



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 02:00 PM UTC

PDB ID : 2Z5Z / pdb_00002z5z
Title : Crystal structure of the complex of buffalo Lactoperoxidase with fluoride ion at 3.5Å resolution
Authors : Sheikh, I.A.; Jain, R.; Singh, N.; Sharma, S.; Bhushan, A.; Kaur, P.; Srinivasan, A.; Singh, T.P.
Deposited on : 2007-07-20
Resolution : 3.50 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

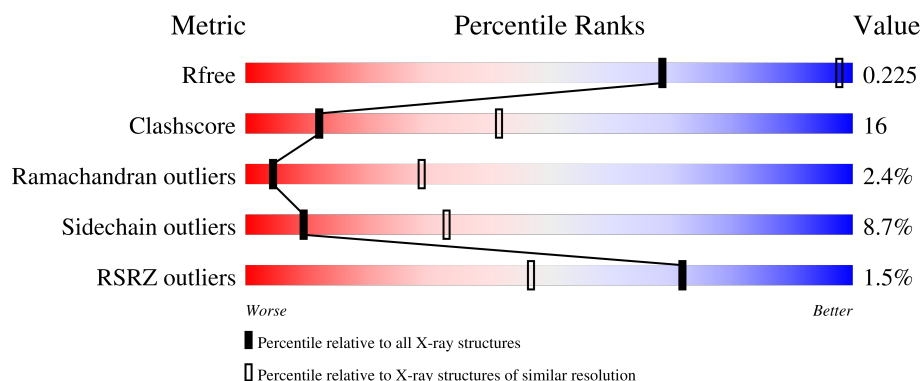
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.50 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.



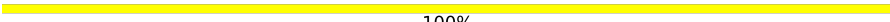
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	1085 (3.54-3.46)
Clashscore	190562	1140 (3.54-3.46)
Ramachandran outliers	187476	1113 (3.54-3.46)
Sidechain outliers	187428	1114 (3.54-3.46)
RSRZ outliers	180081	1084 (3.54-3.46)

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% 65% 28% 6%
2	B	3	33% 67%
2	D	3	33% 33% 33%
3	C	2	50% 50%

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Mol	Chain	Length	Quality of chain
3	E	2	 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
1	SEP	A	198	-	-	X	-
5	IOD	A	607	-	-	X	-
5	IOD	A	611	-	-	X	-
5	IOD	A	613	-	-	X	-
5	IOD	A	615	-	-	X	-
6	F	A	616	-	-	-	X
6	F	A	618	-	-	-	X
8	OSM	A	701	-	-	X	-

2 Entry composition [i](#)

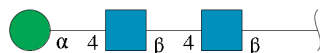
There are 9 unique types of molecules in this entry. The entry contains 4990 atoms, of which 0 are hydrogens and 0 are deuteriums.

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
			4770	3032	845	865	1	27			

- 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 1 1	0	0

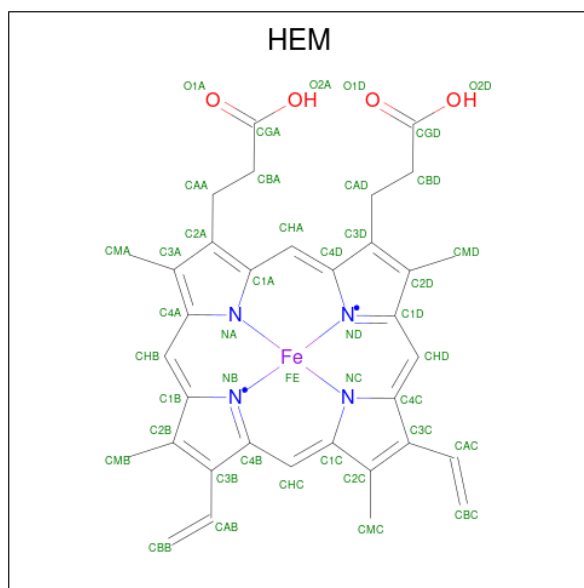
- Molecule 5 is IODIDE ION (CCD ID: IOD) (formula: I).

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

- Molecule 6 is FLUORIDE ION (CCD ID: F) (formula: F).

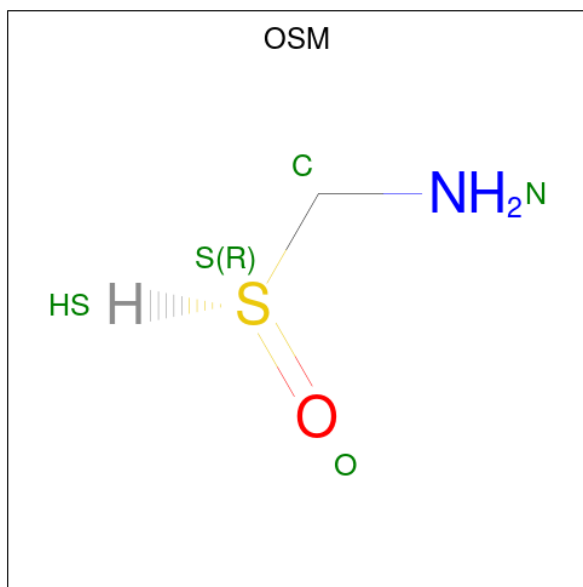
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	A	3	Total F 3 3	0	0

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



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

- Molecule 8 is 1-(OXIDOSULFANYL)METHANAMINE (CCD ID: OSM) (formula: CH_5NOS).



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

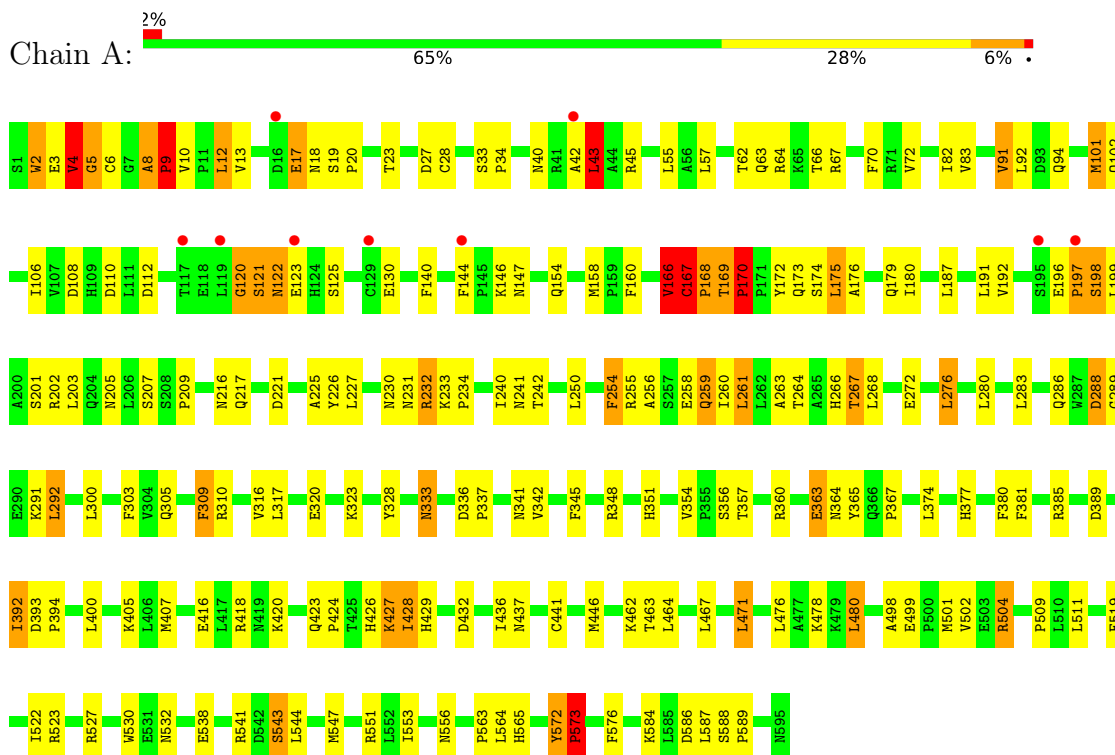
- Molecule 9 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
9	A	26	Total	O	0	0
			26	26		

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:  100%

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	54.13Å 80.74Å 77.70Å 90.00° 102.94° 90.00°	Depositor
Resolution (Å)	75.81 – 3.50 75.73 – 3.50	Depositor EDS
% Data completeness (in resolution range)	83.2 (75.81-3.50) 83.2 (75.73-3.50)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.13	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.85 (at 3.48Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.221 , 0.235 0.211 , 0.225	Depositor DCC
R_{free} test set	324 reflections (3.87%)	wwPDB-VP
Wilson B-factor (Å ²)	45.7	Xtriage
Anisotropy	0.600	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 77.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.89	EDS
Total number of atoms	4990	wwPDB-VP
Average B, all atoms (Å ²)	27.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.34% 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: IOD, MAN, HEM, F, SEP, OSM, NAG, CA

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	0.61	1/4886 (0.0%)	1.48	69/6627 (1.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	573	PRO	N-CD	5.93	1.56	1.47

