



Full wwPDB X-ray Structure Validation Report ⓘ

Nov 6, 2023 – 03:08 pm GMT

PDB ID : 4UZI
Title : Crystal Structure of AauDyP Complexed with Imidazole
Authors : Strittmatter, E.; Liers, C.; Ullrich, R.; Hofrichter, M.; Piontek, K.; Plattner, D.A.
Deposited on : 2014-09-05
Resolution : 2.10 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 : 4.02b-467
Mogul : 1.8.4, CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.36
buster-report : 1.1.7 (2018)
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.36

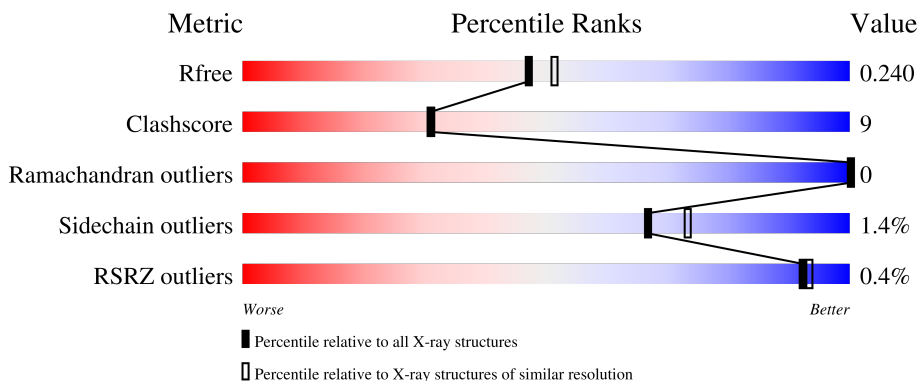
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.10 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}	130704	5197 (2.10-2.10)
Clashscore	141614	5710 (2.10-2.10)
Ramachandran outliers	138981	5647 (2.10-2.10)
Sidechain outliers	138945	5648 (2.10-2.10)
RSRZ outliers	127900	5083 (2.10-2.10)

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	446	 86% 13%
1	B	446	 84% 16%
2	C	2	 50% 50%
3	D	3	 33% 67%

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
10	EPE	A	1460[A]	-	-	X	-
5	NAG	A	1453	X	-	-	-
8	MPD	B	1461	-	-	X	-

2 Entry composition [i](#)

There are 15 unique types of molecules in this entry. The entry contains 7583 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 DYE-DECOLORIZING PEROXIDASE.

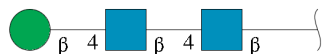
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	446	Total	C	N	O	S	0	4	0
			3333	2096	576	656	5			
1	B	446	Total	C	N	O	S	0	1	0
			3314	2085	573	651	5			

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



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

- Molecule 3 is an oligosaccharide called beta-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
			Total	C	N	O			
3	D	3	Total	C	N	O	0	0	0
			39	22	2	15			

- Molecule 4 is PROTOPORPHYRIN IX CONTAINING FE (three-letter code: HEM) (formula: C₃₄H₃₂FeN₄O₄).



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

- Molecule 5 is 2-acetamido-2-deoxy-beta-D-glucopyranose (three-letter code: NAG) (formula: $C_8H_{15}NO_6$).



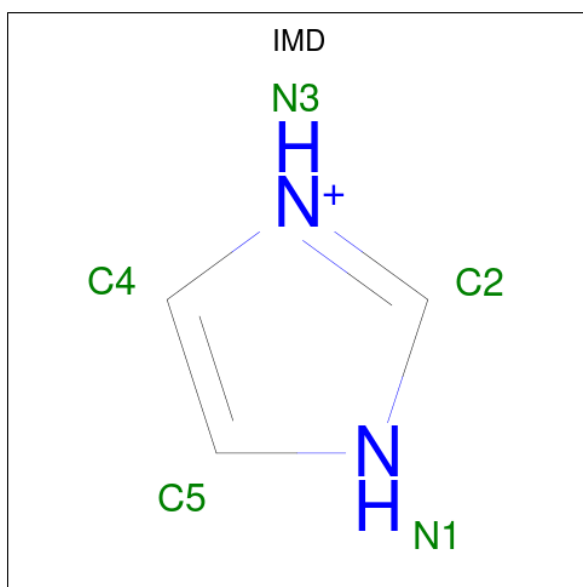
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	
5	A	1	Total	C	N	O	0	0
			14	8	1	5		

Continued on next page...

