56	0

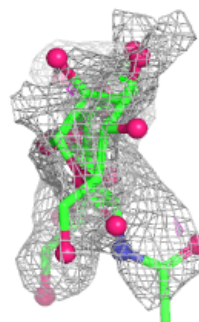
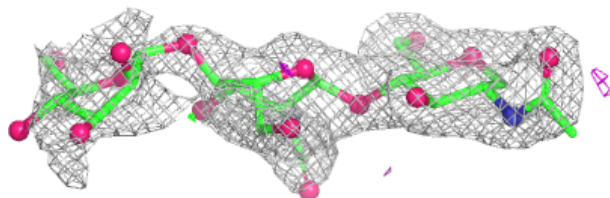
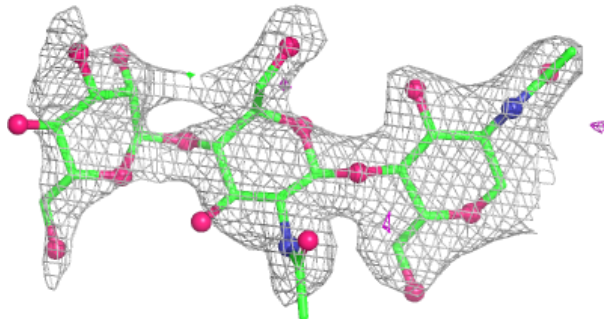
The following is a graphical depiction of the model fit to experimental electron density for oligosaccharide. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around Chain B:

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

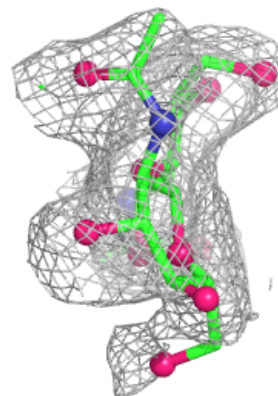
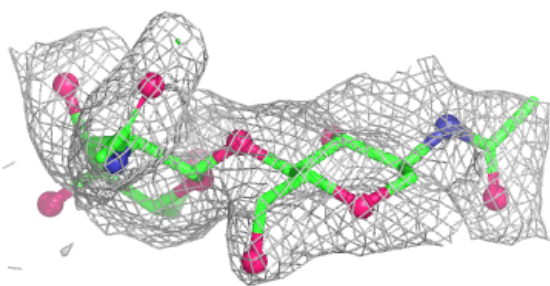
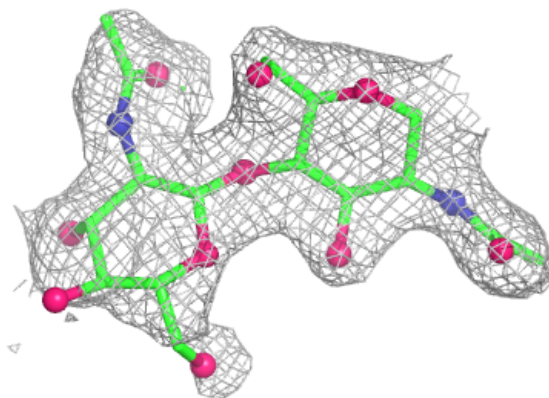
**Electron density around Chain D:**