All (69) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	572	TYR	CA-C-N	44.62	175.62	119.84
1	A	572	TYR	C-N-CA	44.62	175.62	119.84
1	A	573	PRO	CA-N-CD	-18.98	85.42	112.00
1	A	573	PRO	N-CA-CB	16.42	120.49	103.25
1	A	5	GLY	N-CA-C	12.07	126.50	112.50
1	A	573	PRO	N-CD-CG	11.71	120.77	103.20
1	A	121	SER	N-CA-C	-10.73	100.87	113.21
1	A	173	GLN	OE1-CD-NE2	-10.26	112.34	122.60
1	A	392	ILE	N-CA-C	9.51	121.66	111.58
1	A	232	ARG	N-CA-C	9.41	124.57	110.64
1	A	427	LYS	N-CA-C	9.22	124.97	112.90
1	A	316	VAL	N-CA-C	-9.17	103.92	111.81
1	A	572	TYR	C-N-CD	-8.32	90.88	125.00
1	A	289	GLY	N-CA-C	-8.29	102.89	112.50
1	A	429	HIS	N-CA-C	-7.93	97.84	109.11
1	A	205	ASN	OD1-CG-ND2	-7.87	114.73	122.60
1	A	82	ILE	CB-CA-C	-7.72	104.11	111.44
1	A	207	SER	N-CA-C	-7.30	104.17	113.23
1	A	91	VAL	N-CA-C	7.18	122.48	112.50
1	A	205	ASN	CB-CG-ND2	7.02	126.93	116.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	173	GLN	CG-CD-NE2	6.98	126.87	116.40
1	A	82	ILE	N-CA-C	6.95	117.54	111.56
1	A	556	ASN	N-CA-C	6.93	122.72	112.94
1	A	225	ALA	N-CA-C	6.88	119.84	110.55
1	A	112	ASP	N-CA-C	6.88	119.89	109.23
1	A	172	TYR	N-CA-C	6.46	120.06	109.72
1	A	167	CYS	N-CA-C	-6.43	95.59	109.81
1	A	288	ASP	N-CA-C	6.42	119.00	110.53
1	A	34	PRO	N-CA-C	6.41	123.11	113.81
1	A	43	LEU	N-CA-C	-6.39	100.97	110.23
1	A	418	ARG	N-CA-C	6.39	121.15	113.23
1	A	256	ALA	N-CA-C	6.36	119.02	111.33
1	A	67	ARG	N-CA-C	-6.22	98.27	108.41
1	A	167	CYS	CA-C-N	5.99	127.33	119.84
1	A	167	CYS	C-N-CA	5.99	127.33	119.84
1	A	323	LYS	N-CA-C	5.88	118.64	111.82
1	A	170	PRO	N-CA-C	5.83	117.82	110.70
1	A	55	LEU	N-CA-C	5.82	120.43	112.04
1	A	94	GLN	N-CA-C	5.75	119.77	112.87
1	A	260	ILE	N-CA-C	5.75	116.40	110.36
1	A	363	GLU	N-CA-C	5.67	119.01	111.75
1	A	572	TYR	O-C-N	5.66	134.94	121.93
1	A	420	LYS	N-CA-C	5.62	119.68	112.26
1	A	205	ASN	N-CA-C	-5.62	99.37	108.41
1	A	110	ASP	N-CA-C	-5.61	105.55	112.90
1	A	92	LEU	N-CA-C	5.56	118.51	110.28
1	A	309	PHE	N-CA-C	5.54	118.24	111.82
1	A	530	TRP	N-CA-C	5.53	118.83	111.75
1	A	173	GLN	N-CA-C	5.46	122.42	110.80
1	A	345	PHE	N-CA-C	-5.42	107.21	113.88
1	A	573	PRO	CA-C-N	5.40	131.85	121.54
1	A	573	PRO	C-N-CA	5.40	131.85	121.54
1	A	33	SER	CA-C-N	5.40	125.73	119.47
1	A	33	SER	C-N-CA	5.40	125.73	119.47
1	A	192	VAL	CB-CA-C	-5.34	105.41	111.55
1	A	28	CYS	N-CA-C	5.30	119.58	113.38
1	A	405	LYS	N-CA-C	-5.29	102.64	110.48
1	A	154	GLN	N-CA-C	5.29	120.21	113.55
1	A	356	SER	N-CA-C	5.28	119.72	113.28
1	A	565	HIS	N-CA-C	-5.23	99.26	108.20
1	A	140	PHE	CA-C-N	5.20	125.14	119.78
1	A	140	PHE	C-N-CA	5.20	125.14	119.78

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	254	PHE	N-CA-C	5.14	118.65	112.38
1	A	4	VAL	N-CA-C	5.07	119.89	109.34
1	A	428	ILE	N-CA-C	5.04	119.83	109.34
1	A	197	PRO	CB-CA-C	5.04	119.67	110.10
1	A	336	ASP	CA-C-N	5.02	125.09	119.87
1	A	336	ASP	C-N-CA	5.02	125.09	119.87
1	A	337	PRO	N-CA-C	5.00	120.18	114.03

There are no chirality outliers.

There are no planarity outliers.

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	4770	0	4674	150	0
2	B	39	0	34	0	0
2	D	39	0	34	1	0
3	C	28	0	25	3	0
3	E	28	0	25	0	0
4	A	1	0	0	0	0
5	A	9	0	0	17	0
6	A	3	0	0	0	0
7	A	43	0	30	13	0
8	A	4	0	5	2	0
9	A	26	0	0	8	0
All	All	4990	0	4827	153	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 16.

All (153) 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:198:SEP:CA	1:A:198:SEP:CB	1.76	1.58
1:A:108:ASP:OD2	7:A:619:HEM:HMD1	1.24	1.34