Continued from previous page...

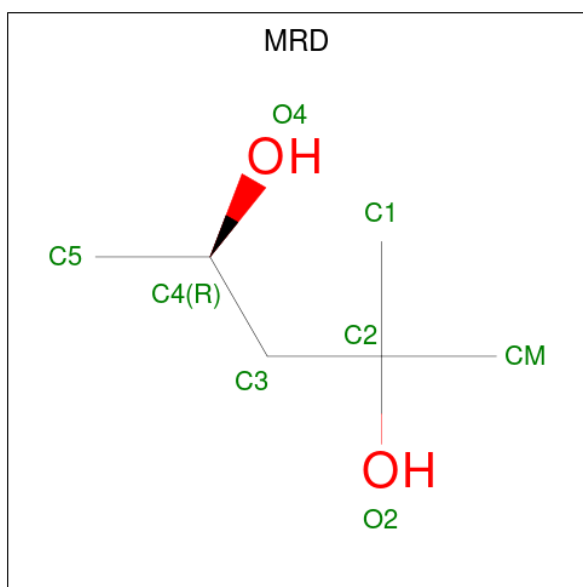
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	A	1	Total	C	N	O	0	0
			14	8	1	5		
5	B	1	Total	C	N	O	0	0
			14	8	1	5		
5	B	1	Total	C	N	O	0	0
			14	8	1	5		

- Molecule 6 is IMIDAZOLE (three-letter code: IMD) (formula: C₃H₅N₂).



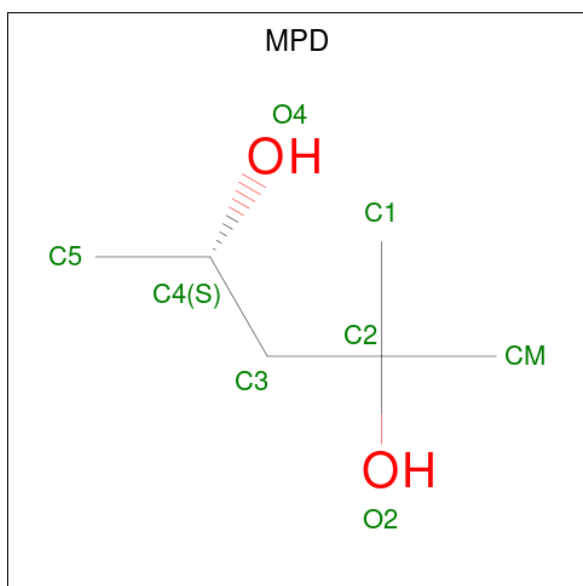
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	A	1	Total	C	N	0	1
			10	6	4		
6	A	1	Total	C	N	0	0
			5	3	2		
6	B	1	Total	C	N	0	1
			10	6	4		
6	B	1	Total	C	N	0	0
			5	3	2		

- Molecule 7 is (4R)-2-METHYLPENTANE-2,4-DIOL (three-letter code: MRD) (formula: C₆H₁₄O₂).



