



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 1, 2026 – 08:51 AM UTC

PDB ID : 2OJV / pdb_00002ojv
Title : Crystal structure of a ternary complex of goat lactoperoxidase with cyanide and iodide ions at 2.4 Å resolution
Authors : Singh, A.K.; Singh, N.; Singh, S.B.; Sharma, S.; Kaur, P.; Singh, T.P.
Deposited on : 2007-01-15
Resolution : 2.40 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) : **NOT EXECUTED**
EDS : **NOT EXECUTED**
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

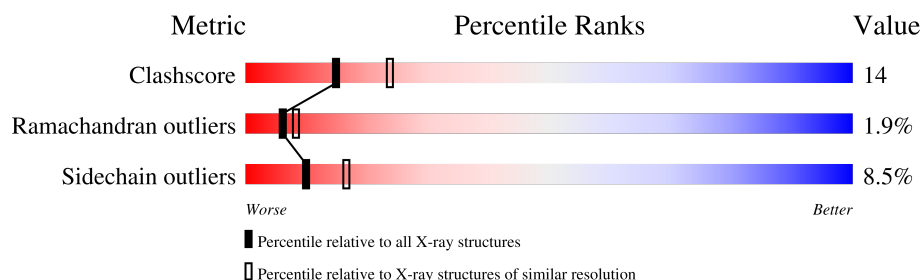
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.40 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))
Clashscore	190562	5391 (2.40-2.40)
Ramachandran outliers	187476	5320 (2.40-2.40)
Sidechain outliers	187428	5321 (2.40-2.40)

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 $\leq 5\%$.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	595	
2	B	3	
2	C	3	
2	D	3	
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
5	CYN	A	2844	-	-	X	-

2 Entry composition [i](#)

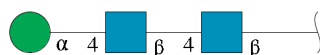
There are 9 unique types of molecules in this entry. The entry contains 5362 atoms, of which 0 are hydrogens and 0 are deuteriums.

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.

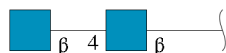
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	595	Total	C	N	O	S	0	0	0
			4753	3021	844	862	26			

- 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	C	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.

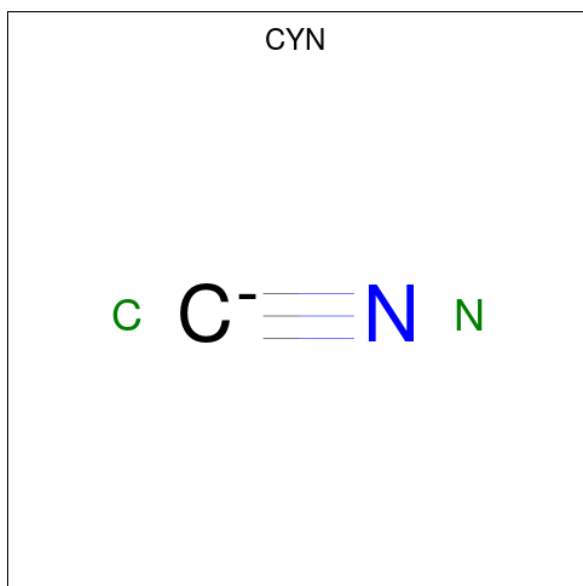


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	E	2	Total	C	N	O	0	0	0
			28	16	2	10			

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

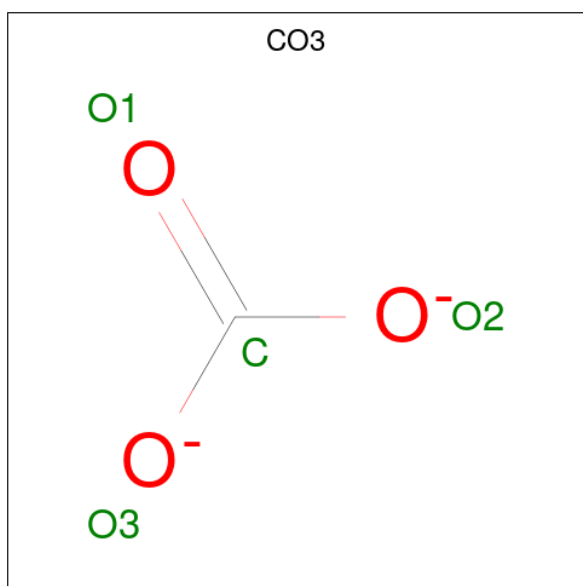
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total	Ca	0	0
			1	1		

- Molecule 5 is CYANIDE ION (CCD ID: CYN) (formula: CN).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	N	0	0
			2	1	1		

- Molecule 6 is CARBONATE ION (CCD ID: CO3) (formula: CO₃).

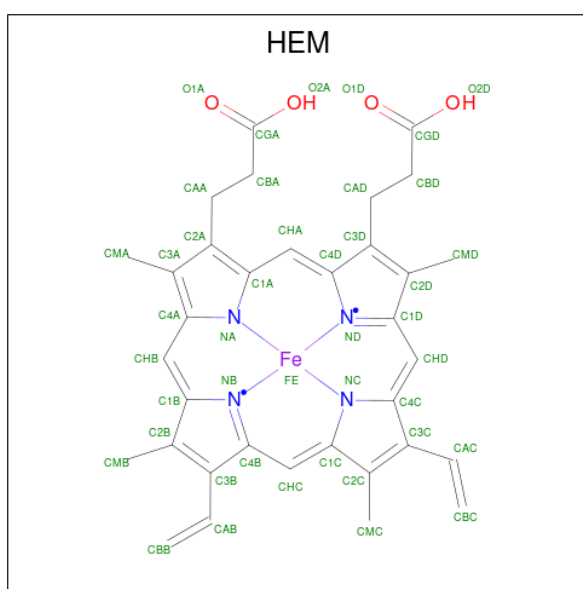


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	A	1	Total	C	O	0	0
			4	1	3		

- Molecule 7 is IODIDE ION (CCD ID: IOD) (formula: I).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	8	Total	I	0	0
			8	8		

- Molecule 8 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	Fe	N	O	
			43	34	1	4	4	0

- Molecule 9 is water.