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:167:CYS:HB3	1:A:168:PRO:HD2	1.23	1.18
1:A:198:SEP:N	1:A:198:SEP:OG	1.82	1.13
1:A:198:SEP:CB	1:A:198:SEP:N	2.12	1.12
1:A:446:MET:HE1	5:A:607:IOD:I	2.26	1.06
1:A:258:GLU:OE2	7:A:619:HEM:HMB1	1.58	1.03
1:A:198:SEP:CB	1:A:198:SEP:C	2.35	1.03
1:A:258:GLU:OE2	7:A:619:HEM:CMB	2.08	1.01
1:A:217:GLN:HE21	3:C:1:NAG:H83	1.27	1.00
1:A:108:ASP:OD2	7:A:619:HEM:CMD	2.11	0.99
1:A:351:HIS:HD1	1:A:437:ASN:HD21	1.08	0.98
1:A:196:GLU:HB3	1:A:198:SEP:O3P	1.64	0.95
1:A:216:ASN:HA	5:A:611:IOD:I	2.36	0.94
5:A:615:IOD:I	9:A:724:HOH:O	2.54	0.94
1:A:8:ALA:HB1	1:A:9:PRO:HD2	1.47	0.92
1:A:341:ASN:HB3	5:A:607:IOD:I	2.40	0.91
1:A:169:THR:HB	1:A:170:PRO:HD2	1.52	0.89
1:A:8:ALA:HB1	1:A:9:PRO:CD	2.02	0.88
1:A:462:LYS:N	5:A:615:IOD:I	2.77	0.86
1:A:8:ALA:CB	1:A:9:PRO:HD2	2.05	0.84
1:A:263:ALA:O	1:A:267:THR:HG22	1.78	0.83
1:A:167:CYS:HB3	1:A:168:PRO:CD	2.06	0.83
1:A:120:GLY:C	1:A:122:ASN:H	1.84	0.81
1:A:12:LEU:HD23	9:A:712:HOH:O	1.81	0.81
1:A:169:THR:HB	1:A:170:PRO:CD	2.12	0.79
1:A:198:SEP:CB	1:A:199:LEU:N	2.47	0.78
1:A:63:GLN:HG2	9:A:720:HOH:O	1.86	0.76
1:A:8:ALA:O	1:A:10:VAL:HG22	1.87	0.75
1:A:2:TRP:HB2	1:A:4:VAL:HG22	1.69	0.75
1:A:463:THR:N	5:A:615:IOD:I	2.89	0.74
1:A:463:THR:OG1	5:A:615:IOD:I	2.72	0.73
1:A:333:ASN:C	1:A:333:ASN:HD22	1.96	0.73
1:A:101:MET:HE2	1:A:102:GLN:HA	1.71	0.73
1:A:2:TRP:CB	1:A:4:VAL:HG22	2.20	0.72
1:A:198:SEP:HB3	1:A:199:LEU:H	1.56	0.71
1:A:42:ALA:HB2	1:A:166:VAL:HG21	1.73	0.69
1:A:504:ARG:NH2	9:A:705:HOH:O	2.25	0.69
1:A:108:ASP:CG	7:A:619:HEM:HMD1	2.16	0.68
1:A:328:TYR:CD1	1:A:523:ARG:HD3	2.29	0.68
1:A:198:SEP:HB3	1:A:199:LEU:N	2.08	0.66
1:A:198:SEP:CB	1:A:198:SEP:H	1.91	0.66
1:A:354:VAL:HG11	7:A:619:HEM:CBB	2.27	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:288:ASP:HB2	1:A:291:LYS:CB	2.28	0.64
1:A:258:GLU:OE2	7:A:619:HEM:HMB2	1.96	0.64
1:A:101:MET:HE2	1:A:102:GLN:CA	2.28	0.64
1:A:221:ASP:HB2	1:A:226:TYR:CZ	2.32	0.64
1:A:216:ASN:CG	5:A:611:IOD:I	3.11	0.63
1:A:202:ARG:HD3	1:A:250:LEU:HD21	1.81	0.63
1:A:144:PHE:HE2	1:A:158:MET:HE2	1.64	0.62
1:A:563:PRO:HD3	1:A:576:PHE:CE2	2.34	0.62
1:A:169:THR:O	1:A:170:PRO:O	2.18	0.62
1:A:354:VAL:HG11	7:A:619:HEM:HBB1	1.82	0.61
1:A:351:HIS:HD1	1:A:437:ASN:ND2	1.88	0.61
1:A:276:LEU:HD12	1:A:587:LEU:HD11	1.83	0.61
1:A:328:TYR:HD1	1:A:523:ARG:HD3	1.66	0.60
1:A:393:ASP:HB2	1:A:394:PRO:HD3	1.83	0.60
1:A:62:THR:HG22	1:A:64:ARG:HG2	1.83	0.60
1:A:146:LYS:O	1:A:147:ASN:HB2	1.99	0.60
1:A:106:ILE:HG23	1:A:191:LEU:HD11	1.83	0.59
1:A:167:CYS:CB	1:A:168:PRO:HD2	2.13	0.59
1:A:538:GLU:HG3	1:A:541:ARG:NH2	2.19	0.58
1:A:169:THR:CB	1:A:170:PRO:CD	2.82	0.57
1:A:197:PRO:O	1:A:198:SEP:C	2.31	0.57
1:A:232:ARG:O	1:A:232:ARG:HG3	2.04	0.57
1:A:385:ARG:O	1:A:389:ASP:HB3	2.05	0.57
1:A:175:LEU:HD23	1:A:176:ALA:H	1.71	0.56
1:A:333:ASN:C	1:A:333:ASN:ND2	2.62	0.56
1:A:120:GLY:C	1:A:122:ASN:N	2.55	0.55
1:A:544:LEU:HD23	1:A:547:MET:CE	2.37	0.55
1:A:216:ASN:CA	5:A:611:IOD:I	3.21	0.54
1:A:381:PHE:CZ	1:A:424:PRO:HB3	2.43	0.54
1:A:476:LEU:HD21	1:A:498:ALA:HB1	1.90	0.53
1:A:288:ASP:HB2	1:A:291:LYS:HB2	1.90	0.53
1:A:2:TRP:HB3	1:A:4:VAL:HG22	1.91	0.53
1:A:363:GLU:O	5:A:614:IOD:I	2.97	0.53
1:A:499:GLU:OE1	1:A:509:PRO:HD2	2.09	0.52
1:A:45:ARG:NE	9:A:727:HOH:O	2.41	0.52
1:A:572:TYR:CD2	1:A:573:PRO:HD3	2.44	0.52
1:A:288:ASP:HB2	1:A:291:LYS:HB3	1.92	0.52
1:A:198:SEP:CA	1:A:198:SEP:OG	2.36	0.52
1:A:544:LEU:HD23	1:A:547:MET:HE3	1.93	0.51
1:A:423:GLN:HE21	1:A:423:GLN:HA	1.77	0.50
1:A:432:ASP:OD1	1:A:432:ASP:C	2.55	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:217:GLN:NE2	3:C:1:NAG:H83	2.10	0.50
1:A:527:ARG:O	1:A:532:ASN:ND2	2.44	0.50
1:A:3:GLU:O	1:A:5:GLY:N	2.44	0.50
1:A:63:GLN:CG	9:A:720:HOH:O	2.53	0.50
1:A:288:ASP:O	1:A:292:LEU:HD23	2.12	0.50
1:A:241:ASN:ND2	2:D:1:NAG:C7	2.74	0.50
1:A:259:GLN:HE22	1:A:261:LEU:HB2	1.76	0.49
1:A:272:GLU:OE1	1:A:272:GLU:HA	2.11	0.49
1:A:400:LEU:HD13	1:A:563:PRO:HD2	1.93	0.49
1:A:8:ALA:CB	1:A:9:PRO:CD	2.67	0.49
1:A:187:LEU:HD13	1:A:305:GLN:HA	1.95	0.49
1:A:467:LEU:HG	1:A:471:LEU:HD22	1.94	0.48
1:A:8:ALA:O	1:A:9:PRO:C	2.56	0.48
1:A:101:MET:CE	1:A:102:GLN:HA	2.41	0.48
1:A:166:VAL:C	1:A:167:CYS:O	2.52	0.48
1:A:309:PHE:HB3	5:A:613:IOD:I	2.83	0.48
1:A:360:ARG:HA	1:A:394:PRO:O	2.14	0.48
1:A:160:PHE:HE2	1:A:436:ILE:HG23	1.80	0.47
1:A:377:HIS:HA	1:A:380:PHE:CE2	2.50	0.47
1:A:9:PRO:O	1:A:10:VAL:HG13	2.15	0.47
1:A:198:SEP:O	1:A:199:LEU:C	2.57	0.47
1:A:258:GLU:HG3	8:A:701:OSM:HN1	1.79	0.46
1:A:320:GLU:HG3	1:A:502:VAL:HG11	1.96	0.46
1:A:342:VAL:HB	5:A:607:IOD:I	2.85	0.46
1:A:551:ARG:HD3	1:A:584:LYS:HA	1.97	0.46
1:A:407:MET:HB3	1:A:501:MET:HE3	1.97	0.46
1:A:23:THR:OG1	1:A:27:ASP:O	2.33	0.46
1:A:2:TRP:O	1:A:3:GLU:HB2	2.16	0.45
1:A:300:LEU:O	1:A:303:PHE:HB3	2.17	0.45
1:A:254:PHE:HE2	9:A:706:HOH:O	1.99	0.45
1:A:43:LEU:HD23	1:A:179:GLN:HB2	1.98	0.45
1:A:217:GLN:HE21	3:C:1:NAG:C8	2.12	0.45
1:A:10:VAL:HG12	1:A:40:ASN:O	2.17	0.45
1:A:167:CYS:O	1:A:168:PRO:O	2.35	0.45
1:A:255:ARG:HG2	8:A:701:OSM:H2	1.99	0.44
1:A:446:MET:HA	1:A:446:MET:HE2	1.99	0.44
1:A:259:GLN:HE21	1:A:259:GLN:HB2	1.54	0.44
1:A:123:GLU:HG3	1:A:125:SER:H	1.82	0.44
1:A:230:ASN:OD1	1:A:231:ASN:N	2.50	0.44
5:A:609:IOD:I	9:A:705:HOH:O	2.92	0.44
1:A:66:THR:HB	1:A:70:PHE:N	2.32	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:351:HIS:CD2	7:A:619:HEM:NC	2.86	0.44
1:A:480:LEU:HD12	1:A:480:LEU:HA	1.85	0.43
1:A:543:SER:OG	1:A:586:ASP:O	2.36	0.43
7:A:619:HEM:HBC2	7:A:619:HEM:HMC2	2.01	0.43
1:A:101:MET:HE2	1:A:102:GLN:N	2.34	0.43
1:A:227:LEU:HD11	1:A:266:HIS:HB3	2.00	0.43
1:A:342:VAL:N	5:A:607:IOD:I	3.22	0.43
1:A:588:SER:N	1:A:589:PRO:CD	2.82	0.42
1:A:108:ASP:CG	7:A:619:HEM:CMD	2.86	0.42
1:A:310:ARG:HB2	5:A:613:IOD:I	2.89	0.42
1:A:130:GLU:OE2	1:A:426:HIS:HB3	2.20	0.42
1:A:424:PRO:O	5:A:612:IOD:I	3.07	0.42
1:A:519:PHE:HA	1:A:522:ILE:HG12	2.02	0.42
1:A:364:ASN:O	1:A:365:TYR:HB2	2.20	0.41
1:A:407:MET:CB	1:A:501:MET:HE3	2.50	0.41
1:A:19:SER:HA	1:A:20:PRO:HD3	1.85	0.41
1:A:348:ARG:HG2	7:A:619:HEM:C2D	2.55	0.41
1:A:166:VAL:HG13	1:A:180:ILE:HG12	2.03	0.41
1:A:233:LYS:HA	1:A:234:PRO:C	2.45	0.41
1:A:377:HIS:HB3	1:A:416:GLU:OE1	2.21	0.41
1:A:63:GLN:HG2	1:A:63:GLN:H	1.58	0.41
1:A:146:LYS:O	1:A:147:ASN:CB	2.69	0.40
1:A:462:LYS:HB2	5:A:615:IOD:I	2.92	0.40
1:A:354:VAL:HG21	7:A:619:HEM:CBB	2.52	0.40
1:A:357:THR:HA	1:A:374:LEU:O	2.21	0.40
1:A:221:ASP:HB2	1:A:226:TYR:CE2	2.55	0.40
1:A:264:THR:HG23	1:A:392:ILE:HD12	2.02	0.40
1:A:400:LEU:HD11	1:A:553:ILE:HD13	2.02	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%)	538 (91%)	40 (7%)	14 (2%)	4	29