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

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



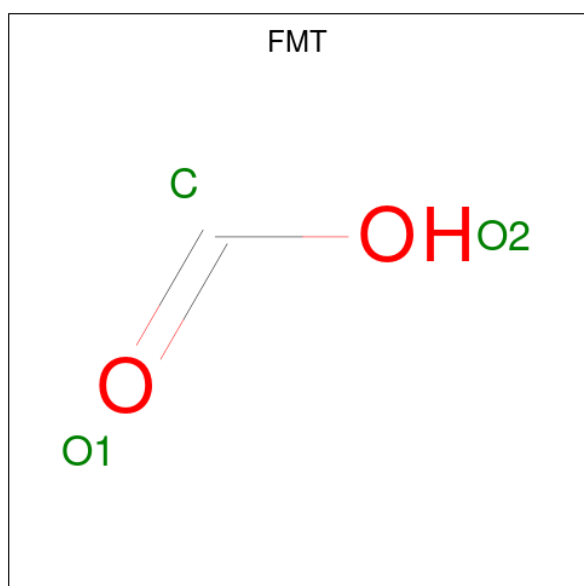
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
8	A	1	Total C O 8 6 2	0	0
8	A	1	Total C O 8 6 2	0	0

Continued on next page...

Continued from previous page...

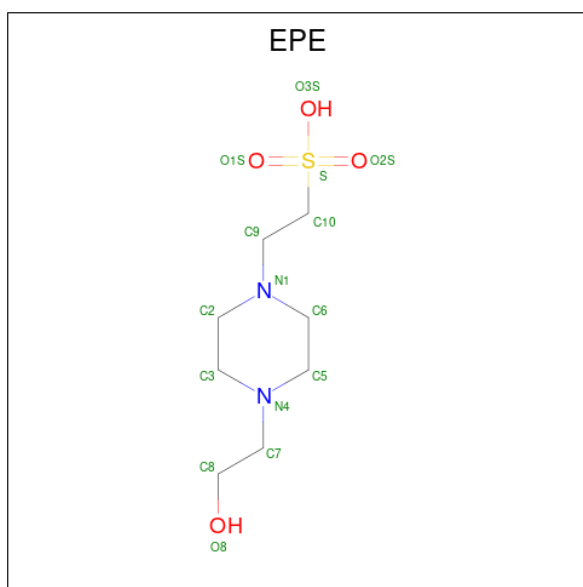
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
8	B	1	Total	C	O	0	0
			8	6	2		
8	B	1	Total	C	O	0	0
			8	6	2		
8	B	1	Total	C	O	0	0
			8	6	2		
8	B	1	Total	C	O	0	0
			8	6	2		

- Molecule 9 is FORMIC ACID (three-letter code: FMT) (formula: CH₂O₂).



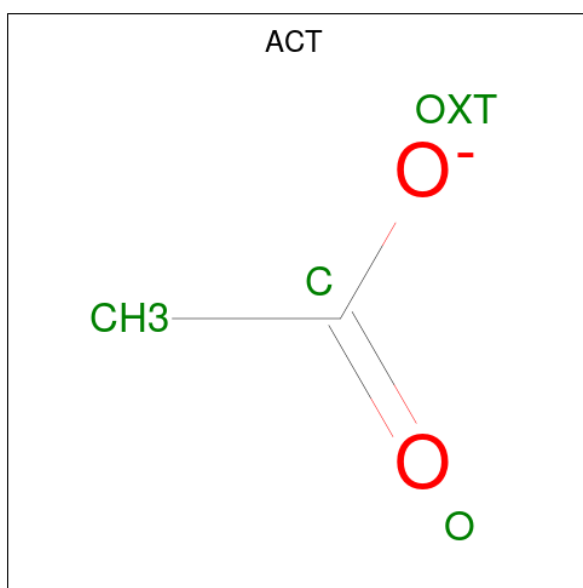
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
9	A	1	Total	C	O	0	0
			3	1	2		

- Molecule 10 is 4-(2-HYDROXYETHYL)-1-PIPERAZINE ETHANESULFONIC ACID (three-letter code: EPE) (formula: C₈H₁₈N₂O₄S).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	
			Total	C	N	O	S			
10	A	1	Total	30	16	4	8	2	0	1
10	B	1	Total	30	16	4	8	2	0	1

- Molecule 11 is ACETATE ION (three-letter code: ACT) (formula: $C_2H_3O_2$).



