



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 11:51 AM UTC

PDB ID : 2O86 / pdb_00002o86
Title : Crystal structure of a ternary complex of buffalo lactoperoxidase with nitrate and iodide at 2.8 Å resolution
Authors : Sheikh, I.A.; Singh, N.; Singh, A.K.; Sharma, S.; Kaur, P.; Singh, T.P.
Deposited on : 2006-12-12
Resolution : 2.80 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

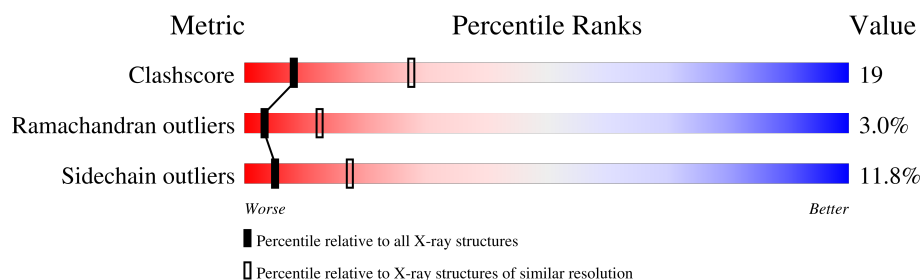
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 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	4276 (2.80-2.80)
Ramachandran outliers	187476	4196 (2.80-2.80)
Sidechain outliers	187428	4198 (2.80-2.80)

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	61% 30% 6% .
2	B	3	33% 67%
2	D	3	100%
3	C	2	50% 50%
3	E	2	50% 50%

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
6	IOD	A	2003	-	-	X	-
7	NO3	A	3007	-	-	X	-

2 Entry composition [i](#)

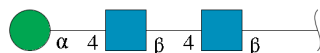
There are 9 unique types of molecules in this entry. The entry contains 5219 atoms, of which 0 are hydrogens and 0 are deuteriums.

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
			4766	3032	845	862	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.

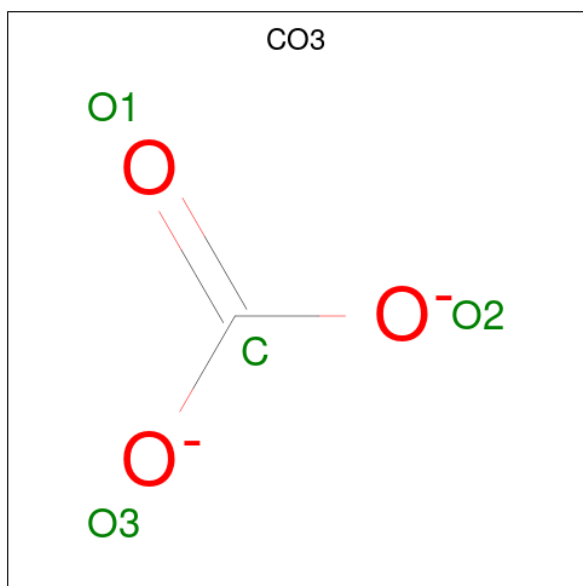


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	C	2	Total	C	N	O	0	0	0
			28	16	2	10			
3	E	2	Total	C	N	O	0	0	0
			28	16	2	10			

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

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

- Molecule 5 is CARBONATE ION (CCD ID: CO3) (formula: CO₃).

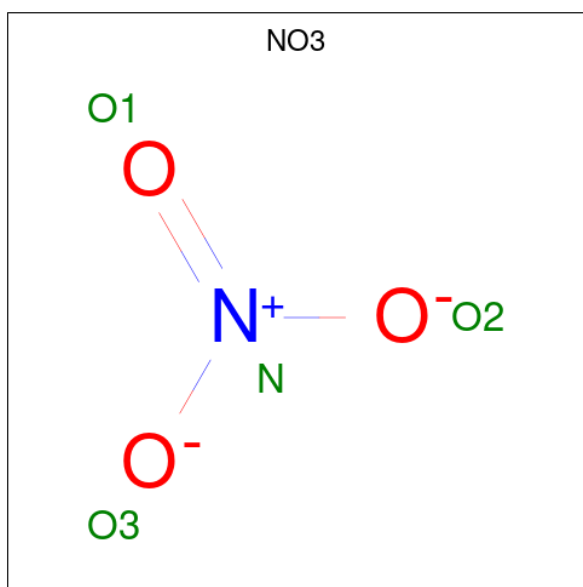


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			4	1	3		

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

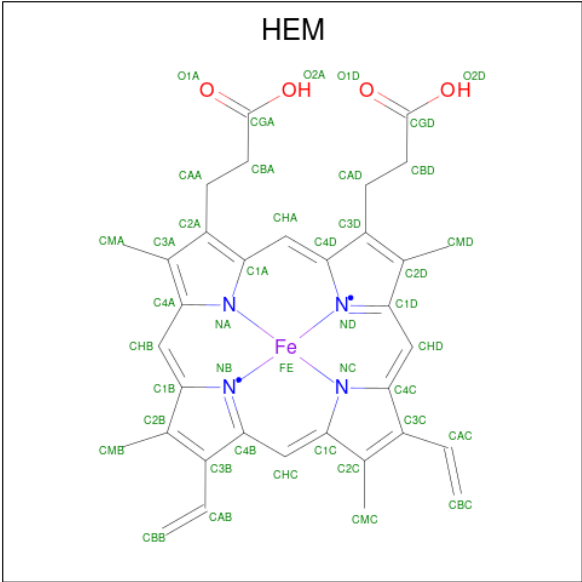
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	7	Total	I	0	0
			7	7		

- Molecule 7 is NITRATE ION (CCD ID: NO3) (formula: NO₃).



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

- Molecule 8 is PROTOPORPHYRIN IX CONTAINING FE (CCD ID: HEM) (formula: C₃₄H₃₂FeN₄O₄).



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

- Molecule 9 is water.

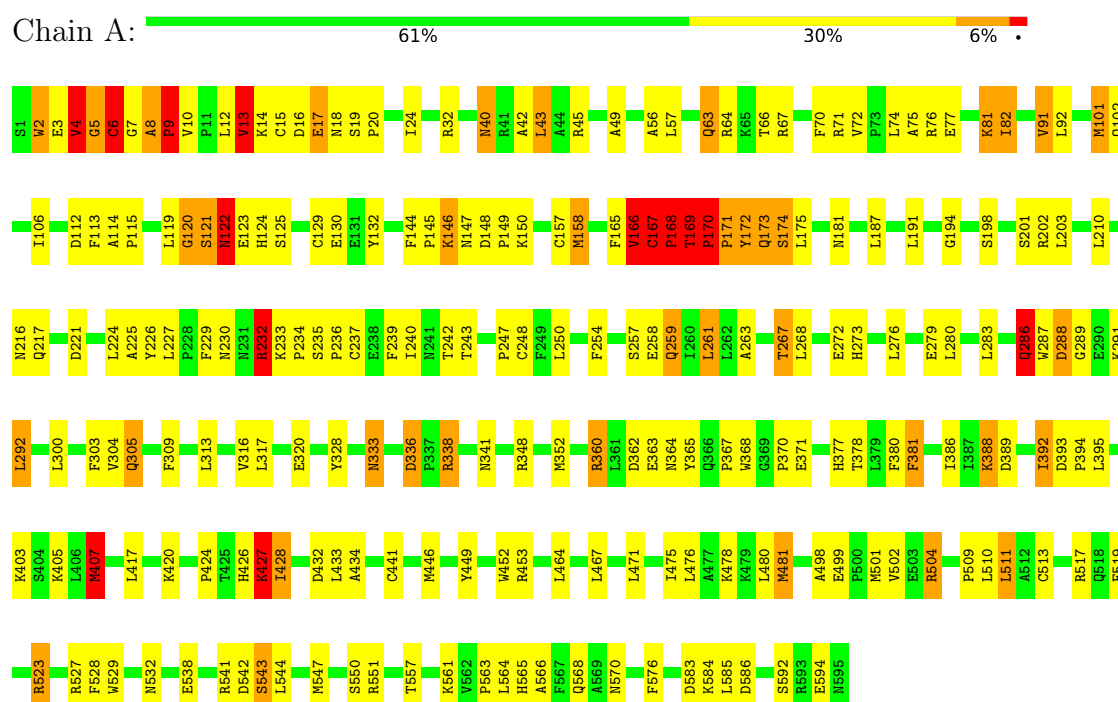
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
9	A	236	Total	O	0	0
			236	236		