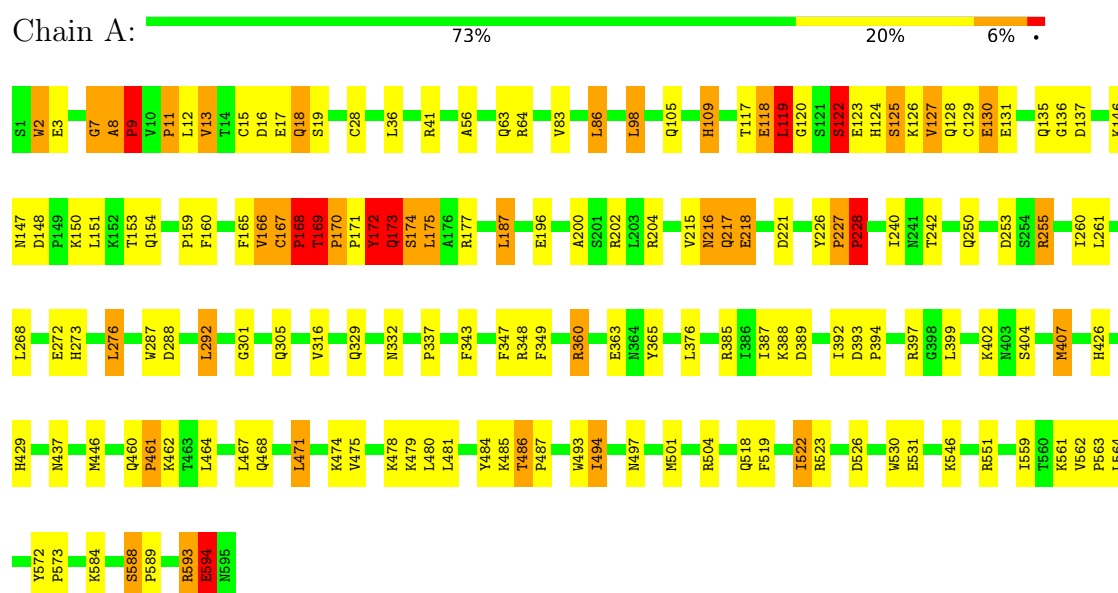
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
9	A	406	Total	O	0	0
			406	406		

3 Residue-property plots

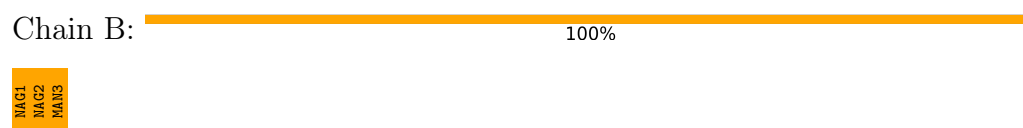
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry. Residues are color-coded 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. 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.

Note EDS was not executed.

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

Chain D:  100%

MAG1
MAG2
MAG3

- 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

Xtriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	54.20Å 80.81Å 77.05Å 90.00° 102.96° 90.00°	Depositor
Resolution (Å)	74.54 – 2.40	Depositor
% Data completeness (in resolution range)	97.8 (74.54-2.40)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	REFMAC 5.0	Depositor
R, R_{free}	0.178 , 0.228	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	5362	wwPDB-VP
Average B, all atoms (Å ²)	27.0	wwPDB-VP

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: CYN, CA, NAG, HEM, IOD, CO3, MAN

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	0.94	13/4882 (0.3%)	1.35	60/6632 (0.9%)

All (13) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	170	PRO	CA-C	7.89	1.59	1.52
1	A	169	THR	C-N	6.76	1.41	1.33
1	A	18	GLN	N-CA	-6.62	1.41	1.47
1	A	8	ALA	CA-C	6.55	1.60	1.53
1	A	18	GLN	CA-C	6.39	1.58	1.52
1	A	7	GLY	C-N	6.23	1.41	1.33
1	A	173	GLN	C-N	5.82	1.41	1.34
1	A	169	THR	CA-C	5.53	1.59	1.52
1	A	123	GLU	N-CA	-5.47	1.38	1.45
1	A	170	PRO	N-CA	5.40	1.57	1.46
1	A	7	GLY	CA-C	5.27	1.57	1.52
1	A	3	GLU	N-CA	5.19	1.52	1.46
1	A	494	ILE	CA-CB	5.06	1.61	1.54

All (60) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	11	PRO	CA-C-N	-14.47	99.66	122.67
1	A	11	PRO	C-N-CA	-14.47	99.66	122.67
1	A	13	VAL	N-CA-C	13.88	127.49	109.30
1	A	7	GLY	N-CA-C	13.51	126.17	111.85
1	A	9	PRO	CA-C-N	-12.61	113.51	122.59
1	A	9	PRO	C-N-CA	-12.61	113.51	122.59
1	A	172	TYR	CA-CB-CG	-10.65	94.73	113.90
1	A	173	GLN	OE1-CD-NE2	-10.24	112.36	122.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	12	LEU	N-CA-C	10.17	124.99	112.59
1	A	135	GLN	OE1-CD-NE2	-9.58	113.02	122.60
1	A	18	GLN	OE1-CD-NE2	-9.57	113.03	122.60
1	A	460	GLN	OE1-CD-NE2	-9.52	113.08	122.60
1	A	11	PRO	N-CA-C	9.17	131.36	112.47
1	A	217	GLN	OE1-CD-NE2	-8.90	113.69	122.60
1	A	170	PRO	N-CA-C	8.33	120.87	110.70
1	A	169	THR	CA-C-N	8.18	128.81	120.38
1	A	169	THR	C-N-CA	8.18	128.81	120.38
1	A	119	LEU	N-CA-C	-7.18	102.05	111.74
1	A	227	PRO	O-C-N	7.09	129.83	121.46
1	A	2	TRP	N-CA-C	-7.08	99.17	110.14
1	A	173	GLN	CG-CD-NE2	6.87	126.71	116.40
1	A	217	GLN	CG-CD-NE2	6.81	126.62	116.40
1	A	18	GLN	CG-CD-NE2	6.80	126.60	116.40
1	A	172	TYR	N-CA-C	6.79	125.26	110.80
1	A	135	GLN	CG-CD-NE2	6.68	126.43	116.40
1	A	126	LYS	N-CA-C	-6.59	104.34	112.38
1	A	228	PRO	CA-N-CD	-6.54	102.85	112.00
1	A	460	GLN	CG-CD-NE2	6.33	125.89	116.40
1	A	461	PRO	CA-N-CD	-6.20	103.31	112.00
1	A	171	PRO	CA-N-CD	-6.16	103.38	112.00
1	A	174	SER	CB-CA-C	-6.11	107.62	116.54
1	A	125	SER	N-CA-C	-6.11	104.90	112.90
1	A	83	VAL	N-CA-C	6.07	119.57	111.44
1	A	260	ILE	N-CA-C	6.05	116.79	110.62
1	A	11	PRO	O-C-N	-6.04	114.48	122.64
1	A	316	VAL	N-CA-C	-5.94	105.92	111.45
1	A	122	SER	CA-C-N	-5.77	112.71	123.03
1	A	122	SER	C-N-CA	-5.77	112.71	123.03
1	A	404	SER	N-CA-C	-5.52	103.41	110.53
1	A	13	VAL	N-CA-CB	-5.49	102.81	110.05
1	A	3	GLU	N-CA-C	5.47	122.45	110.80
1	A	11	PRO	CA-N-CD	-5.46	104.35	112.00
1	A	8	ALA	CA-C-N	5.44	126.64	119.84
1	A	8	ALA	C-N-CA	5.44	126.64	119.84
1	A	18	GLN	N-CA-C	-5.43	105.61	113.21
1	A	124	HIS	CA-CB-CG	-5.38	108.42	113.80
1	A	7	GLY	CA-C-N	5.35	129.58	123.21
1	A	7	GLY	C-N-CA	5.35	129.58	123.21
1	A	18	GLN	CB-CG-CD	5.33	121.65	112.60
1	A	461	PRO	N-CA-C	5.31	119.55	110.95