All (14) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	4	VAL
1	A	6	CYS
1	A	8	ALA
1	A	9	PRO
1	A	17	GLU
1	A	122	ASN
1	A	168	PRO
1	A	169	THR
1	A	170	PRO
1	A	174	SER
1	A	573	PRO
1	A	167	CYS
1	A	120	GLY
1	A	166	VAL

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	516/516 (100%)	471 (91%)	45 (9%)	9	33

All (45) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	2	TRP
1	A	4	VAL
1	A	9	PRO
1	A	12	LEU
1	A	13	VAL
1	A	17	GLU
1	A	18	ASN

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Mol	Chain	Res	Type
1	A	43	LEU
1	A	57	LEU
1	A	72	VAL
1	A	83	VAL
1	A	91	VAL
1	A	101	MET
1	A	121	SER
1	A	166	VAL
1	A	175	LEU
1	A	201	SER
1	A	203	LEU
1	A	209	PRO
1	A	240	ILE
1	A	242	THR
1	A	259	GLN
1	A	261	LEU
1	A	267	THR
1	A	268	LEU
1	A	276	LEU
1	A	280	LEU
1	A	283	LEU
1	A	286	GLN
1	A	292	LEU
1	A	317	LEU
1	A	333	ASN
1	A	367	PRO
1	A	427	LYS
1	A	428	ILE
1	A	441	CYS
1	A	464	LEU
1	A	471	LEU
1	A	478	LYS
1	A	480	LEU
1	A	504	ARG
1	A	511	LEU
1	A	543	SER
1	A	564	LEU
1	A	573	PRO

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	173	GLN
1	A	217	GLN
1	A	259	GLN
1	A	266	HIS
1	A	333	ASN
1	A	364	ASN
1	A	423	GLN
1	A	429	HIS
1	A	460	GLN
1	A	468	GLN
1	A	497	ASN
1	A	520	GLN
1	A	568	GLN

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	4.55	3 (37%)	7,12,14	6.81	4 (57%)

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	-	1/6/8/10	-

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	198	SEP	CB-CA	8.70	1.76	1.52
1	A	198	SEP	O-C	7.63	1.48	1.20
1	A	198	SEP	CA-N	-4.88	1.34	1.48

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	198	SEP	OG-CB-CA	-15.88	92.70	108.14
1	A	198	SEP	OG-P-O1P	6.28	123.43	106.44
1	A	198	SEP	O2P-P-OG	4.12	117.42	106.67
1	A	198	SEP	O3P-P-O1P	-3.33	97.84	110.83

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	198	SEP	N-CA-CB-OG

There are no ring outliers.

1 monomer is involved in 12 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	A	198	SEP	12	0

5.5 Carbohydrates

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	0.65	0	17,19,21	1.75	6 (35%)
2	NAG	B	2	2	14,14,15	0.92	0	17,19,21	1.28	2 (11%)
2	MAN	B	3	2	11,11,12	0.88	0	15,15,17	0.57	0
3	NAG	C	1	3,1	14,14,15	1.59	3 (21%)	17,19,21	2.29	3 (17%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	NAG	C	2	3	14,14,15	0.84	1 (7%)	17,19,21	2.49	5 (29%)
2	NAG	D	1	2,1	14,14,15	0.62	0	17,19,21	1.05	1 (5%)
2	NAG	D	2	2	14,14,15	1.03	1 (7%)	17,19,21	2.41	4 (23%)
2	MAN	D	3	2	11,11,12	0.84	0	15,15,17	0.39	0
3	NAG	E	1	3,1	14,14,15	0.70	0	17,19,21	1.58	3 (17%)
3	NAG	E	2	3	14,14,15	0.57	0	17,19,21	1.01	1 (5%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	NAG	B	1	2,1	-	0/6/23/26	0/1/1/1
2	NAG	B	2	2	-	0/6/23/26	0/1/1/1
2	MAN	B	3	2	-	1/2/19/22	0/1/1/1
3	NAG	C	1	3,1	-	2/6/23/26	0/1/1/1
3	NAG	C	2	3	-	3/6/23/26	0/1/1/1
2	NAG	D	1	2,1	-	4/6/23/26	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	0/1/1/1
3	NAG	E	1	3,1	-	4/6/23/26	0/1/1/1
3	NAG	E	2	3	-	2/6/23/26	0/1/1/1

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	1	NAG	O5-C5	4.14	1.51	1.43
2	D	2	NAG	O5-C5	2.90	1.49	1.43
3	C	1	NAG	O5-C1	2.58	1.48	1.43
3	C	2	NAG	O5-C5	2.17	1.47	1.43
3	C	1	NAG	C2-N2	-2.10	1.42	1.46

All (25) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	2	NAG	C4-C3-C2	-7.73	99.68	111.02
3	C	2	NAG	C1-O5-C5	7.25	121.91	112.19
3	C	1	NAG	C1-O5-C5	5.90	120.10	112.19
3	C	1	NAG	C2-N2-C7	-5.54	115.48	122.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	E	1	NAG	C4-C3-C2	4.55	117.69	111.02
3	C	2	NAG	O5-C1-C2	4.33	117.99	111.29
3	C	2	NAG	C2-N2-C7	-3.69	117.96	122.90
2	D	2	NAG	C2-N2-C7	-3.24	118.56	122.90
2	B	2	NAG	C4-C3-C2	3.19	115.69	111.02
2	B	1	NAG	C2-N2-C7	-3.13	118.71	122.90
2	D	2	NAG	C1-O5-C5	3.05	116.28	112.19
2	B	2	NAG	C2-N2-C7	-3.01	118.86	122.90
3	E	1	NAG	C3-C4-C5	2.96	115.60	110.23
3	C	2	NAG	C3-C4-C5	2.77	115.25	110.23
2	B	1	NAG	C4-C3-C2	2.73	115.02	111.02
3	C	1	NAG	C3-C4-C5	2.69	115.11	110.23
2	B	1	NAG	C1-O5-C5	2.67	115.77	112.19
2	D	2	NAG	C1-C2-N2	2.63	114.58	110.43
2	B	1	NAG	O5-C5-C6	-2.63	102.54	107.66
3	C	2	NAG	O5-C5-C4	2.54	117.01	110.83
3	E	2	NAG	C2-N2-C7	-2.50	119.55	122.90
2	B	1	NAG	C6-C5-C4	-2.48	106.94	113.02
3	E	1	NAG	C2-N2-C7	-2.39	119.70	122.90
2	D	1	NAG	C3-C4-C5	2.30	114.40	110.23
2	B	1	NAG	C3-C4-C5	2.21	114.23	110.23

There are no chirality outliers.

All (16) torsion outliers are listed below:

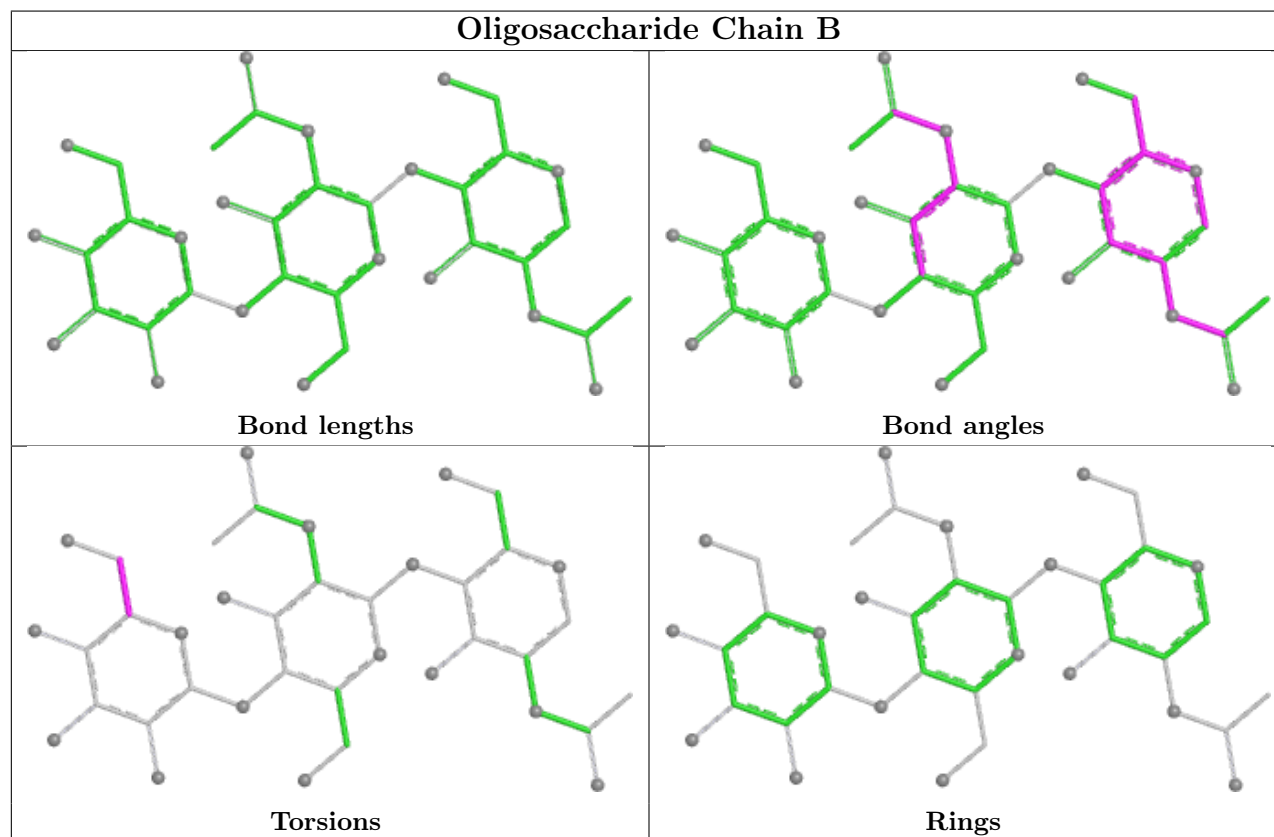
Mol	Chain	Res	Type	Atoms
3	C	2	NAG	C8-C7-N2-C2
3	C	2	NAG	O7-C7-N2-C2
3	E	1	NAG	C8-C7-N2-C2
3	E	1	NAG	O7-C7-N2-C2
2	D	1	NAG	C8-C7-N2-C2
2	D	1	NAG	O7-C7-N2-C2
3	E	1	NAG	C4-C5-C6-O6
3	E	2	NAG	C4-C5-C6-O6
3	E	1	NAG	O5-C5-C6-O6
3	C	2	NAG	O5-C5-C6-O6
3	E	2	NAG	O5-C5-C6-O6
2	B	3	MAN	O5-C5-C6-O6
3	C	1	NAG	O5-C5-C6-O6
2	D	1	NAG	O5-C5-C6-O6
3	C	1	NAG	C8-C7-N2-C2
2	D	1	NAG	C1-C2-N2-C7