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. 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 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

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	54.53Å 80.83Å 77.66Å 90.00° 102.96° 90.00°	Depositor
Resolution (Å)	20.00 – 2.80	Depositor
% Data completeness (in resolution range)	96.5 (20.00-2.80)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	REFMAC 5.0	Depositor
R, R_{free}	0.209 , 0.218	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	5219	wwPDB-VP
Average B, all atoms (Å ²)	39.0	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: NO3, CA, IOD, HEM, NAG, 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	1.05	4/4893 (0.1%)	1.33	51/6638 (0.8%)

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	1

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	158	MET	SD-CE	-6.57	1.63	1.79
1	A	170	PRO	CA-C	6.42	1.58	1.52
1	A	101	MET	SD-CE	-5.51	1.65	1.79
1	A	305	GLN	C-O	-5.20	1.17	1.24

All (51) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	169	THR	N-CA-C	-10.70	86.16	109.81
1	A	286	GLN	OE1-CD-NE2	-10.67	111.93	122.60
1	A	121	SER	N-CA-C	-9.26	102.11	112.57
1	A	336	ASP	CB-CA-C	8.90	118.98	110.17
1	A	170	PRO	N-CA-C	7.77	120.18	110.70
1	A	428	ILE	N-CA-C	7.52	118.25	110.05
1	A	288	ASP	N-CA-C	7.49	122.22	110.17
1	A	287	TRP	CA-C-N	7.32	135.08	122.64
1	A	287	TRP	C-N-CA	7.32	135.08	122.64

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	288	ASP	CB-CA-C	-7.21	97.32	109.50
1	A	427	LYS	N-CA-C	7.12	122.23	112.90
1	A	82	ILE	CB-CA-C	-6.99	104.06	110.91
1	A	286	GLN	CG-CD-NE2	6.80	126.61	116.40
1	A	566	ALA	N-CA-C	-6.78	104.69	114.12
1	A	288	ASP	CA-CB-CG	6.74	119.34	112.60
1	A	232	ARG	N-CA-C	6.68	120.32	110.30
1	A	565	HIS	CA-C-N	-6.67	111.76	122.23
1	A	565	HIS	C-N-CA	-6.67	111.76	122.23
1	A	225	ALA	N-CA-C	6.46	119.27	110.55
1	A	316	VAL	N-CA-C	-6.41	106.30	111.81
1	A	392	ILE	N-CA-C	6.35	119.95	111.44
1	A	170	PRO	CB-CA-C	6.11	118.37	110.92
1	A	594	GLU	N-CA-C	5.92	117.43	110.97
1	A	194	GLY	CA-C-O	-5.88	118.17	122.23
1	A	173	GLN	N-CA-C	5.72	122.99	110.80
1	A	169	THR	O-C-N	-5.68	114.78	121.32
1	A	171	PRO	CA-C-N	-5.62	113.77	122.59
1	A	171	PRO	C-N-CA	-5.62	113.77	122.59
1	A	92	LEU	N-CA-C	5.61	118.36	110.23
1	A	4	VAL	N-CA-C	5.57	120.93	109.34
1	A	363	GLU	N-CA-C	5.57	118.87	111.75
1	A	112	ASP	N-CA-C	5.55	117.83	109.23
1	A	513	CYS	N-CA-C	5.54	117.76	111.11
1	A	407	MET	N-CA-C	-5.52	102.23	110.23
1	A	122	ASN	N-CA-C	-5.42	99.25	110.80
1	A	172	TYR	N-CA-C	5.40	117.89	109.52
1	A	388	LYS	N-CA-C	5.40	119.39	112.26
1	A	40	ASN	N-CA-C	5.40	119.03	111.74
1	A	565	HIS	O-C-N	-5.34	117.24	123.27
1	A	224	LEU	N-CA-C	-5.25	102.00	110.14
1	A	168	PRO	N-CA-C	5.22	123.22	112.47
1	A	169	THR	CA-C-N	5.18	125.71	120.38
1	A	169	THR	C-N-CA	5.18	125.71	120.38
1	A	170	PRO	CA-C-N	5.17	126.31	119.84
1	A	170	PRO	C-N-CA	5.17	126.31	119.84
1	A	5	GLY	N-CA-C	5.17	125.42	113.18
1	A	167	CYS	CB-CA-C	5.12	120.25	110.17
1	A	157	CYS	N-CA-C	5.11	115.27	108.38
1	A	336	ASP	N-CA-CB	-5.07	102.61	110.07
1	A	386	ILE	N-CA-C	-5.05	105.67	110.42
1	A	169	THR	CA-C-O	-5.04	113.26	120.16

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	169	THR	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	4766	0	4676	183	0
2	B	39	0	34	2	0
2	D	39	0	34	0	0
3	C	28	0	25	1	0
3	E	28	0	25	1	0
4	A	1	0	0	0	0
5	A	4	0	0	0	0
6	A	7	0	0	5	0
7	A	28	0	0	4	0
8	A	43	0	30	6	0
9	A	236	0	0	37	0
All	All	5219	0	4824	189	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 19.

All (189) 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:7:GLY:HA3	9:A:3187:HOH:O	1.51	1.09
1:A:289:GLY:HA3	9:A:3196:HOH:O	1.57	1.04
1:A:8:ALA:HB1	1:A:9:PRO:CD	1.88	1.03
1:A:258:GLU:OE2	8:A:3008:HEM:HMB1	0.85	1.02
1:A:167:CYS:HB3	1:A:168:PRO:HD2	1.43	0.98
9:A:3206:HOH:O	3:E:2:NAG:H5	1.62	0.98
1:A:292:LEU:HG	9:A:3225:HOH:O	1.64	0.97
1:A:8:ALA:O	1:A:10:VAL:HG22	1.65	0.96
1:A:8:ALA:HB1	1:A:9:PRO:HD2	1.45	0.95