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	170	PRO	CB-CA-C	5.27	117.35	110.92
1	A	109	HIS	CA-CB-CG	-5.26	108.54	113.80
1	A	168	PRO	N-CA-C	5.23	123.24	112.47
1	A	166	VAL	N-CA-C	5.22	120.20	109.34
1	A	170	PRO	CA-N-CD	-5.17	104.77	112.00
1	A	8	ALA	O-C-N	5.12	126.38	121.31
1	A	559	ILE	CA-C-N	-5.11	114.56	122.49
1	A	559	ILE	C-N-CA	-5.11	114.56	122.49
1	A	429	HIS	N-CA-C	-5.07	102.56	109.71
1	A	493	TRP	N-CA-C	5.03	116.45	111.07

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	4753	0	4647	127	0
2	B	39	0	34	2	0
2	C	39	0	34	0	0
2	D	39	0	34	2	0
3	E	28	0	25	0	0
4	A	1	0	0	0	0
5	A	2	0	0	2	0
6	A	4	0	0	0	0
7	A	8	0	0	2	0
8	A	43	0	30	3	0
9	A	406	0	0	18	0
All	All	5362	0	4804	134	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 14.

All (134) 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:168:PRO:HG2	1:A:172:TYR:CE1	1.85	1.11
1:A:63:GLN:HG2	9:A:3178:HOH:O	1.50	1.10
1:A:167:CYS:HB3	1:A:168:PRO:HD2	1.34	1.09
1:A:426:HIS:HB3	9:A:3202:HOH:O	1.55	1.05
1:A:2:TRP:HB2	1:A:175:LEU:HD12	1.35	1.03
1:A:169:THR:CG2	1:A:170:PRO:HD3	1.89	1.02
1:A:64:ARG:HG3	9:A:3237:HOH:O	1.64	0.96
1:A:119:LEU:HD11	1:A:169:THR:HG21	1.48	0.96
1:A:167:CYS:HB3	1:A:168:PRO:CD	1.97	0.95
1:A:169:THR:HG22	1:A:170:PRO:HD3	1.49	0.93
1:A:169:THR:HG22	1:A:170:PRO:CD	2.00	0.91
1:A:118:GLU:HB3	9:A:3243:HOH:O	1.72	0.90
1:A:551:ARG:HD3	1:A:584:LYS:HA	1.55	0.87
1:A:216:ASN:ND2	1:A:218:GLU:H	1.72	0.87
1:A:167:CYS:CB	1:A:168:PRO:HD2	2.07	0.84
1:A:481:LEU:HD21	1:A:487:PRO:HD3	1.56	0.84
5:A:2844:CYN:N	7:A:2852:IOD:I	2.82	0.82
1:A:202:ARG:NH2	7:A:2848:IOD:I	2.83	0.81
1:A:119:LEU:HD11	1:A:169:THR:CG2	2.11	0.81
1:A:169:THR:HG23	1:A:170:PRO:HD3	1.61	0.81
1:A:19:SER:HB3	9:A:3189:HOH:O	1.79	0.81
1:A:130:GLU:HG2	1:A:159:PRO:HG3	1.62	0.80
1:A:172:TYR:O	1:A:173:GLN:HB3	1.81	0.80
1:A:146:LYS:HE2	9:A:3257:HOH:O	1.84	0.77
1:A:216:ASN:HD22	1:A:218:GLU:H	1.31	0.76
1:A:169:THR:CG2	1:A:170:PRO:CD	2.59	0.76
1:A:253:ASP:OD2	1:A:255:ARG:HD3	1.87	0.75
1:A:172:TYR:HD2	1:A:173:GLN:H	1.34	0.74
1:A:481:LEU:CD2	1:A:487:PRO:HD3	2.18	0.73
1:A:167:CYS:CB	1:A:168:PRO:CD	2.63	0.72
1:A:561:LYS:HE2	9:A:2905:HOH:O	1.89	0.72
1:A:365:TYR:CZ	1:A:397:ARG:HG2	2.25	0.72
1:A:146:LYS:O	1:A:147:ASN:HB2	1.88	0.71
1:A:130:GLU:HG3	9:A:3201:HOH:O	1.93	0.69
1:A:174:SER:OG	1:A:175:LEU:N	2.25	0.69
1:A:216:ASN:HD22	1:A:217:GLN:N	1.93	0.67
1:A:407:MET:HB3	1:A:501:MET:HE3	1.77	0.67
1:A:467:LEU:HG	1:A:471:LEU:HD22	1.77	0.66
1:A:169:THR:HG22	1:A:170:PRO:HD2	1.77	0.66
1:A:288:ASP:O	1:A:292:LEU:HD22	1.96	0.66
1:A:169:THR:H	1:A:170:PRO:HD2	1.61	0.64
1:A:8:ALA:HB3	9:A:3247:HOH:O	1.97	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:129:CYS:HB2	9:A:3188:HOH:O	1.98	0.63
1:A:119:LEU:CD1	1:A:169:THR:HG21	2.25	0.63
1:A:168:PRO:CG	1:A:172:TYR:CE1	2.75	0.63
8:A:2853:HEM:HBC2	8:A:2853:HEM:HMC2	1.81	0.62
1:A:216:ASN:HD22	1:A:216:ASN:C	2.07	0.62
1:A:2:TRP:HB2	1:A:175:LEU:CD1	2.22	0.61
2:B:1:NAG:H61	2:B:2:NAG:C1	2.31	0.61
1:A:168:PRO:HG2	1:A:172:TYR:HE1	1.58	0.61
1:A:475:VAL:HG12	1:A:479:LYS:HE3	1.81	0.61
1:A:594:GLU:OE2	1:A:594:GLU:HA	2.02	0.59
1:A:8:ALA:N	1:A:9:PRO:CD	2.65	0.59
1:A:64:ARG:HD3	9:A:3249:HOH:O	2.02	0.59
1:A:7:GLY:HA2	1:A:167:CYS:HA	1.84	0.59
1:A:8:ALA:N	1:A:9:PRO:HD2	2.18	0.58
1:A:15:CYS:HG	1:A:28:CYS:HG	0.60	0.58
1:A:125:SER:HA	1:A:128:GLN:HB3	1.87	0.57