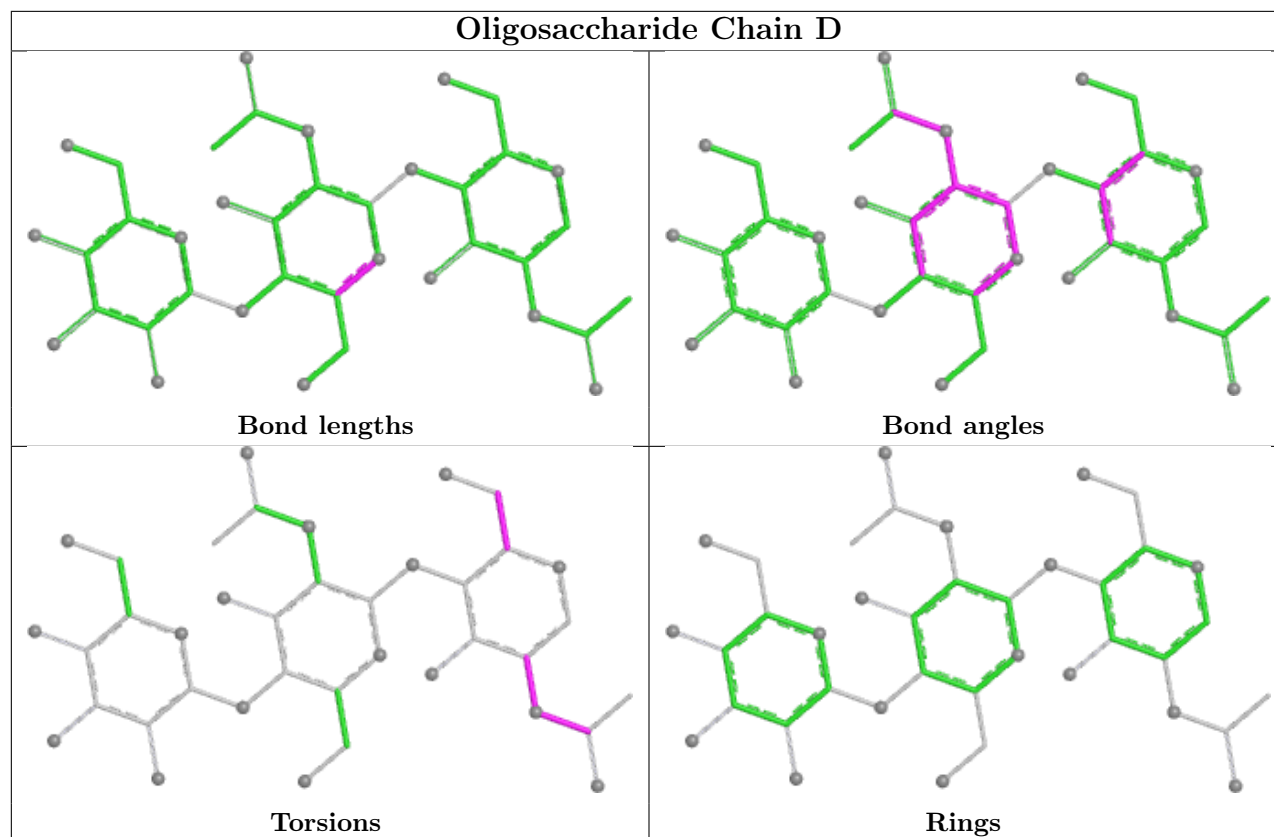
There are no ring outliers.

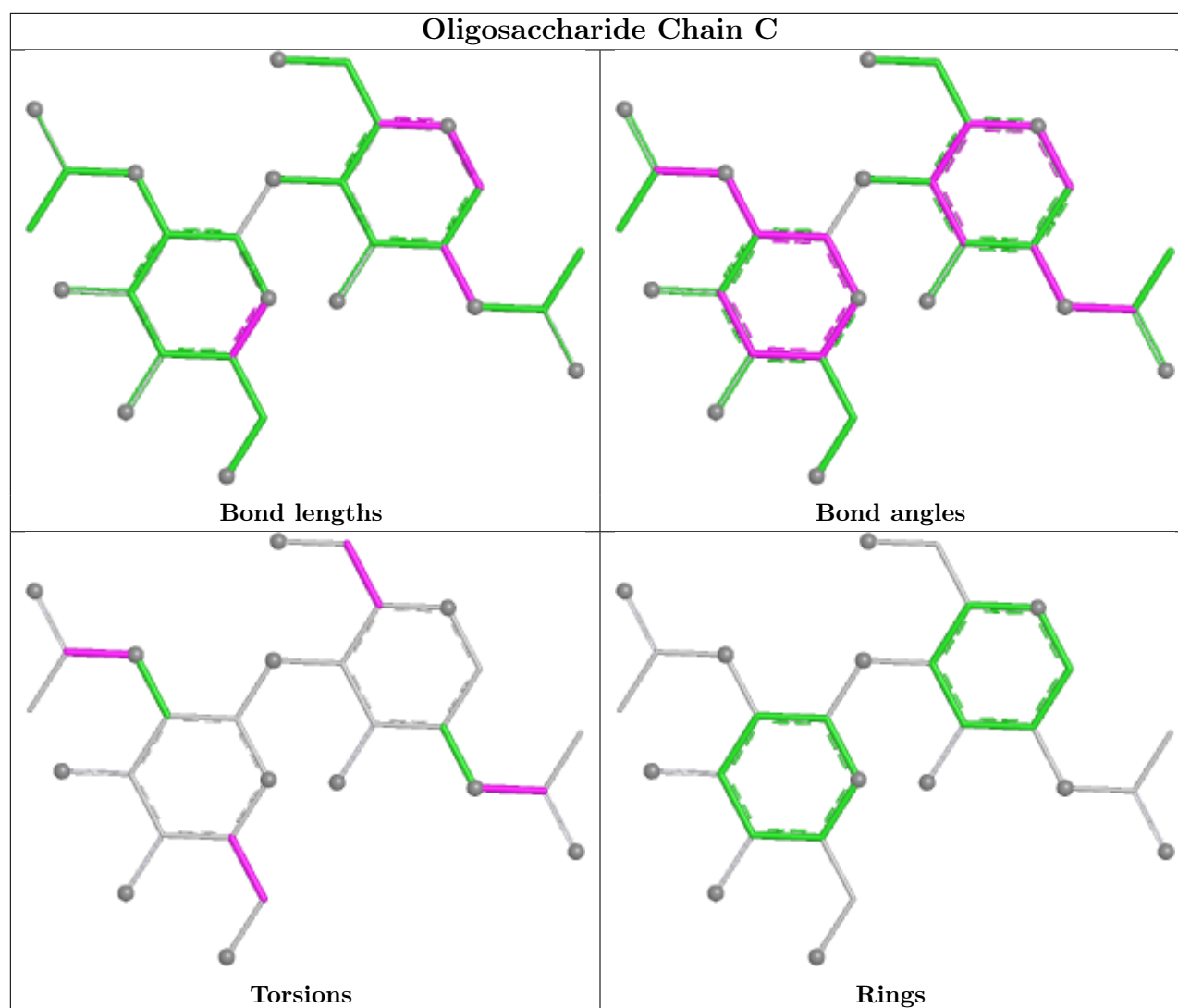
2 monomers are involved in 4 short contacts:

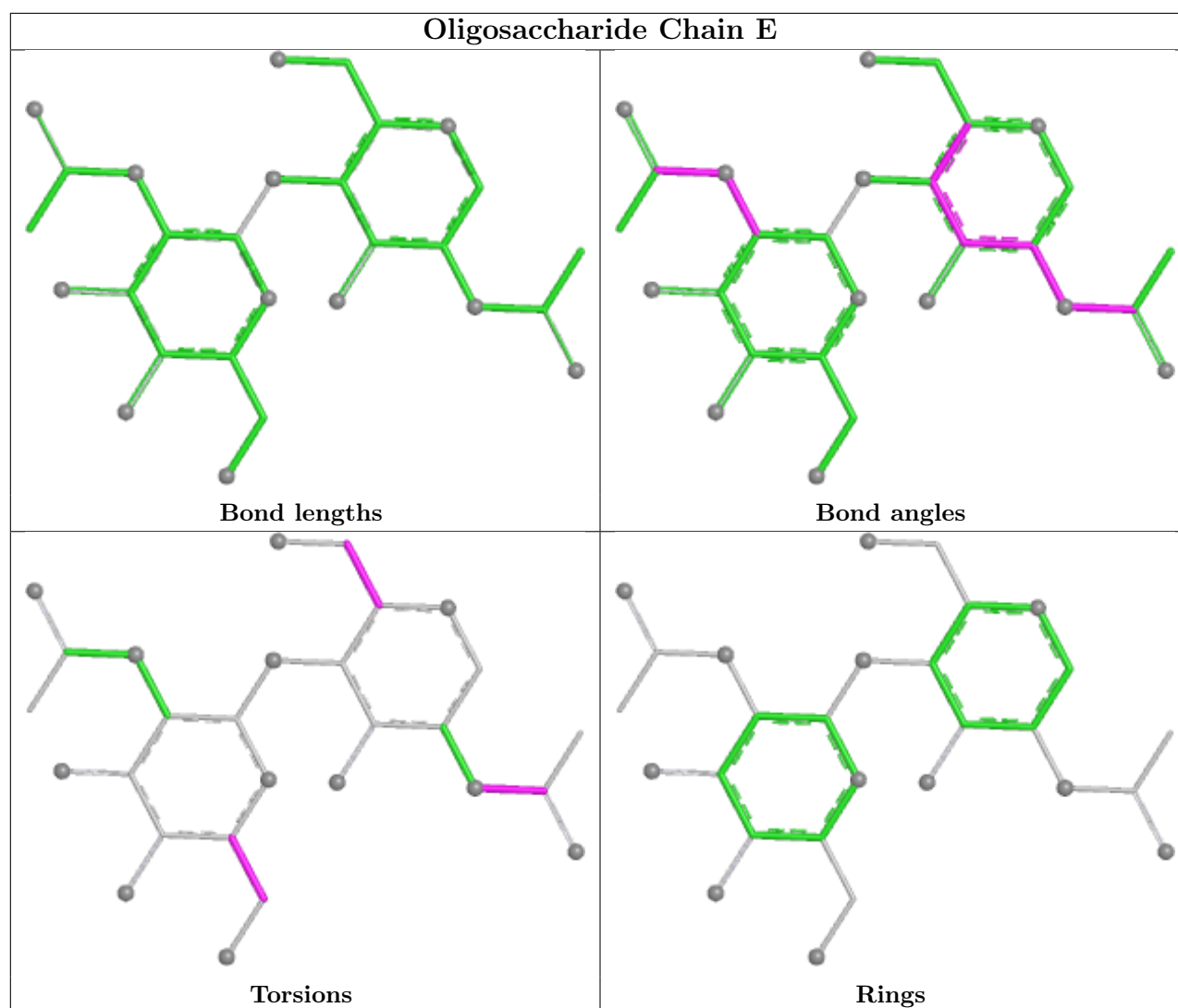
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	C	1	NAG	3	0
2	D	1	NAG	1	0

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









5.6 Ligand geometry [i](#)

Of 15 ligands modelled in this entry, 13 are monoatomic - leaving 2 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	$\# Z > 2$
7	HEM	A	619	1	50,50,50	2.20	15 (30%)	67,82,82	1.26	8 (11%)
8	OSM	A	701	-	1,3,3	0.38	0	0,2,2	-	-

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	619	1	-	6/14/54/54	-
8	OSM	A	701	-	-	0/0/1/1	-

All (15) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
7	A	619	HEM	C3D-C2D	7.24	1.52	1.36
7	A	619	HEM	FE-NA	6.45	2.16	1.95
7	A	619	HEM	O1D-CGD	3.73	1.34	1.22
7	A	619	HEM	CAC-C3C	3.30	1.56	1.47
7	A	619	HEM	CMC-C2C	3.09	1.57	1.50
7	A	619	HEM	FE-ND	-2.75	1.86	1.94
7	A	619	HEM	CHA-C1A	-2.69	1.33	1.39
7	A	619	HEM	CAB-C3B	2.53	1.54	1.47
7	A	619	HEM	C3B-C2B	-2.47	1.32	1.37
7	A	619	HEM	C3C-C4C	-2.24	1.42	1.46
7	A	619	HEM	C3B-C4B	-2.22	1.40	1.44
7	A	619	HEM	CHB-C4A	-2.03	1.34	1.39
7	A	619	HEM	C4A-NA	-2.02	1.35	1.39
7	A	619	HEM	O1A-CGA	2.00	1.28	1.22
7	A	619	HEM	O2D-CGD	-2.00	1.24	1.30

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	A	619	HEM	CHD-C1D-ND	3.80	128.51	124.42
7	A	619	HEM	CAD-C3D-C4D	3.47	130.75	124.70
7	A	619	HEM	C4D-C3D-C2D	-2.75	102.89	106.89
7	A	619	HEM	C4D-ND-C1D	2.71	108.41	105.21
7	A	619	HEM	CMD-C2D-C1D	2.59	129.09	125.03
7	A	619	HEM	CHC-C4B-NB	2.40	127.00	124.42
7	A	619	HEM	C3D-C4D-ND	2.21	112.60	110.17
7	A	619	HEM	CHC-C1C-NC	-2.14	122.13	124.45

There are no chirality outliers.

All (6) torsion outliers are listed below:

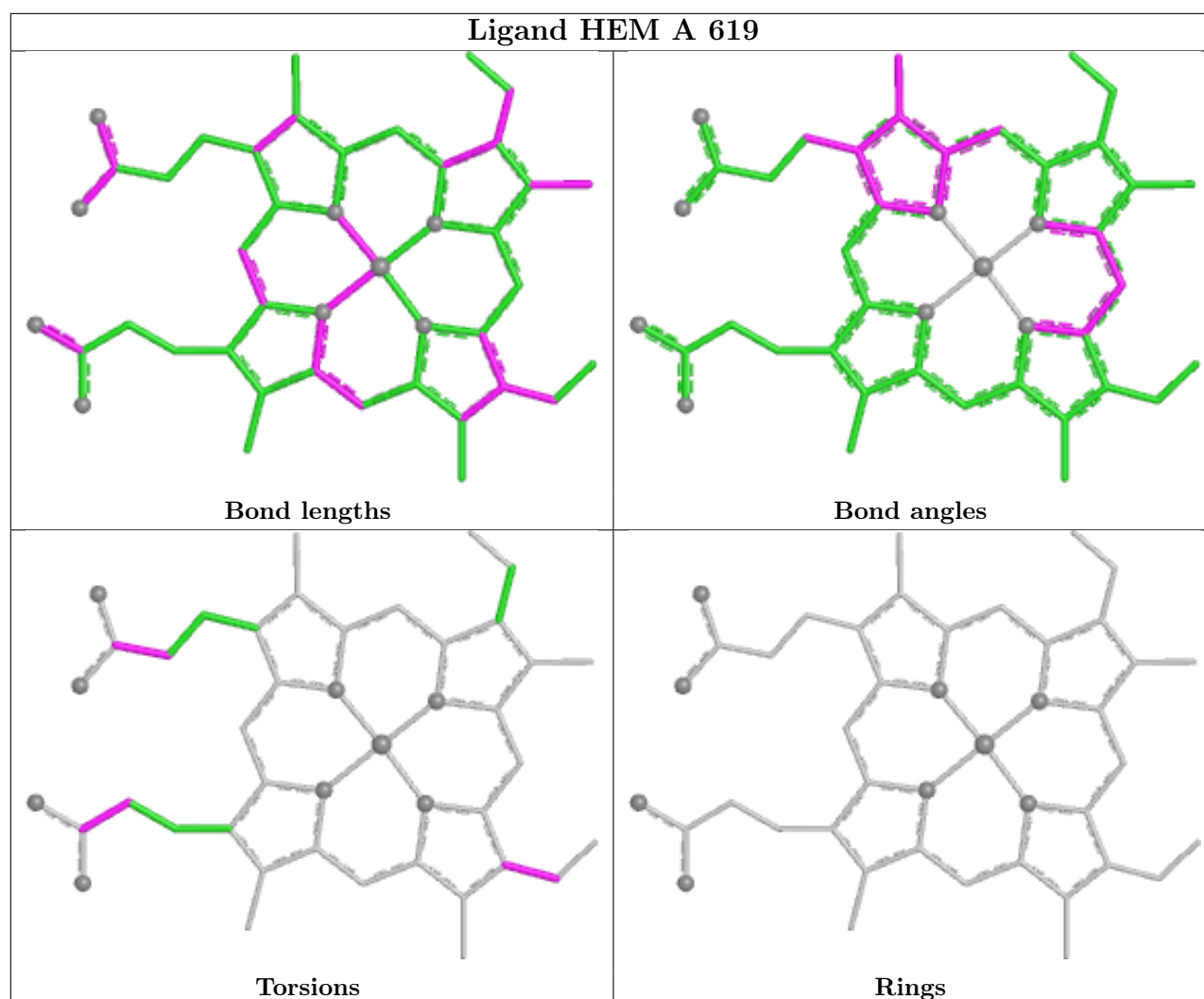
Mol	Chain	Res	Type	Atoms
7	A	619	HEM	C2B-C3B-CAB-CBB
7	A	619	HEM	C4B-C3B-CAB-CBB
7	A	619	HEM	CAD-CBD-CGD-O1D
7	A	619	HEM	CAA-CBA-CGA-O1A
7	A	619	HEM	CAD-CBD-CGD-O2D
7	A	619	HEM	CAA-CBA-CGA-O2A

There are no ring outliers.