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
8:A:3008:HEM:HMC2	8:A:3008:HEM:HBC2	1.51	0.91
1:A:333:ASN:C	1:A:333:ASN:HD22	1.84	0.86
1:A:42:ALA:HB2	1:A:166:VAL:HG21	1.57	0.86
1:A:263:ALA:O	1:A:267:THR:HG22	1.76	0.86
1:A:8:ALA:CB	1:A:9:PRO:HD2	2.04	0.85
1:A:288:ASP:HB2	1:A:291:LYS:HB3	1.56	0.85
1:A:8:ALA:CB	1:A:9:PRO:CD	2.55	0.84
1:A:173:GLN:HG3	1:A:174:SER:H	1.43	0.83
1:A:32:ARG:HG3	9:A:3209:HOH:O	1.78	0.82
1:A:232:ARG:O	1:A:232:ARG:HG3	1.78	0.82
1:A:167:CYS:HB3	1:A:168:PRO:CD	2.09	0.82
1:A:106:ILE:HG23	1:A:191:LEU:HD11	1.63	0.80
1:A:132:TYR:HE1	9:A:3228:HOH:O	1.66	0.78
1:A:254:PHE:HD2	9:A:3160:HOH:O	1.68	0.75
1:A:63:GLN:HG2	9:A:3135:HOH:O	1.87	0.75
1:A:288:ASP:HB2	1:A:291:LYS:CB	2.20	0.72
1:A:360:ARG:HD2	9:A:3200:HOH:O	1.90	0.71
1:A:367:PRO:HG3	6:A:2007:IOD:I	2.60	0.71
1:A:146:LYS:O	1:A:147:ASN:HB2	1.90	0.71
1:A:77:GLU:HG3	7:A:3004:NO3:O2	1.91	0.70
8:A:3008:HEM:HBC2	8:A:3008:HEM:CMC	2.21	0.70
1:A:129:CYS:HB2	9:A:3033:HOH:O	1.92	0.70
1:A:258:GLU:OE2	8:A:3008:HEM:C2B	2.47	0.68
1:A:403:LYS:NZ	9:A:3118:HOH:O	2.26	0.68
1:A:551:ARG:HD3	1:A:584:LYS:HA	1.77	0.67
1:A:171:PRO:HD3	9:A:3084:HOH:O	1.92	0.67
1:A:221:ASP:HB2	1:A:226:TYR:CZ	2.30	0.67
1:A:360:ARG:HA	1:A:394:PRO:O	1.95	0.67
1:A:527:ARG:O	1:A:532:ASN:ND2	2.28	0.66
1:A:169:THR:HB	1:A:170:PRO:HD3	1.78	0.66
1:A:120:GLY:C	1:A:122:ASN:H	2.02	0.65
1:A:2:TRP:CB	1:A:4:VAL:HG22	2.27	0.65
1:A:2:TRP:HB3	1:A:4:VAL:HG22	1.79	0.64
1:A:258:GLU:CD	8:A:3008:HEM:CMB	2.66	0.63
1:A:8:ALA:HB1	1:A:9:PRO:HD3	1.76	0.62
1:A:145:PRO:HG2	1:A:148:ASP:HB2	1.80	0.62
1:A:328:TYR:HD1	1:A:523:ARG:HD3	1.63	0.62
1:A:169:THR:HB	1:A:170:PRO:CD	2.29	0.61
1:A:173:GLN:HB3	9:A:3221:HOH:O	2.00	0.61
1:A:243:THR:HG22	9:A:3215:HOH:O	2.00	0.60
1:A:9:PRO:O	1:A:10:VAL:HG13	2.00	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:407:MET:HG3	9:A:3177:HOH:O	2.01	0.60
1:A:64:ARG:HA	9:A:3219:HOH:O	2.02	0.59
1:A:333:ASN:C	1:A:333:ASN:ND2	2.56	0.59
1:A:547:MET:HE2	1:A:585:LEU:HD22	1.83	0.59
1:A:187:LEU:HD13	1:A:305:GLN:HA	1.85	0.59
1:A:542:ASP:HB3	9:A:3223:HOH:O	2.02	0.59
1:A:544:LEU:HD23	1:A:547:MET:HE3	1.85	0.58
1:A:114:ALA:HB3	9:A:3027:HOH:O	2.03	0.58
1:A:167:CYS:CB	1:A:168:PRO:HD2	2.27	0.58
1:A:393:ASP:OD1	1:A:557:THR:HB	2.05	0.57
1:A:273:HIS:CD2	1:A:273:HIS:C	2.83	0.56
1:A:328:TYR:CD1	1:A:523:ARG:HD3	2.40	0.56
1:A:232:ARG:HG2	9:A:3188:HOH:O	2.05	0.55
1:A:9:PRO:C	1:A:10:VAL:HG13	2.32	0.55
1:A:187:LEU:HD22	1:A:304:VAL:HG12	1.87	0.55
1:A:259:GLN:HE22	1:A:261:LEU:HB2	1.71	0.54
1:A:43:LEU:HD13	1:A:181:ASN:HB2	1.88	0.54
1:A:288:ASP:HB3	1:A:291:LYS:H	1.72	0.54
1:A:393:ASP:HB2	1:A:394:PRO:HD3	1.90	0.54
1:A:232:ARG:HH21	1:A:248:CYS:CB	2.21	0.53
1:A:544:LEU:HD23	1:A:547:MET:CE	2.38	0.53
1:A:336:ASP:OD2	1:A:338:ARG:NH2	2.33	0.53
2:B:2:NAG:H3	2:B:3:MAN:H61	1.91	0.53
1:A:352:MET:HB2	1:A:407:MET:HG2	1.91	0.52
1:A:167:CYS:CB	1:A:168:PRO:CD	2.81	0.52
9:A:3151:HOH:O	2:B:3:MAN:H3	2.09	0.52
1:A:45:ARG:CZ	1:A:49:ALA:HB2	2.40	0.52
1:A:173:GLN:HG3	1:A:174:SER:N	2.20	0.51
1:A:481:MET:HA	1:A:481:MET:HE2	1.92	0.51
1:A:286:GLN:H	1:A:286:GLN:CD	2.19	0.51
1:A:76:ARG:NE	1:A:148:ASP:OD2	2.43	0.51
1:A:504:ARG:HD2	9:A:3210:HOH:O	2.09	0.51
1:A:166:VAL:C	1:A:167:CYS:O	2.53	0.51
1:A:10:VAL:HG12	1:A:40:ASN:O	2.11	0.51
1:A:66:THR:HB	1:A:70:PHE:N	2.26	0.51
1:A:101:MET:O	1:A:101:MET:HG2	2.09	0.51
1:A:216:ASN:HA	6:A:2001:IOD:I	2.80	0.51
1:A:64:ARG:HB3	9:A:3024:HOH:O	2.11	0.50
1:A:2:TRP:HB2	1:A:4:VAL:HG22	1.93	0.50
1:A:77:GLU:OE2	1:A:81:LYS:NZ	2.36	0.50
1:A:467:LEU:HD22	1:A:481:MET:HE1	1.93	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:239:PHE:CZ	1:A:427:LYS:HG2	2.46	0.50
1:A:173:GLN:CG	1:A:174:SER:H	2.09	0.49
8:A:3008:HEM:HMC2	8:A:3008:HEM:CBC	2.35	0.49
1:A:144:PHE:CE2	1:A:158:MET:HG3	2.48	0.49
1:A:538:GLU:HG3	1:A:541:ARG:NH2	2.28	0.49
1:A:42:ALA:HB2	1:A:166:VAL:CG2	2.38	0.49
1:A:481:MET:HE2	1:A:481:MET:CA	2.43	0.48
1:A:43:LEU:HB2	9:A:3045:HOH:O	2.11	0.48