1:A:348:ARG:HH11	1:A:437:ASN:ND2	2.03	0.57
1:A:169:THR:H	1:A:170:PRO:CD	2.18	0.57
1:A:202:ARG:HD2	1:A:250:GLN:HE22	1.69	0.57
1:A:343:PHE:CD1	1:A:518:GLN:HG2	2.40	0.56
1:A:119:LEU:CD1	1:A:169:THR:CG2	2.80	0.56
1:A:125:SER:C	1:A:127:VAL:N	2.60	0.56
1:A:172:TYR:HD2	1:A:173:GLN:N	2.04	0.54
1:A:462:LYS:HD2	9:A:3136:HOH:O	2.09	0.53
1:A:588:SER:N	1:A:589:PRO:CD	2.71	0.53
1:A:530:TRP:CE2	1:A:531:GLU:HG3	2.44	0.53
1:A:16:ASP:O	1:A:17:GLU:HB2	2.09	0.52
1:A:349:PHE:CB	1:A:497:ASN:HD21	2.23	0.52
2:D:1:NAG:H62	2:D:2:NAG:C1	2.40	0.52
1:A:125:SER:C	1:A:127:VAL:H	2.19	0.51
1:A:172:TYR:CD2	1:A:173:GLN:N	2.77	0.51
1:A:120:GLY:C	1:A:122:SER:H	2.19	0.51
2:B:2:NAG:H3	2:B:3:MAN:O5	2.11	0.51
1:A:187:LEU:HD13	1:A:305:GLN:HA	1.92	0.51
1:A:468:GLN:HG2	1:A:474:LYS:HA	1.92	0.51
1:A:16:ASP:HB3	9:A:3189:HOH:O	2.11	0.50
1:A:216:ASN:HB2	1:A:227:PRO:O	2.11	0.50
1:A:221:ASP:HB2	1:A:226:TYR:CZ	2.45	0.50
1:A:407:MET:HE3	1:A:407:MET:C	2.37	0.49
1:A:365:TYR:CE1	1:A:397:ARG:HG2	2.48	0.49
1:A:287:TRP:HE3	1:A:292:LEU:HD13	1.78	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:301:GLY:O	1:A:305:GLN:HG3	2.14	0.47
1:A:349:PHE:HA	1:A:497:ASN:HD21	1.80	0.47
1:A:593:ARG:HD3	1:A:593:ARG:HA	1.79	0.47
1:A:216:ASN:ND2	1:A:217:GLN:N	2.63	0.47
1:A:393:ASP:HB2	1:A:394:PRO:HD3	1.96	0.47
1:A:36:LEU:HG	1:A:337:PRO:HD2	1.96	0.47
1:A:485:LYS:O	1:A:486:THR:HB	2.15	0.47
1:A:7:GLY:C	1:A:9:PRO:HD2	2.40	0.46
1:A:105:GLN:HE21	5:A:2844:CYN:C	2.28	0.46
1:A:109:HIS:ND1	9:A:2881:HOH:O	2.28	0.46
1:A:148:ASP:O	1:A:151:LEU:HB2	2.15	0.46
1:A:348:ARG:HH11	1:A:437:ASN:HD22	1.63	0.45
2:D:2:NAG:H4	2:D:3:MAN:H2	1.65	0.45
1:A:169:THR:N	1:A:170:PRO:HD2	2.31	0.45
1:A:215:VAL:HA	1:A:228:PRO:HD3	1.97	0.45
1:A:15:CYS:CB	1:A:28:CYS:HG	2.24	0.44
1:A:131:GLU:OE2	1:A:131:GLU:HA	2.17	0.44
1:A:136:GLY:O	1:A:137:ASP:C	2.58	0.44
1:A:150:LYS:O	1:A:154:GLN:N	2.46	0.44
1:A:478:LYS:HE3	1:A:478:LYS:HB2	1.50	0.44
1:A:273:HIS:CD2	1:A:273:HIS:C	2.95	0.44
1:A:332:ASN:ND2	9:A:3196:HOH:O	2.41	0.43
1:A:360:ARG:NH2	1:A:389:ASP:OD2	2.51	0.43
1:A:202:ARG:HD2	1:A:250:GLN:NE2	2.34	0.43
1:A:15:CYS:CB	1:A:28:CYS:SG	3.06	0.43
1:A:196:GLU:CD	9:A:3142:HOH:O	2.60	0.43
1:A:572:TYR:HA	1:A:573:PRO:HA	1.62	0.43
1:A:119:LEU:HD21	1:A:169:THR:HG23	2.00	0.43
1:A:272:GLU:HG3	1:A:276:LEU:HD22	2.01	0.43
1:A:562:VAL:HB	1:A:563:PRO:HD2	2.01	0.43
1:A:484:TYR:O	1:A:485:LYS:HB2	2.19	0.42
1:A:9:PRO:CG	1:A:41:ARG:CZ	2.96	0.42
1:A:160:PHE:HA	9:A:3144:HOH:O	2.19	0.42
1:A:385:ARG:O	1:A:389:ASP:HB3	2.19	0.42
1:A:481:LEU:HD23	1:A:481:LEU:HA	1.74	0.42
8:A:2853:HEM:HBC2	8:A:2853:HEM:CMC	2.48	0.42
1:A:594:GLU:OE2	1:A:594:GLU:CA	2.66	0.42
1:A:519:PHE:CD1	1:A:522:ILE:HD11	2.55	0.42
1:A:446:MET:HA	1:A:446:MET:HE2	2.02	0.42
1:A:387:ILE:HG22	1:A:388:LYS:HG3	2.02	0.42
1:A:588:SER:N	1:A:589:PRO:HD2	2.35	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:98:LEU:HD13	1:A:399:LEU:HD23	2.01	0.41
1:A:200:ALA:O	1:A:204:ARG:HG3	2.21	0.41
1:A:347:PHE:HB3	8:A:2853:HEM:HMD3	2.02	0.41
1:A:86:LEU:HD13	9:A:3037:HOH:O	2.21	0.41
1:A:522:ILE:O	1:A:526:ASP:HB2	2.20	0.41
1:A:9:PRO:O	1:A:11:PRO:HD3	2.20	0.40
1:A:165:PHE:CD1	1:A:165:PHE:N	2.90	0.40
1:A:402:LYS:HA	1:A:402:LYS:HD3	1.95	0.40
1:A:407:MET:HB3	1:A:501:MET:CE	2.47	0.40
1:A:120:GLY:C	1:A:122:SER:N	2.77	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	593/595 (100%)	557 (94%)	25 (4%)	11 (2%)	6 8