2 monomers are involved in 15 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
7	A	619	HEM	13	0
8	A	701	OSM	2	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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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%)	0.17	9 (1%)	72 47	2, 20, 73, 126	0

All (9) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	144	PHE	3.2
1	A	16	ASP	2.6
1	A	197	PRO	2.6
1	A	123	GLU	2.5
1	A	129	CYS	2.5
1	A	42	ALA	2.3
1	A	117	THR	2.1
1	A	119	LEU	2.1
1	A	195	SER	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.

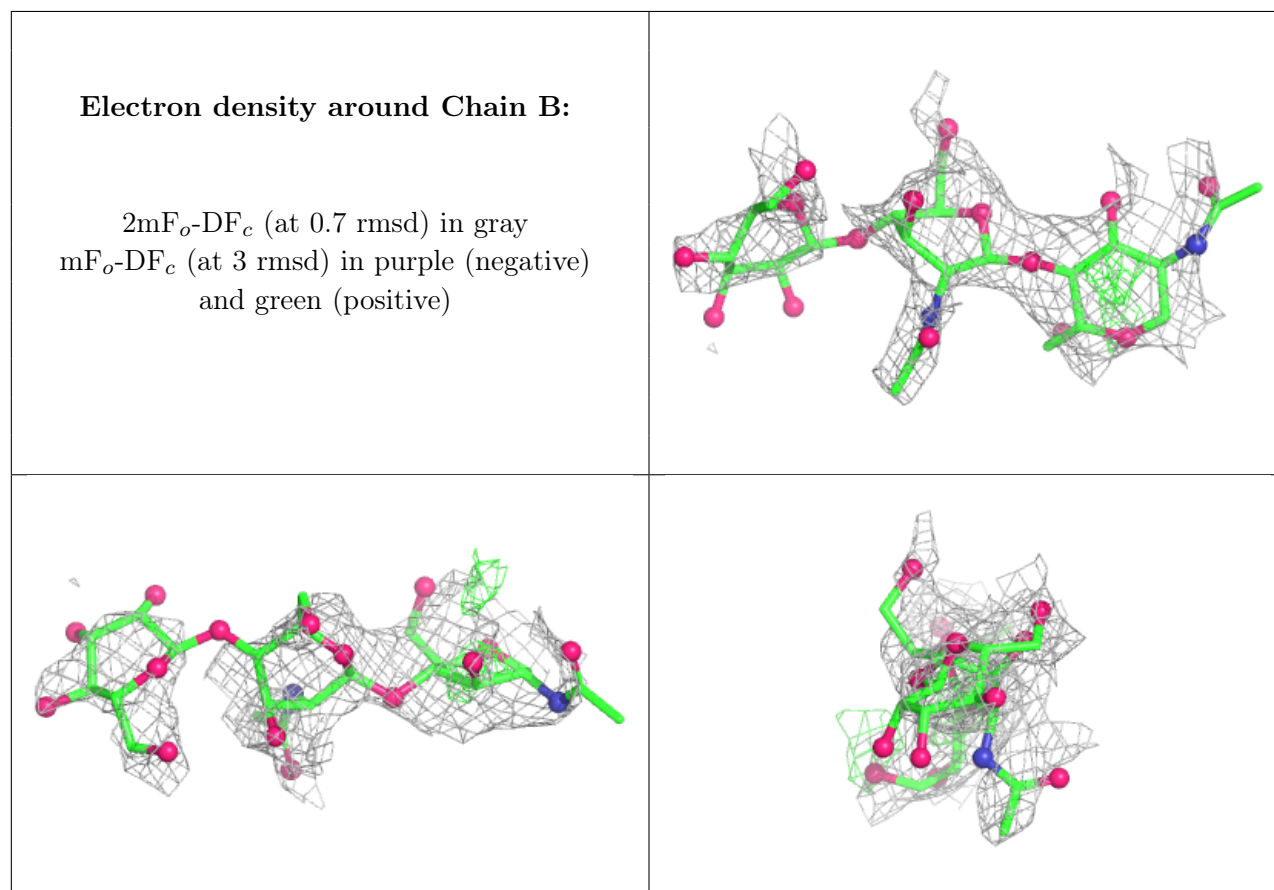
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
1	SEP	A	198	10/11	0.84	0.16	43,55,57,61	0

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.

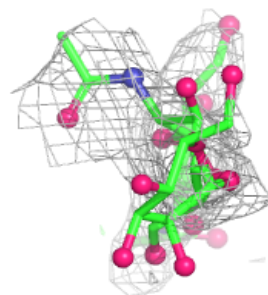
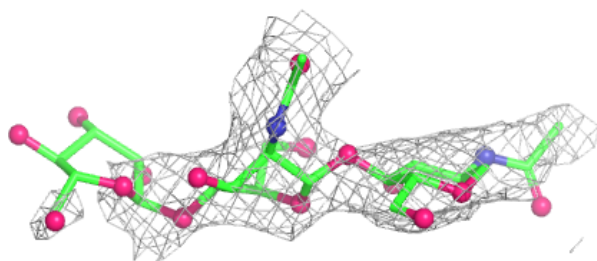
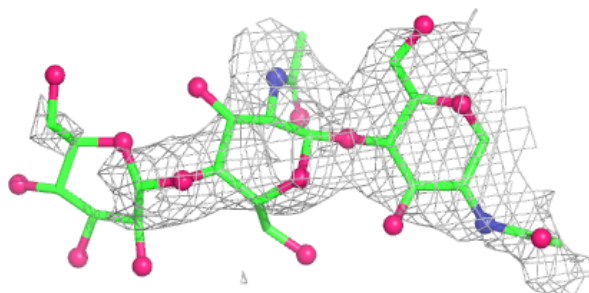
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	MAN	B	3	11/12	0.45	0.21	30,32,32,32	11
2	MAN	D	3	11/12	0.68	0.21	65,66,67,68	11
3	NAG	E	1	14/15	0.74	0.23	55,59,61,61	14
3	NAG	C	2	14/15	0.75	0.16	16,17,18,18	14
2	NAG	B	2	14/15	0.78	0.12	15,16,18,18	14
3	NAG	C	1	14/15	0.78	0.20	3,7,9,9	14
3	NAG	E	2	14/15	0.78	0.25	39,40,41,41	14
2	NAG	D	2	14/15	0.83	0.12	4,6,8,9	14
2	NAG	B	1	14/15	0.89	0.18	2,2,2,2	14
2	NAG	D	1	14/15	0.90	0.16	2,2,2,2	14

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.

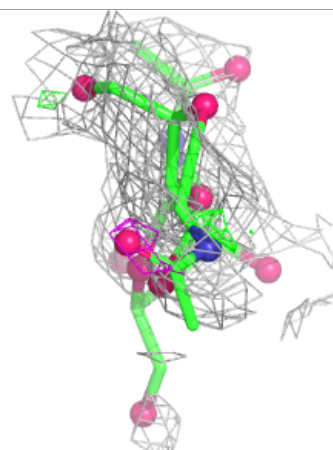
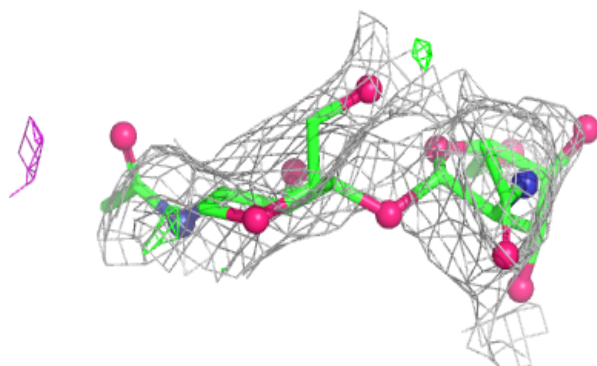
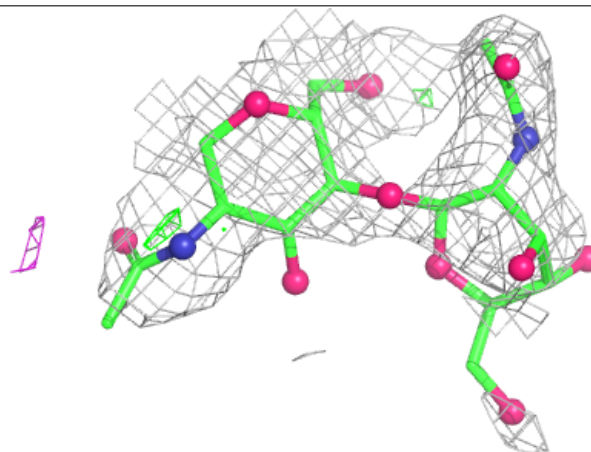