1:A:198:SER:HB2	7:A:3003:NO3:O2	2.13	0.48
1:A:511:LEU:HB2	9:A:3101:HOH:O	2.14	0.48
1:A:43:LEU:HG	1:A:341:ASN:HA	1.96	0.48
1:A:165:PHE:HE2	1:A:172:TYR:HB3	1.78	0.48
1:A:170:PRO:HB3	1:A:171:PRO:HD2	1.95	0.47
1:A:544:LEU:CD2	1:A:547:MET:HE1	2.43	0.47
1:A:393:ASP:HB2	1:A:394:PRO:CD	2.44	0.47
1:A:424:PRO:O	6:A:2003:IOD:I	3.01	0.47
1:A:216:ASN:HB2	1:A:227:LEU:O	2.14	0.47
1:A:9:PRO:O	1:A:10:VAL:CG1	2.63	0.47
1:A:217:GLN:HE21	3:C:1:NAG:C7	2.28	0.46
1:A:453:ARG:NH1	1:A:499:GLU:OE2	2.46	0.46
1:A:19:SER:HA	1:A:20:PRO:HD3	1.81	0.46
1:A:551:ARG:CZ	1:A:584:LYS:HG2	2.45	0.46
1:A:336:ASP:CG	1:A:338:ARG:HH21	2.22	0.46
1:A:101:MET:CE	1:A:102:GLN:HA	2.46	0.46
1:A:123:GLU:HG3	1:A:124:HIS:N	2.31	0.46
1:A:237:CYS:HA	1:A:381:PHE:O	2.15	0.46
1:A:272:GLU:OE1	1:A:272:GLU:HA	2.15	0.46
1:A:3:GLU:O	1:A:5:GLY:N	2.49	0.46
1:A:132:TYR:CE1	9:A:3228:HOH:O	2.53	0.46
1:A:259:GLN:HE21	1:A:259:GLN:HB2	1.46	0.46
1:A:446:MET:HA	1:A:446:MET:HE2	1.98	0.45
1:A:91:VAL:HG13	1:A:405:LYS:HG3	1.98	0.45
1:A:130:GLU:OE2	1:A:426:HIS:HB3	2.17	0.45
1:A:15:CYS:O	1:A:16:ASP:C	2.59	0.45
1:A:388:LYS:HA	9:A:3216:HOH:O	2.16	0.45
1:A:119:LEU:HD22	9:A:3084:HOH:O	2.16	0.45
1:A:432:ASP:OD1	1:A:432:ASP:C	2.60	0.45
1:A:67:ARG:NH1	9:A:3034:HOH:O	2.47	0.45
1:A:377:HIS:HA	1:A:380:PHE:CE2	2.52	0.45
1:A:407:MET:HB3	1:A:501:MET:HE3	1.99	0.45
1:A:568:GLN:HE21	1:A:570:ASN:HD21	1.64	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:254:PHE:CD2	9:A:3160:HOH:O	2.55	0.45
1:A:364:ASN:O	1:A:365:TYR:HB2	2.17	0.44
1:A:368:TRP:CH2	1:A:389:ASP:O	2.70	0.44
1:A:5:GLY:O	1:A:6:CYS:C	2.60	0.44
1:A:202:ARG:HD3	1:A:250:LEU:HD21	1.98	0.44
1:A:148:ASP:O	1:A:149:PRO:C	2.59	0.44
1:A:476:LEU:HD21	1:A:498:ALA:HB1	1.99	0.44
1:A:360:ARG:NH2	1:A:371:GLU:O	2.51	0.44
1:A:300:LEU:O	1:A:303:PHE:HB3	2.18	0.43
1:A:563:PRO:HD3	1:A:576:PHE:CE2	2.53	0.43
1:A:2:TRP:C	1:A:4:VAL:H	2.24	0.43
1:A:543:SER:OG	1:A:586:ASP:O	2.36	0.43
1:A:71:ARG:CZ	9:A:3219:HOH:O	2.66	0.43
1:A:74:LEU:O	1:A:75:ALA:C	2.61	0.43
1:A:171:PRO:HG3	9:A:3242:HOH:O	2.17	0.43
1:A:233:LYS:HA	1:A:234:PRO:C	2.42	0.43
1:A:288:ASP:CB	1:A:291:LYS:H	2.31	0.43
1:A:313:LEU:HD11	1:A:519:PHE:CG	2.54	0.43
1:A:544:LEU:CD2	1:A:547:MET:CE	2.96	0.43
1:A:113:PHE:C	1:A:115:PRO:HD3	2.44	0.43
1:A:258:GLU:HG3	6:A:2002:IOD:I	2.89	0.43
1:A:561:LYS:HE2	9:A:3197:HOH:O	2.18	0.43
1:A:279:GLU:OE1	1:A:279:GLU:HA	2.18	0.43
1:A:407:MET:HB3	1:A:407:MET:HE3	1.88	0.42
1:A:13:VAL:HB	1:A:14:LYS:H	1.71	0.42
1:A:232:ARG:CG	9:A:3188:HOH:O	2.67	0.42
1:A:348:ARG:NH2	9:A:3179:HOH:O	2.51	0.42
1:A:452:TRP:HB3	1:A:510:LEU:HD11	2.01	0.42
1:A:167:CYS:O	1:A:168:PRO:O	2.37	0.42
1:A:258:GLU:O	1:A:380:PHE:HA	2.20	0.42
1:A:499:GLU:OE1	1:A:509:PRO:HD2	2.20	0.42
1:A:362:ASP:OD1	1:A:362:ASP:C	2.63	0.42
1:A:378:THR:HG22	7:A:3007:NO3:O1	2.19	0.42
1:A:432:ASP:OD1	1:A:434:ALA:N	2.53	0.42
1:A:550:SER:O	1:A:551:ARG:C	2.62	0.42
1:A:166:VAL:O	1:A:167:CYS:C	2.61	0.42
1:A:377:HIS:NE2	7:A:3007:NO3:O1	2.52	0.42
1:A:101:MET:HE2	1:A:102:GLN:N	2.35	0.42
1:A:561:LYS:CE	9:A:3197:HOH:O	2.67	0.42
1:A:229:PHE:CG	1:A:247:PRO:HG2	2.55	0.42
1:A:235:SER:HA	1:A:236:PRO:HD3	1.89	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:125:SER:HB3	9:A:3226:HOH:O	2.20	0.41
1:A:309:PHE:CD1	1:A:529:TRP:HH2	2.39	0.41
1:A:146:LYS:O	1:A:147:ASN:CB	2.57	0.41
1:A:169:THR:CB	1:A:170:PRO:CD	2.98	0.41
1:A:236:PRO:HG3	6:A:2003:IOD:I	2.90	0.41
1:A:320:GLU:HG3	1:A:502:VAL:HG11	2.03	0.41
1:A:370:PRO:HD2	1:A:371:GLU:H	1.86	0.41
1:A:417:LEU:HD13	1:A:433:LEU:HD23	2.03	0.41
1:A:338:ARG:HE	1:A:338:ARG:HB2	1.67	0.40
1:A:101:MET:HE2	1:A:102:GLN:CA	2.51	0.40
1:A:232:ARG:HH21	1:A:248:CYS:HB2	1.87	0.40
1:A:528:PHE:O	1:A:529:TRP:C	2.64	0.40
1:A:449:TYR:CZ	1:A:453:ARG:HD3	2.56	0.40
1:A:230:ASN:HB3	9:A:3188:HOH:O	2.21	0.40
1:A:393:ASP:C	1:A:395:LEU:H	2.29	0.40
1:A:393:ASP:CB	1:A:394:PRO:CD	3.00	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
1	A	593/595 (100%)	527 (89%)	48 (8%)	18 (3%)	3 12