All (11) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	167	CYS
1	A	168	PRO
1	A	169	THR
1	A	173	GLN
1	A	594	GLU
1	A	242	THR
1	A	486	THR
1	A	56	ALA
1	A	172	TYR
1	A	9	PRO
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	517/517 (100%)	473 (92%)	44 (8%)	10	16

All (44) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	9	PRO
1	A	13	VAL
1	A	18	GLN
1	A	86	LEU
1	A	98	LEU
1	A	117	THR
1	A	118	GLU
1	A	119	LEU
1	A	122	SER
1	A	127	VAL
1	A	130	GLU
1	A	153	THR
1	A	173	GLN
1	A	175	LEU
1	A	177	ARG
1	A	187	LEU
1	A	216	ASN
1	A	218	GLU
1	A	228	PRO
1	A	240	ILE
1	A	255	ARG
1	A	261	LEU
1	A	268	LEU
1	A	276	LEU
1	A	292	LEU
1	A	329	GLN
1	A	360	ARG
1	A	363	GLU
1	A	376	LEU
1	A	392	ILE

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Mol	Chain	Res	Type
1	A	407	MET
1	A	461	PRO
1	A	464	LEU
1	A	471	LEU
1	A	480	LEU
1	A	494	ILE
1	A	504	ARG
1	A	522	ILE
1	A	523	ARG
1	A	546	LYS
1	A	564	LEU
1	A	588	SER
1	A	593	ARG
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	18	GLN
1	A	147	ASN
1	A	216	ASN
1	A	222	HIS
1	A	245	HIS
1	A	250	GLN
1	A	329	GLN
1	A	341	ASN
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 ⓘ

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

5.5 Carbohydrates ⓘ

11 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	1,2	14,14,15	0.57	0	17,19,21	1.50	3 (17%)
2	NAG	B	2	2	14,14,15	0.54	0	17,19,21	1.52	2 (11%)
2	MAN	B	3	2	11,11,12	0.64	0	15,15,17	1.35	2 (13%)
2	NAG	C	1	1,2	14,14,15	0.60	0	17,19,21	0.81	0
2	NAG	C	2	2	14,14,15	0.72	0	17,19,21	1.37	2 (11%)
2	MAN	C	3	2	11,11,12	0.56	0	15,15,17	0.98	1 (6%)
2	NAG	D	1	1,2	14,14,15	0.78	1 (7%)	17,19,21	1.41	2 (11%)
2	NAG	D	2	2	14,14,15	0.56	0	17,19,21	1.78	4 (23%)
2	MAN	D	3	2	11,11,12	0.51	0	15,15,17	1.58	2 (13%)
3	NAG	E	1	3,1	14,14,15	0.64	0	17,19,21	2.10	5 (29%)
3	NAG	E	2	3	14,14,15	0.53	0	17,19,21	0.87	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
2	NAG	B	1	1,2	-	2/6/23/26	0/1/1/1
2	NAG	B	2	2	-	3/6/23/26	0/1/1/1
2	MAN	B	3	2	-	2/2/19/22	0/1/1/1
2	NAG	C	1	1,2	-	0/6/23/26	0/1/1/1
2	NAG	C	2	2	-	2/6/23/26	0/1/1/1
2	MAN	C	3	2	-	2/2/19/22	0/1/1/1
2	NAG	D	1	1,2	-	0/6/23/26	0/1/1/1
2	NAG	D	2	2	-	1/6/23/26	0/1/1/1
2	MAN	D	3	2	-	2/2/19/22	1/1/1/1
3	NAG	E	1	3,1	-	0/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	NAG	E	2	3	-	2/6/23/26	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	1	NAG	O5-C1	-2.07	1.40	1.43

All (23) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	2	NAG	C1-O5-C5	5.34	119.34	112.19
2	D	3	MAN	C1-O5-C5	5.16	119.10	112.19
3	E	1	NAG	C4-C3-C2	4.42	117.49	111.02
2	B	2	NAG	C4-C3-C2	4.12	117.06	111.02
3	E	1	NAG	O5-C1-C2	-3.91	105.23	111.29
2	C	2	NAG	C1-O5-C5	3.90	117.41	112.19
3	E	1	NAG	O5-C5-C6	3.79	115.04	107.66
3	E	1	NAG	C1-O5-C5	-3.72	107.20	112.19
2	D	1	NAG	O5-C1-C2	-3.58	105.75	111.29
2	B	3	MAN	C1-O5-C5	3.57	116.97	112.19
2	B	1	NAG	C2-N2-C7	-3.47	118.25	122.90
2	B	1	NAG	O5-C1-C2	-3.18	106.38	111.29
2	D	1	NAG	C6-C5-C4	-2.90	105.91	113.02
2	B	3	MAN	C1-C2-C3	2.82	113.76	109.64
3	E	1	NAG	C2-N2-C7	-2.73	119.25	122.90
2	B	2	NAG	O5-C5-C6	2.66	112.85	107.66
2	D	2	NAG	C2-N2-C7	-2.56	119.47	122.90
2	C	3	MAN	C1-O5-C5	2.37	115.37	112.19
2	D	2	NAG	C4-C3-C2	-2.29	107.66	111.02
2	B	1	NAG	O4-C4-C5	2.21	114.76	109.32
2	C	2	NAG	C2-N2-C7	-2.12	120.06	122.90
2	D	3	MAN	C1-C2-C3	2.05	112.62	109.64
2	D	2	NAG	C1-C2-N2	2.01	113.59	110.43

There are no chirality outliers.