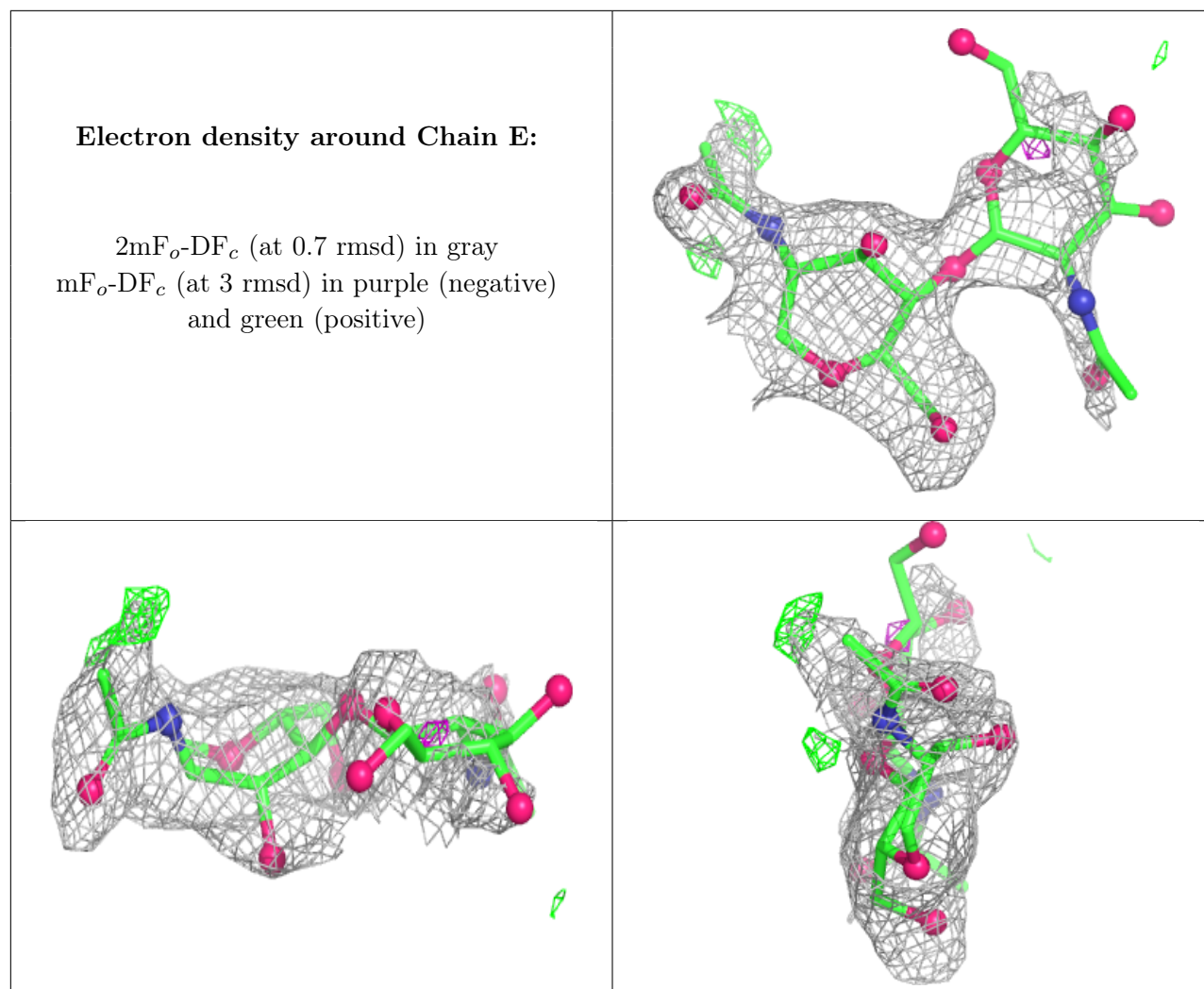
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around Chain C:

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



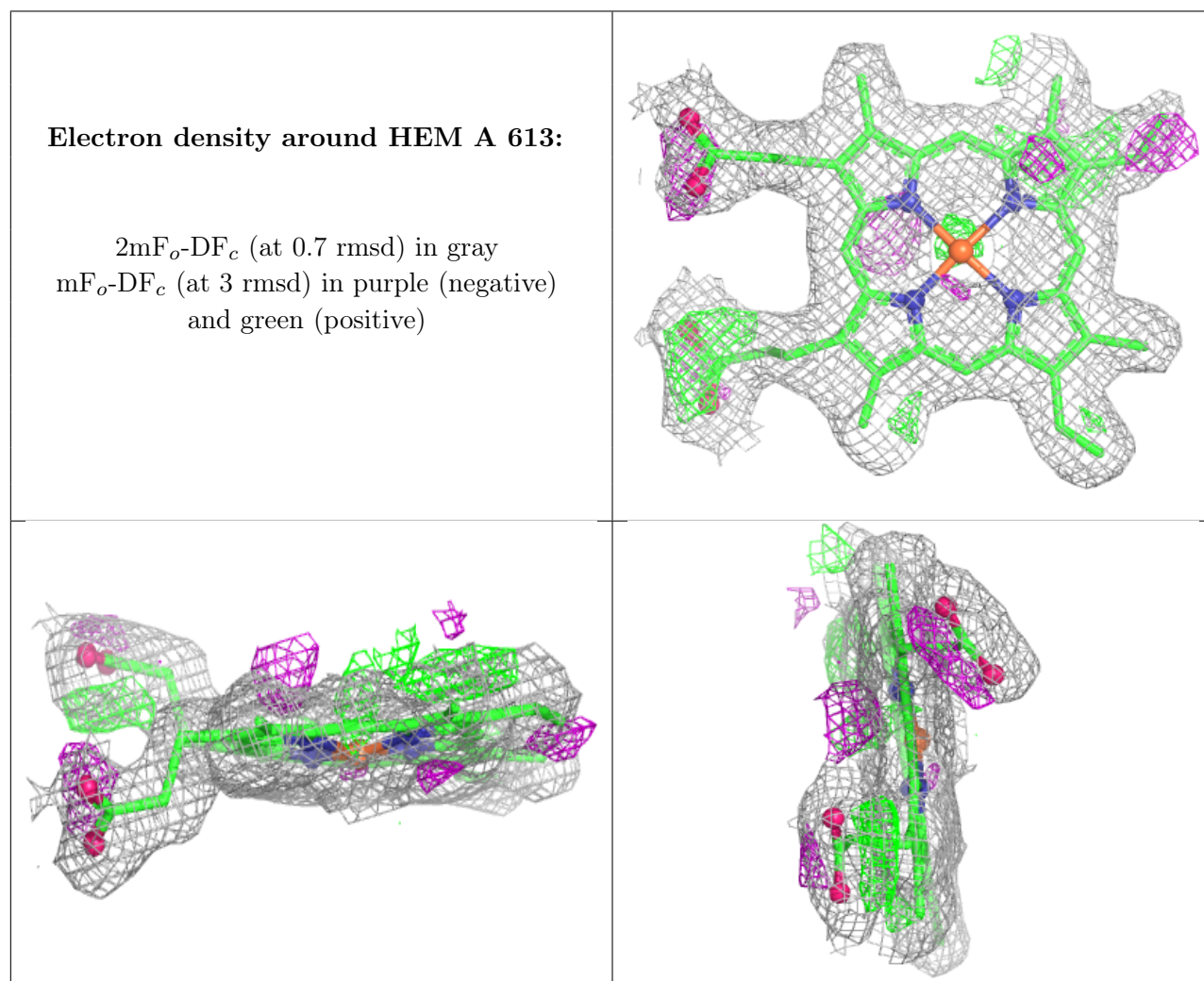


6.4 Ligands ⓘ

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
8	OSM	A	614	4/4	0.70	0.20	40,40,41,48	0
6	NO3	A	612	4/4	0.80	0.22	53,54,54,54	0
6	NO3	A	610	4/4	0.84	0.14	43,44,44,45	0
6	NO3	A	609	4/4	0.88	0.16	59,59,59,59	0
5	SCN	A	607	3/3	0.88	0.13	61,61,62,62	0
6	NO3	A	611	4/4	0.92	0.09	43,44,44,45	0
7	HEM	A	613	43/43	0.95	0.08	8,11,15,17	0
6	NO3	A	608	4/4	0.96	0.09	15,17,20,21	0
4	CA	A	606	1/1	0.98	0.02	12,12,12,12	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.



6.5 Other polymers ⓘ

There are no such residues in this entry.