All (18) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	4	VAL
1	A	8	ALA
1	A	122	ASN
1	A	167	CYS
1	A	168	PRO

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Mol	Chain	Res	Type
1	A	169	THR
1	A	170	PRO
1	A	174	SER
1	A	6	CYS
1	A	9	PRO
1	A	17	GLU
1	A	56	ALA
1	A	150	LYS
1	A	63	GLN
1	A	166	VAL
1	A	381	PHE
1	A	13	VAL
1	A	120	GLY

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%)	456 (88%)	61 (12%)	5 17

All (61) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	2	TRP
1	A	4	VAL
1	A	6	CYS
1	A	9	PRO
1	A	12	LEU
1	A	13	VAL
1	A	17	GLU
1	A	18	ASN
1	A	24	ILE
1	A	43	LEU
1	A	57	LEU
1	A	72	VAL
1	A	81	LYS

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Mol	Chain	Res	Type
1	A	82	ILE
1	A	91	VAL
1	A	121	SER
1	A	146	LYS
1	A	166	VAL
1	A	168	PRO
1	A	170	PRO
1	A	175	LEU
1	A	201	SER
1	A	203	LEU
1	A	210	LEU
1	A	232	ARG
1	A	240	ILE
1	A	242	THR
1	A	257	SER
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	338	ARG
1	A	360	ARG
1	A	392	ILE
1	A	407	MET
1	A	420	LYS
1	A	427	LYS
1	A	428	ILE
1	A	441	CYS
1	A	464	LEU
1	A	471	LEU
1	A	475	ILE
1	A	478	LYS
1	A	480	LEU
1	A	481	MET
1	A	504	ARG
1	A	511	LEU

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Mol	Chain	Res	Type
1	A	517	ARG
1	A	523	ARG
1	A	543	SER
1	A	564	LEU
1	A	583	ASP
1	A	592	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (15) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	40	ASN
1	A	68	ASN
1	A	204	GLN
1	A	217	GLN
1	A	259	GLN
1	A	286	GLN
1	A	333	ASN
1	A	460	GLN
1	A	468	GLN
1	A	497	ASN
1	A	520	GLN
1	A	545	GLN
1	A	558	HIS
1	A	565	HIS
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 ⓘ

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

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	0.73	0	17,19,21	2.13	6 (35%)
2	NAG	B	2	2	14,14,15	0.77	1 (7%)	17,19,21	1.41	1 (5%)
2	MAN	B	3	2	11,11,12	0.57	0	15,15,17	1.00	1 (6%)
3	NAG	C	1	3,1	14,14,15	0.66	1 (7%)	17,19,21	1.29	2 (11%)
3	NAG	C	2	3	14,14,15	0.92	0	17,19,21	2.12	4 (23%)
2	NAG	D	1	2,1	14,14,15	0.68	0	17,19,21	2.18	8 (47%)
2	NAG	D	2	2	14,14,15	0.49	0	17,19,21	2.22	4 (23%)
2	MAN	D	3	2	11,11,12	0.84	0	15,15,17	1.82	2 (13%)
3	NAG	E	1	3,1	14,14,15	0.83	0	17,19,21	1.66	3 (17%)
3	NAG	E	2	3	14,14,15	0.72	0	17,19,21	1.21	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	-	4/6/23/26	0/1/1/1
2	NAG	B	2	2	-	0/6/23/26	0/1/1/1
2	MAN	B	3	2	-	2/2/19/22	1/1/1/1
3	NAG	C	1	3,1	-	0/6/23/26	0/1/1/1
3	NAG	C	2	3	-	2/6/23/26	0/1/1/1
2	NAG	D	1	2,1	-	0/6/23/26	0/1/1/1
2	NAG	D	2	2	-	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	-	2/6/23/26	0/1/1/1
3	NAG	E	2	3	-	2/6/23/26	0/1/1/1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	2	NAG	O5-C1	-2.17	1.40	1.43
3	C	1	NAG	O5-C1	-2.02	1.40	1.43

All (33) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	2	NAG	C1-O5-C5	7.18	121.80	112.19
3	C	2	NAG	C1-O5-C5	6.76	121.25	112.19
2	D	3	MAN	C1-O5-C5	5.59	119.68	112.19
2	B	1	NAG	O5-C1-C2	-5.53	102.73	111.29
3	E	1	NAG	C4-C3-C2	4.22	117.21	111.02
2	D	1	NAG	C1-O5-C5	4.05	117.61	112.19
2	B	2	NAG	C4-C3-C2	3.98	116.85	111.02
2	D	1	NAG	C1-C2-N2	3.56	116.05	110.43
2	B	1	NAG	C2-N2-C7	-3.48	118.24	122.90
3	C	1	NAG	C1-O5-C5	3.36	116.69	112.19
3	E	2	NAG	C1-O5-C5	-3.08	108.06	112.19
2	B	1	NAG	C1-O5-C5	3.07	116.30	112.19
2	D	1	NAG	O7-C7-N2	2.91	127.12	121.98
2	B	1	NAG	C4-C3-C2	2.81	115.14	111.02
2	D	3	MAN	C1-C2-C3	2.73	113.62	109.64
2	D	1	NAG	O5-C1-C2	-2.72	107.08	111.29
2	B	3	MAN	O5-C5-C6	2.71	112.94	107.66
2	D	1	NAG	C3-C4-C5	2.66	115.06	110.23
2	D	1	NAG	C6-C5-C4	-2.66	106.48	113.02
3	E	1	NAG	O5-C1-C2	-2.64	107.21	111.29
3	E	1	NAG	C2-N2-C7	-2.57	119.46	122.90
3	C	1	NAG	O7-C7-C8	-2.47	117.65	122.05
3	C	2	NAG	O5-C5-C4	2.47	116.83	110.83
2	D	2	NAG	C4-C3-C2	-2.46	107.42	111.02
3	E	2	NAG	O5-C5-C6	2.44	112.41	107.66
2	D	1	NAG	O7-C7-C8	-2.43	117.73	122.05
2	D	2	NAG	C1-C2-N2	2.43	114.26	110.43
3	C	2	NAG	O3-C3-C2	2.40	114.39	109.40
2	D	2	NAG	C2-N2-C7	2.27	125.94	122.90
3	C	2	NAG	O5-C1-C2	2.24	114.75	111.29
2	D	1	NAG	C4-C3-C2	2.06	114.04	111.02
2	B	1	NAG	O6-C6-C5	2.04	118.30	111.33
2	B	1	NAG	O4-C4-C5	2.00	114.25	109.32

There are no chirality outliers.