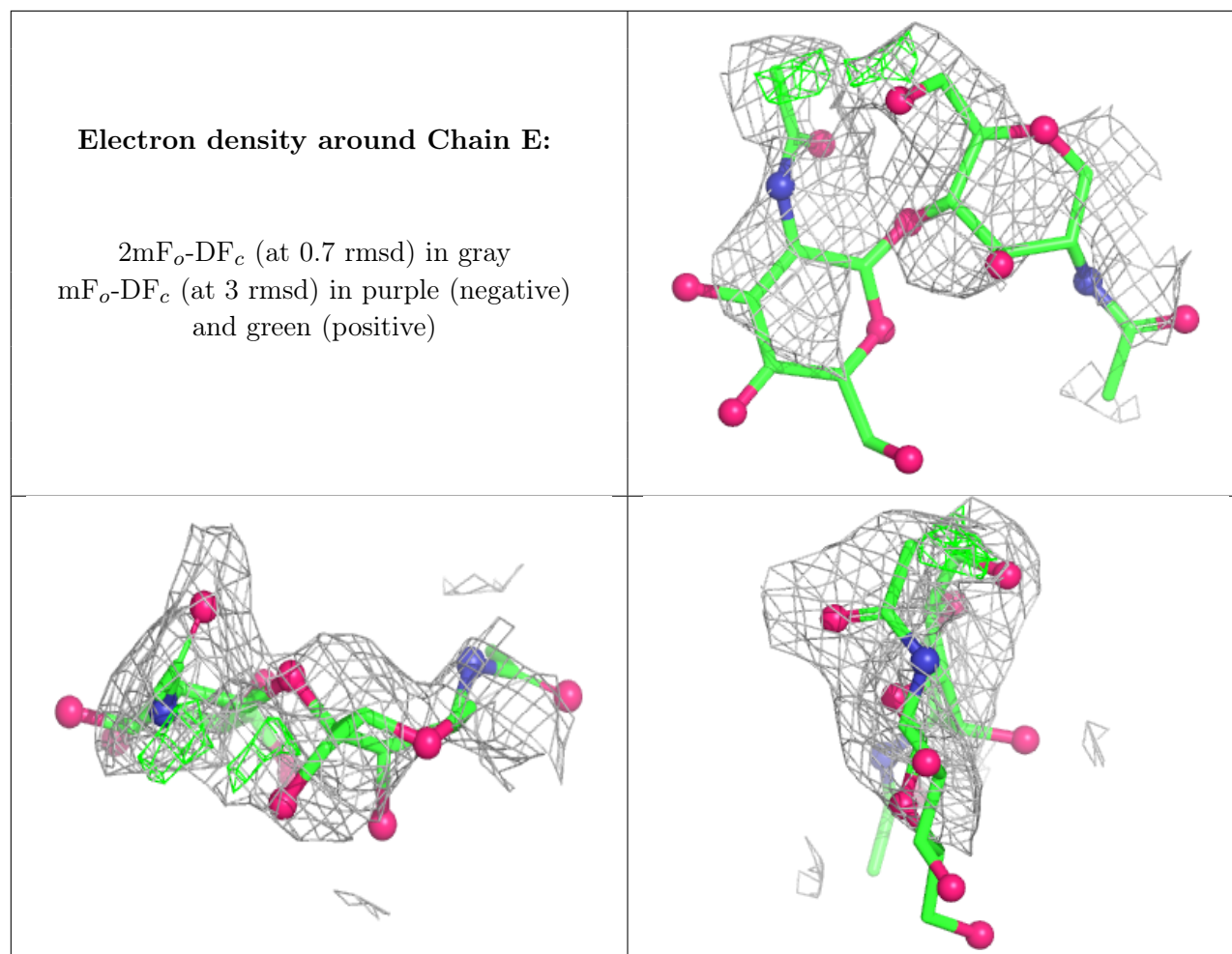
Electron density around Chain D:

$2mF_o-DF_c$ (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 [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.

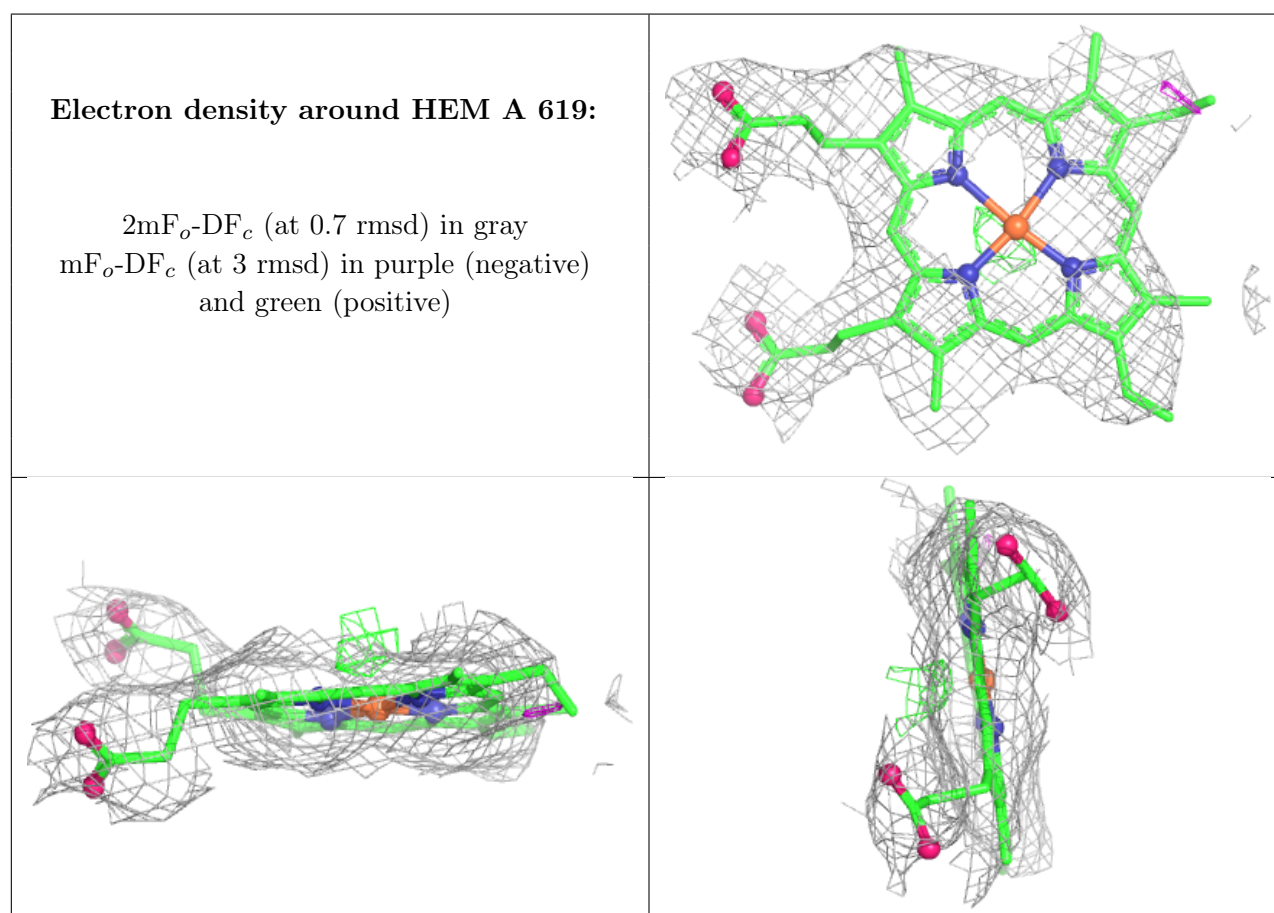
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
6	F	A	616	1/1	0.43	0.48	25,25,25,25	1
6	F	A	617	1/1	0.45	0.35	25,25,25,25	1
6	F	A	618	1/1	0.65	0.45	25,25,25,25	1
5	IOD	A	610	1/1	0.76	0.10	42,42,42,42	1
5	IOD	A	614	1/1	0.77	0.17	41,41,41,41	1
5	IOD	A	613	1/1	0.83	0.15	21,21,21,21	1
5	IOD	A	615	1/1	0.84	0.09	35,35,35,35	1
5	IOD	A	607	1/1	0.84	0.08	14,14,14,14	1
5	IOD	A	611	1/1	0.85	0.11	25,25,25,25	1
5	IOD	A	612	1/1	0.86	0.10	37,37,37,37	1

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
5	IOD	A	609	1/1	0.88	0.12	57,57,57,57	1
8	OSM	A	701	4/4	0.90	0.13	18,18,18,18	0
7	HEM	A	619	43/43	0.94	0.10	4,4,7,31	0
5	IOD	A	608	1/1	0.97	0.06	40,40,40,40	1
4	CA	A	606	1/1	0.99	0.03	5,5,5,5	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.