All (16) torsion outliers are listed below:

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

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

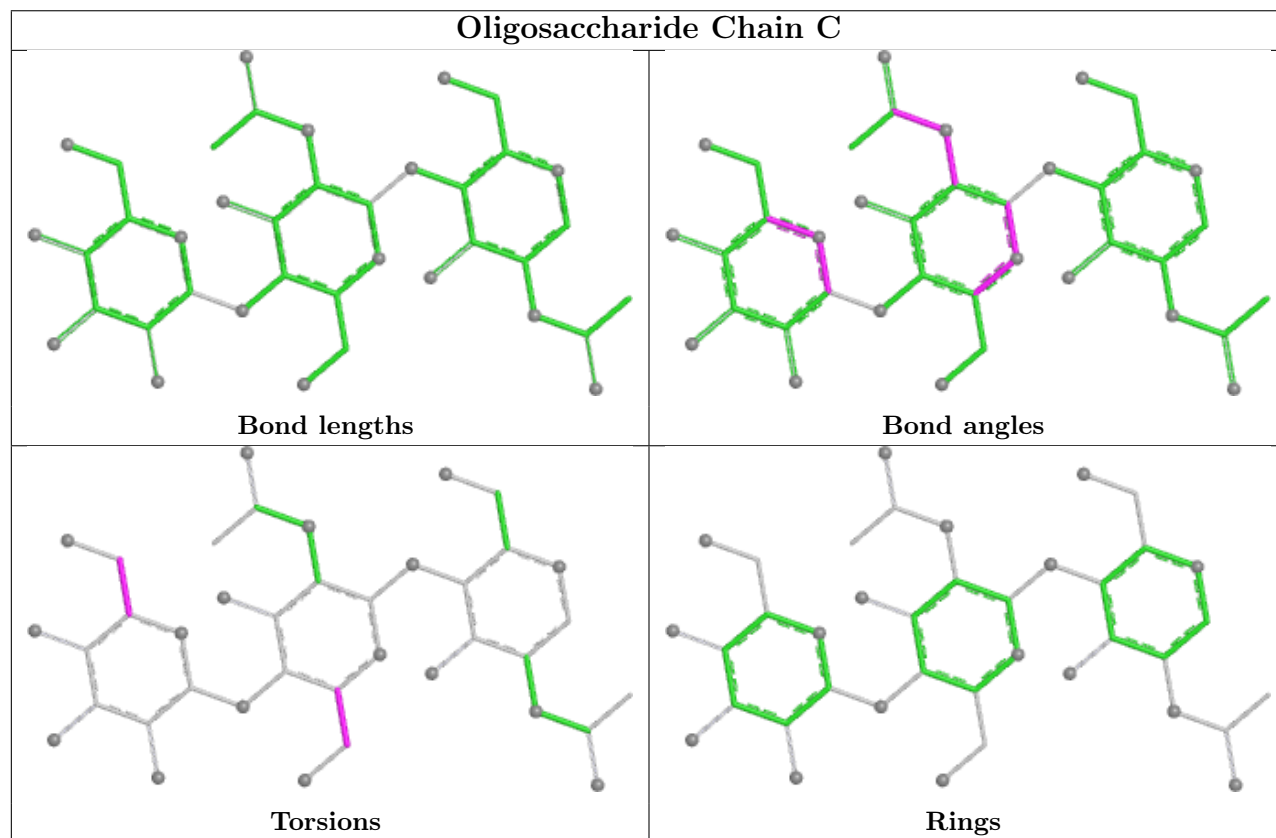
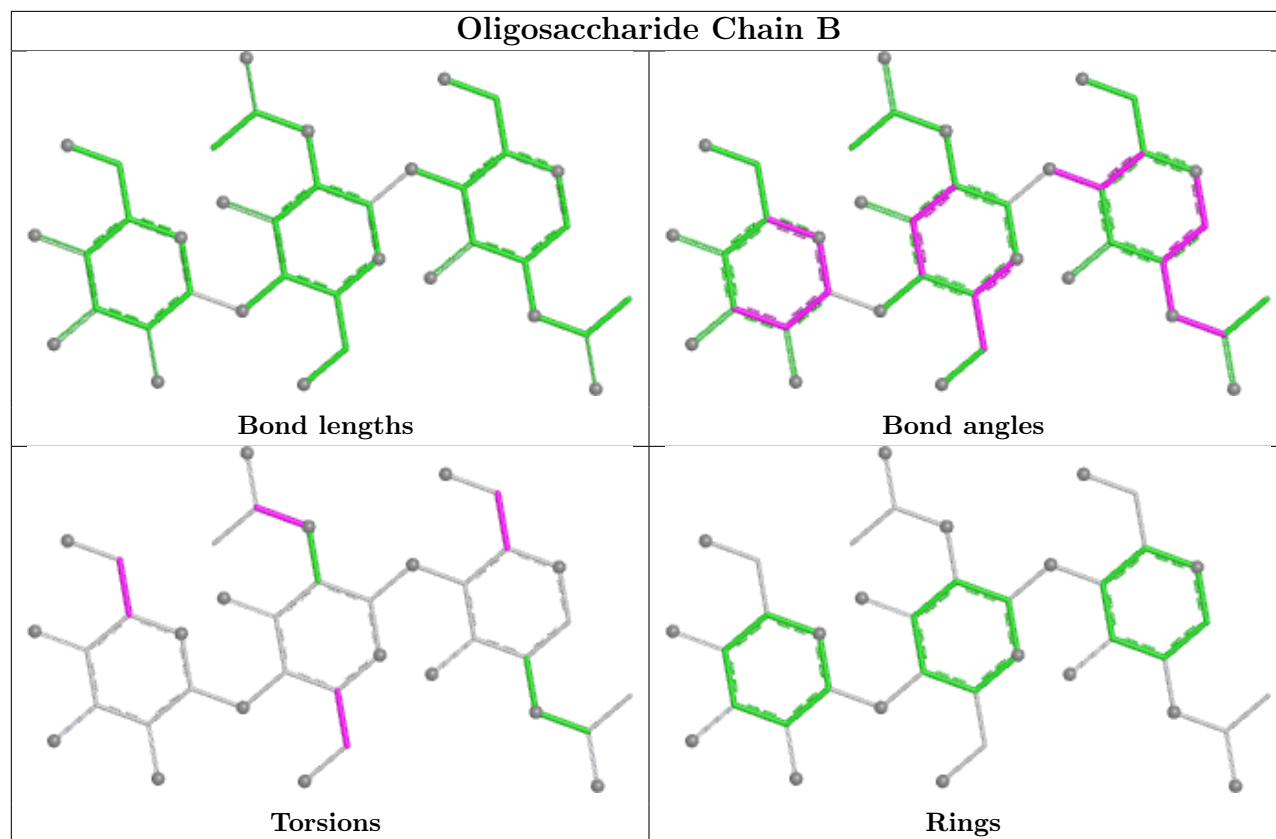
All (1) ring outliers are listed below:

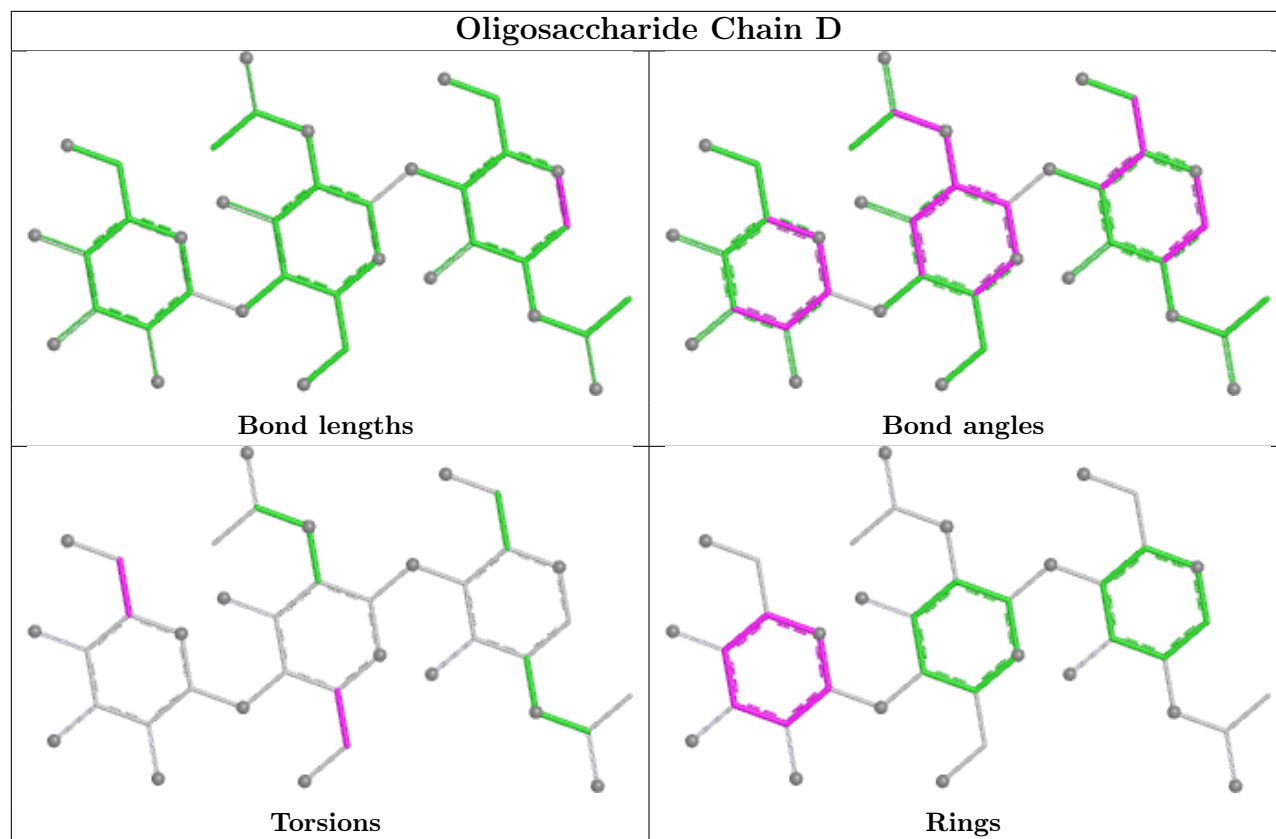
Mol	Chain	Res	Type	Atoms
2	D	3	MAN	C1-C2-C3-C4-C5-O5

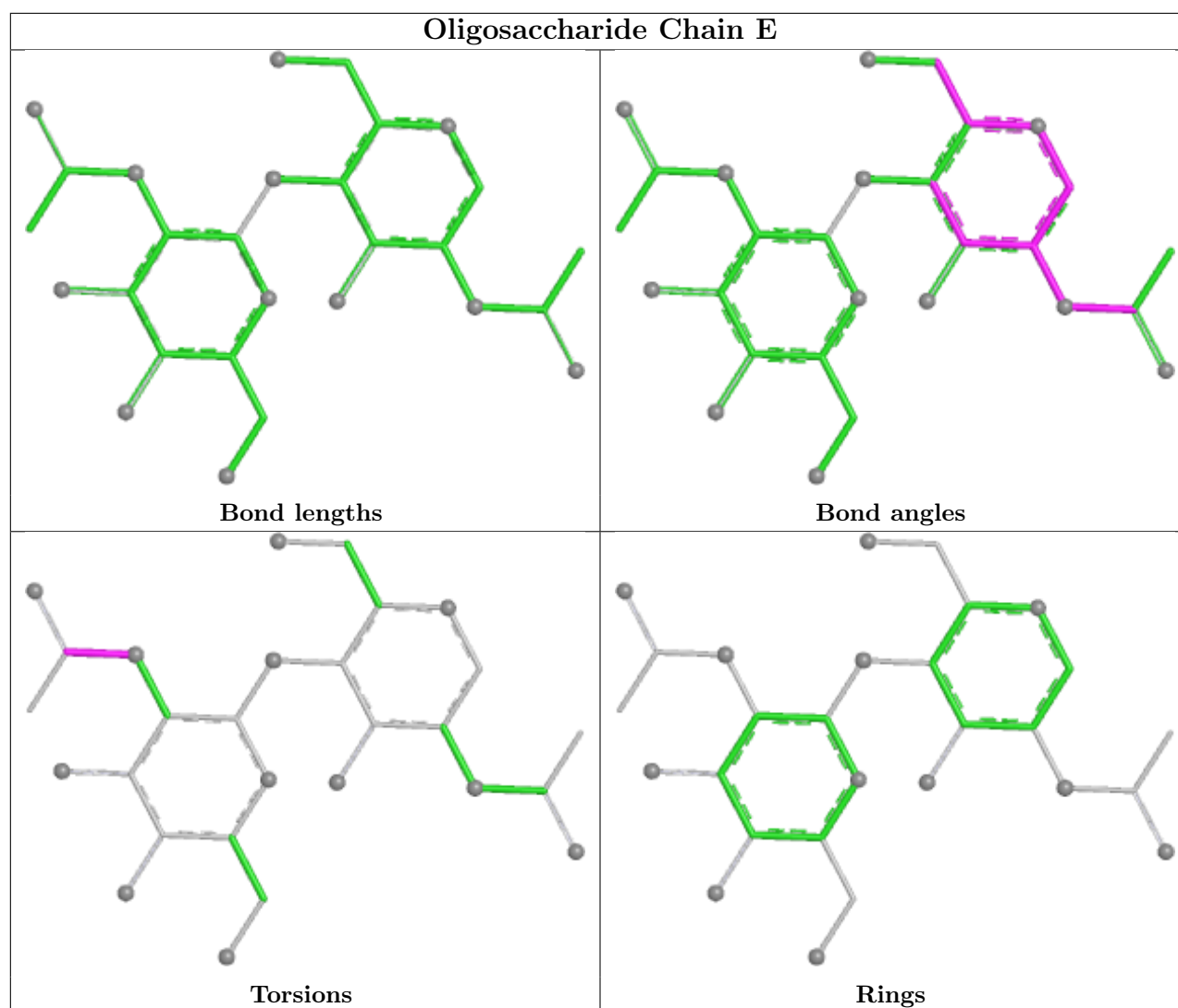
6 monomers are involved in 4 short contacts:

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







5.6 Ligand geometry [i](#)

Of 12 ligands modelled in this entry, 9 are monoatomic - leaving 3 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	$\# Z > 2$
8	HEM	A	2853	1,5	50,50,50	1.81	10 (20%)	67,82,82	1.09	4 (5%)
5	CYN	A	2844	8	1,1,1	0.33	0	-		
6	CO3	A	688	-	3,3,3	0.70	0	2,3,3	0.38	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
8	HEM	A	2853	1,5	-	4/14/54/54	-

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
8	A	2853	HEM	C3D-C2D	6.99	1.51	1.36
8	A	2853	HEM	FE-NC	3.86	2.07	1.95
8	A	2853	HEM	CAC-C3C	3.22	1.56	1.47
8	A	2853	HEM	FE-ND	3.21	2.04	1.94
8	A	2853	HEM	FE-NB	3.15	2.04	1.94
8	A	2853	HEM	CAB-C3B	2.95	1.55	1.47
8	A	2853	HEM	FE-NA	2.53	2.03	1.95
8	A	2853	HEM	CMB-C2B	2.12	1.55	1.50
8	A	2853	HEM	CMC-C2C	2.06	1.55	1.50
8	A	2853	HEM	CMA-C3A	2.03	1.54	1.50

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
8	A	2853	HEM	C4D-ND-C1D	3.82	109.73	105.21
8	A	2853	HEM	CHC-C4B-NB	2.67	127.30	124.42
8	A	2853	HEM	C3B-C2B-C1B	2.41	108.22	106.41
8	A	2853	HEM	CAD-C3D-C4D	2.32	128.75	124.70

There are no chirality outliers.

All (4) torsion outliers are listed below:

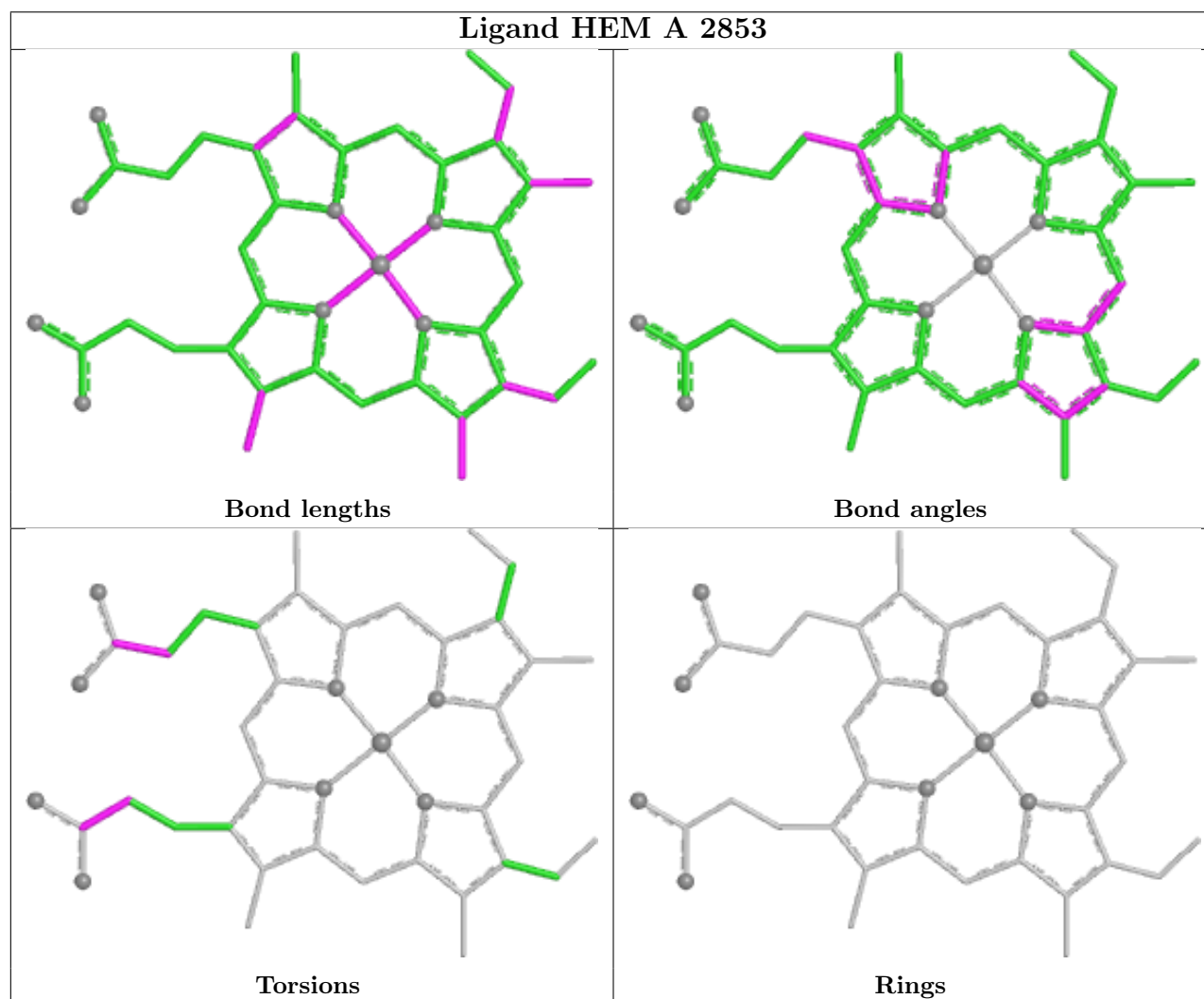
Mol	Chain	Res	Type	Atoms
8	A	2853	HEM	CAD-CBD-CGD-O2D
8	A	2853	HEM	CAD-CBD-CGD-O1D
8	A	2853	HEM	CAA-CBA-CGA-O2A
8	A	2853	HEM	CAA-CBA-CGA-O1A

There are no ring outliers.

2 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
8	A	2853	HEM	3	0
5	A	2844	CYN	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

6.1 Protein, DNA and RNA chains

EDS was not executed - this section is therefore empty.

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

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates

EDS was not executed - this section is therefore empty.

6.4 Ligands

EDS was not executed - this section is therefore empty.

6.5 Other polymers

EDS was not executed - this section is therefore empty.