All (15) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	1	NAG	O5-C5-C6-O6
2	B	1	NAG	C4-C5-C6-O6
2	B	1	NAG	C8-C7-N2-C2

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

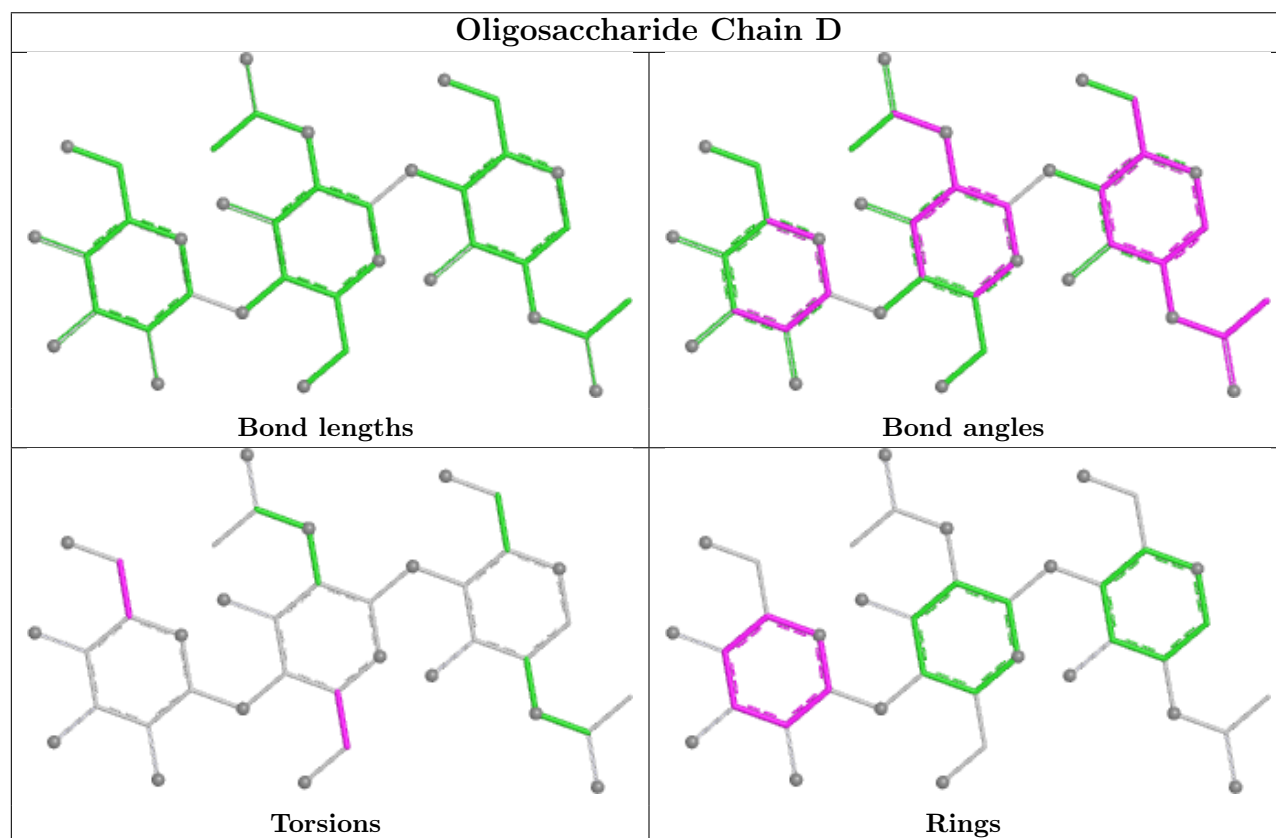
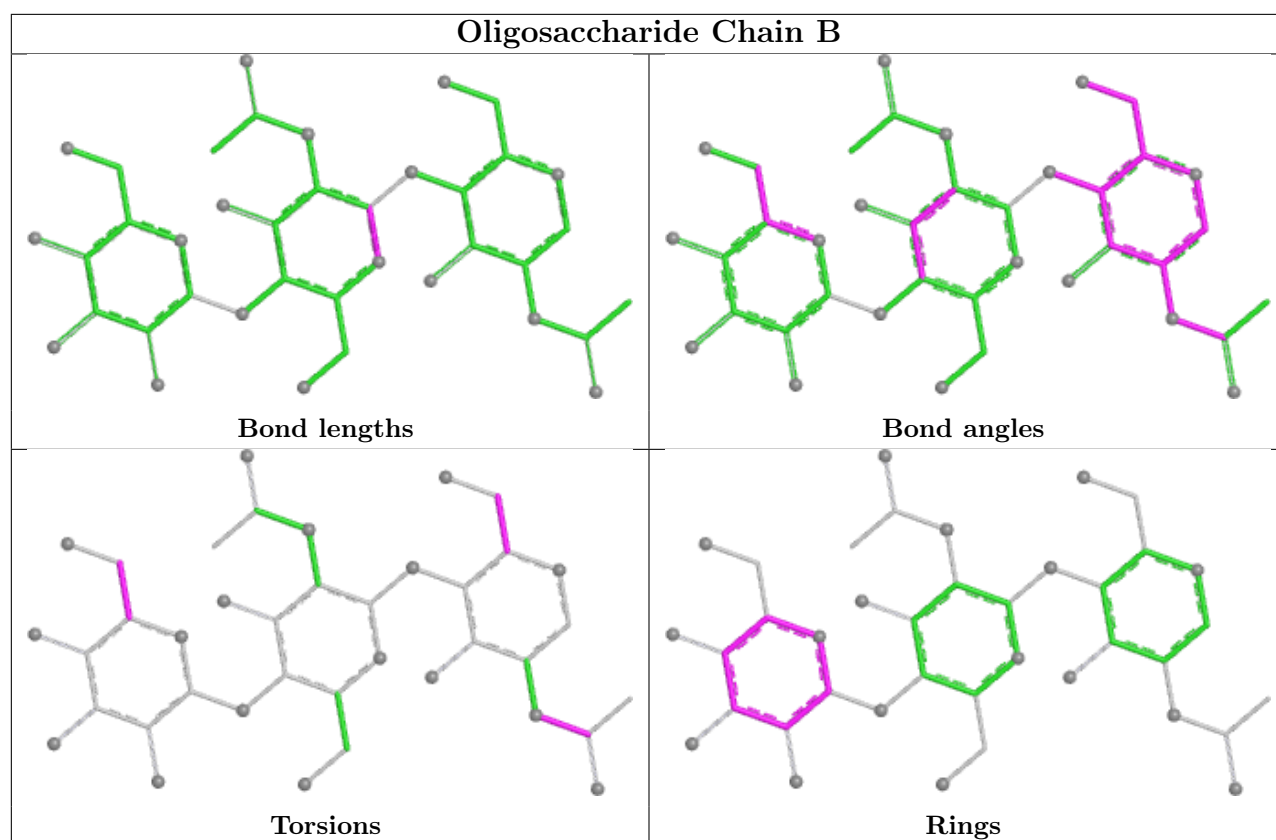
All (2) ring outliers are listed below:

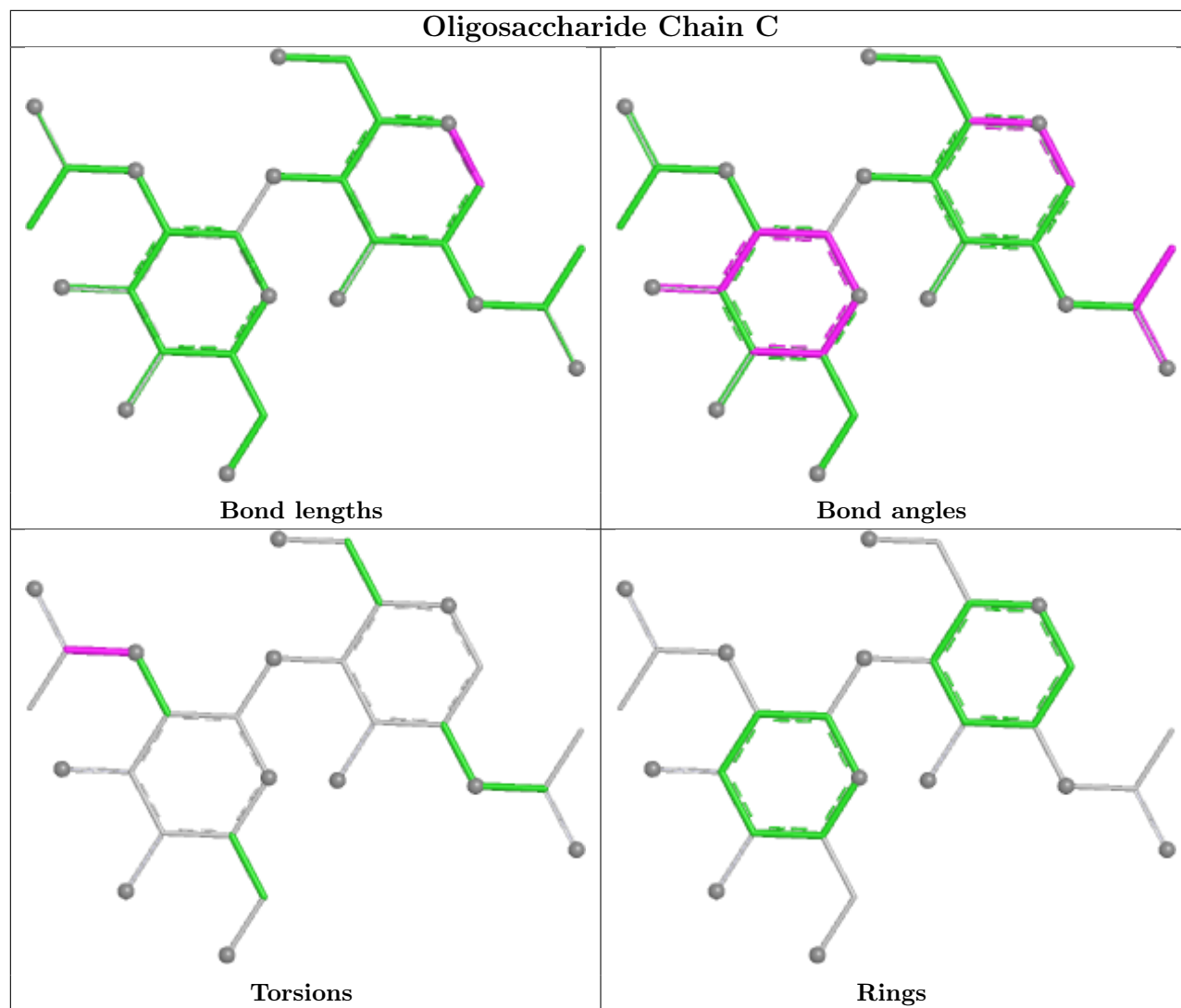
Mol	Chain	Res	Type	Atoms
2	D	3	MAN	C1-C2-C3-C4-C5-O5
2	B	3	MAN	C1-C2-C3-C4-C5-O5

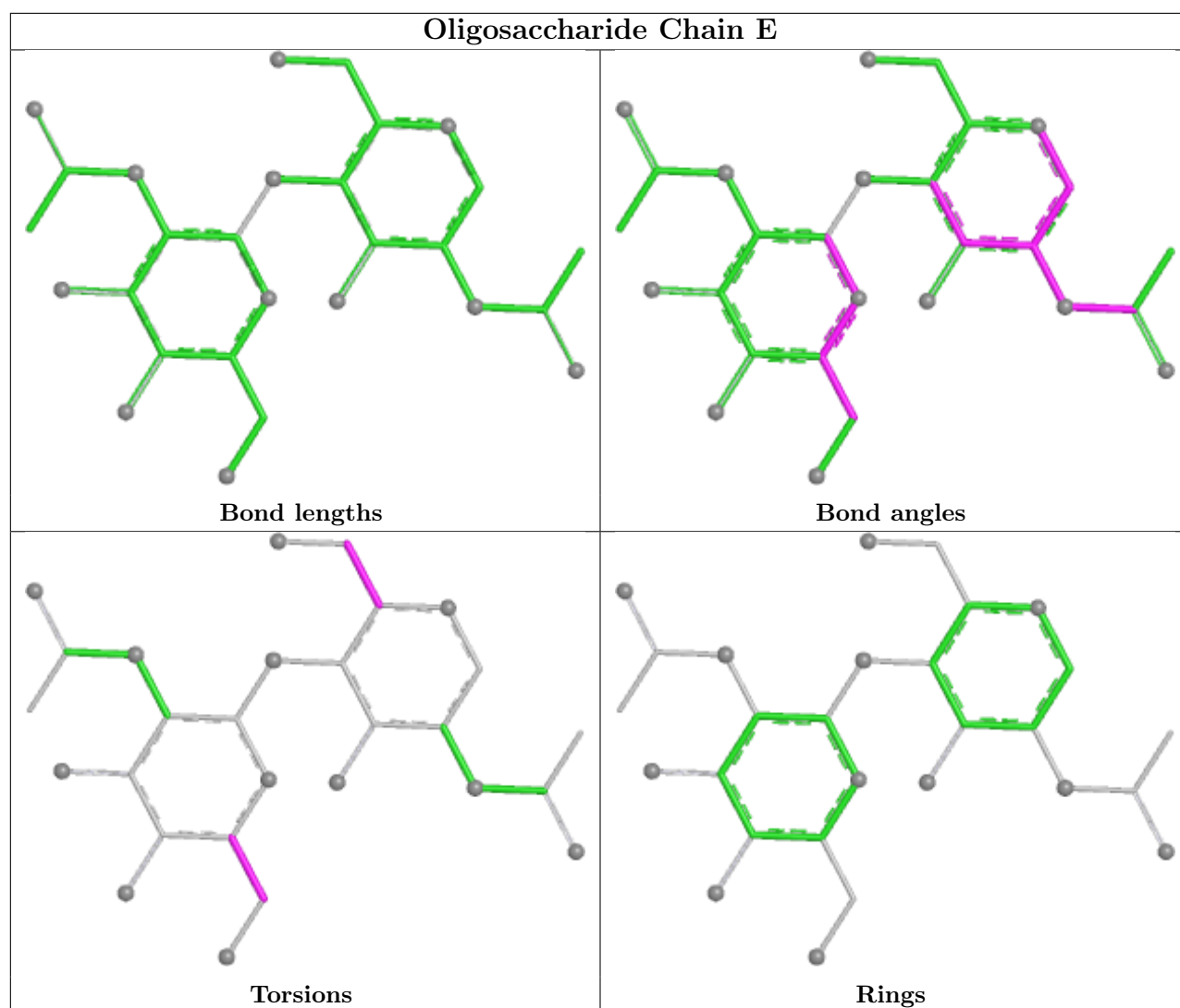
4 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	E	2	NAG	1	0
3	C	1	NAG	1	0
2	B	2	NAG	1	0
2	B	3	MAN	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 17 ligands modelled in this entry, 8 are monoatomic - leaving 9 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	NO3	A	3005	-	1,3,3	2.18	1 (100%)	0,3,3	-	-
5	CO3	A	688	-	3,3,3	0.84	0	2,3,3	2.75	2 (100%)
8	HEM	A	3008	1	50,50,50	2.01	10 (20%)	67,82,82	1.38	9 (13%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
7	NO3	A	3001	-	1,3,3	2.67	1 (100%)	0,3,3	-	-
7	NO3	A	3003	-	1,3,3	4.37	1 (100%)	0,3,3	-	-
7	NO3	A	3006	-	1,3,3	2.62	1 (100%)	0,3,3	-	-
7	NO3	A	3002	-	1,3,3	3.31	1 (100%)	0,3,3	-	-
7	NO3	A	3007	-	1,3,3	2.82	1 (100%)	0,3,3	-	-
7	NO3	A	3004	-	1,3,3	3.32	1 (100%)	0,3,3	-	-

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	3008	1	-	4/14/54/54	-

All (17) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
8	A	3008	HEM	C3D-C2D	7.83	1.53	1.36
8	A	3008	HEM	FE-NC	5.57	2.13	1.95
7	A	3003	NO3	O1-N	4.37	1.45	1.24
7	A	3004	NO3	O1-N	3.32	1.40	1.24
7	A	3002	NO3	O1-N	3.31	1.40	1.24
8	A	3008	HEM	CAB-C3B	3.24	1.56	1.47
8	A	3008	HEM	FE-NA	3.18	2.05	1.95
8	A	3008	HEM	CAC-C3C	3.13	1.55	1.47
7	A	3007	NO3	O1-N	2.82	1.38	1.24
8	A	3008	HEM	FE-ND	2.77	2.03	1.94
7	A	3001	NO3	O1-N	2.67	1.37	1.24
7	A	3006	NO3	O1-N	2.62	1.37	1.24
8	A	3008	HEM	FE-NB	2.42	2.02	1.94
8	A	3008	HEM	CAD-C3D	2.21	1.57	1.51
7	A	3005	NO3	O1-N	2.18	1.35	1.24
8	A	3008	HEM	CAA-C2A	2.13	1.56	1.51
8	A	3008	HEM	CMA-C3A	2.13	1.55	1.50

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
8	A	3008	HEM	C4D-ND-C1D	4.39	110.41	105.21

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
8	A	3008	HEM	C3B-C2B-C1B	3.03	108.69	106.41
8	A	3008	HEM	CHC-C4B-NB	2.94	127.59	124.42
5	A	688	CO3	O2-C-O1	-2.80	112.51	119.68
5	A	688	CO3	O3-C-O1	2.70	126.58	119.68
8	A	3008	HEM	CHD-C1D-ND	2.48	127.09	124.42
8	A	3008	HEM	CMB-C2B-C3B	-2.36	122.72	128.43
8	A	3008	HEM	CAD-C3D-C4D	2.32	128.75	124.70
8	A	3008	HEM	CMB-C2B-C1B	2.27	128.59	125.03
8	A	3008	HEM	C2B-C1B-NB	-2.14	107.38	109.84
8	A	3008	HEM	C1B-NB-C4B	2.13	107.73	105.21

There are no chirality outliers.

All (4) torsion outliers are listed below:

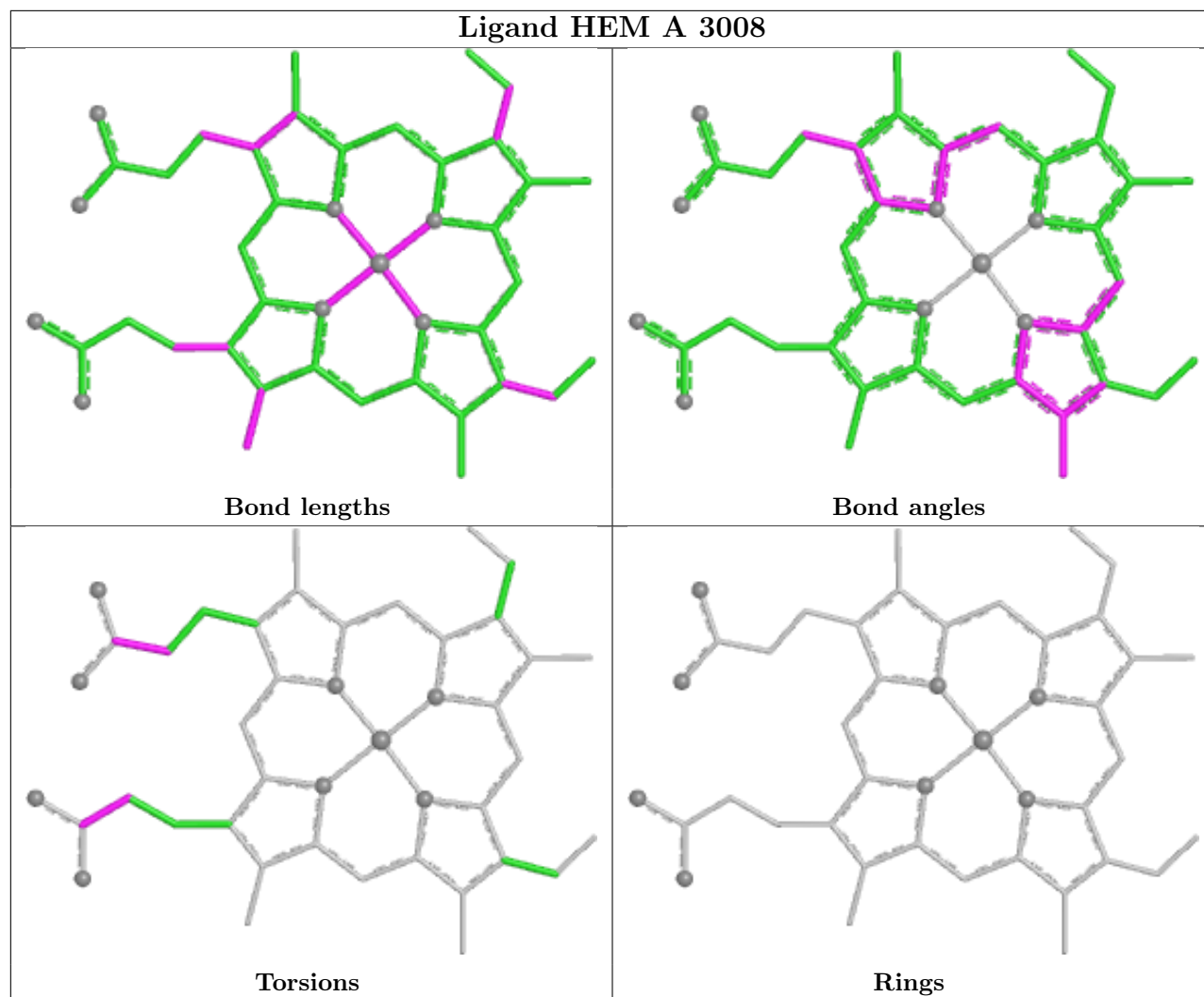
Mol	Chain	Res	Type	Atoms
8	A	3008	HEM	CAA-CBA-CGA-O2A
8	A	3008	HEM	CAD-CBD-CGD-O1D
8	A	3008	HEM	CAD-CBD-CGD-O2D
8	A	3008	HEM	CAA-CBA-CGA-O1A

There are no ring outliers.

4 monomers are involved in 10 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
8	A	3008	HEM	6	0
7	A	3003	NO3	1	0
7	A	3007	NO3	2	0
7	A	3004	NO3	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 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.