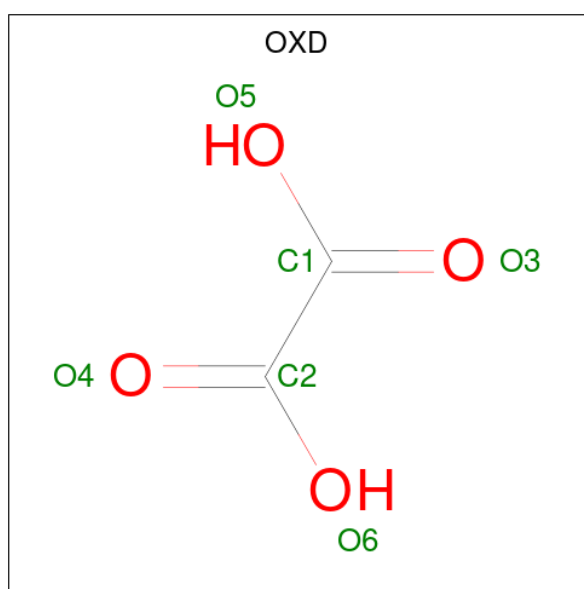
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	
			Total	C	O			
11	A	1	Total	4	2	2	0	0
11	A	1	Total	4	2	2	0	0

Continued on next page...

Continued from previous page...

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

- Molecule 12 is OXALIC ACID (three-letter code: OXD) (formula: C₂H₂O₄).

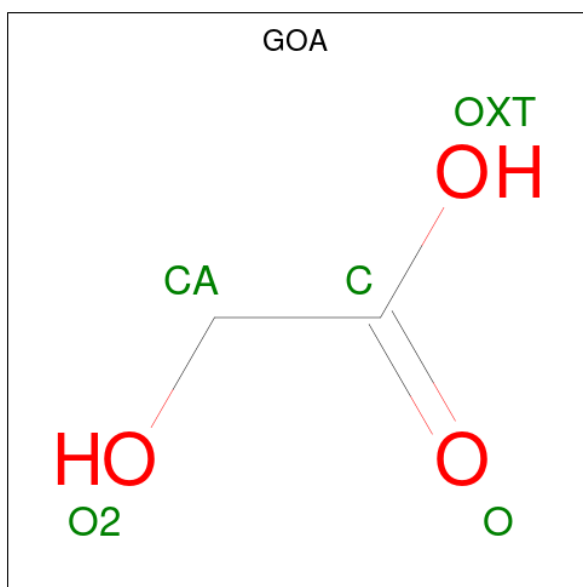


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
12	A	1	Total	C	O	0	0
			6	2	4		

- Molecule 13 is POTASSIUM ION (three-letter code: K) (formula: K).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
13	A	1	Total	K	0	0
			1	1		
13	B	1	Total	K	0	0
			1	1		

- Molecule 14 is GLYCOLIC ACID (three-letter code: GOA) (formula: C₂H₄O₃).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
14	B	1	Total C O 5 2 3	0	0

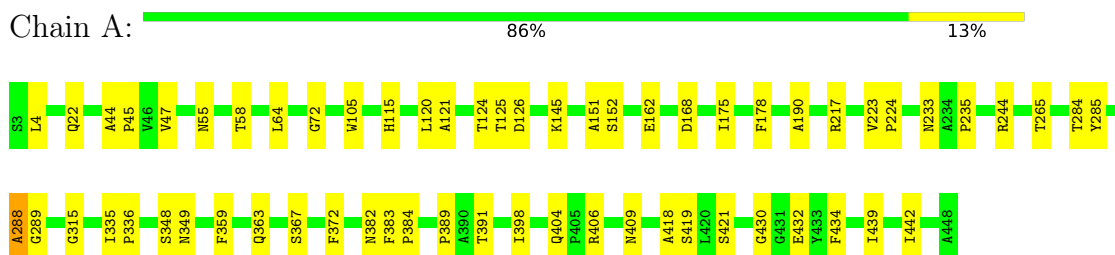
- Molecule 15 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
15	A	317	Total O 318 318	0	1
15	B	215	Total O 215 215	0	0

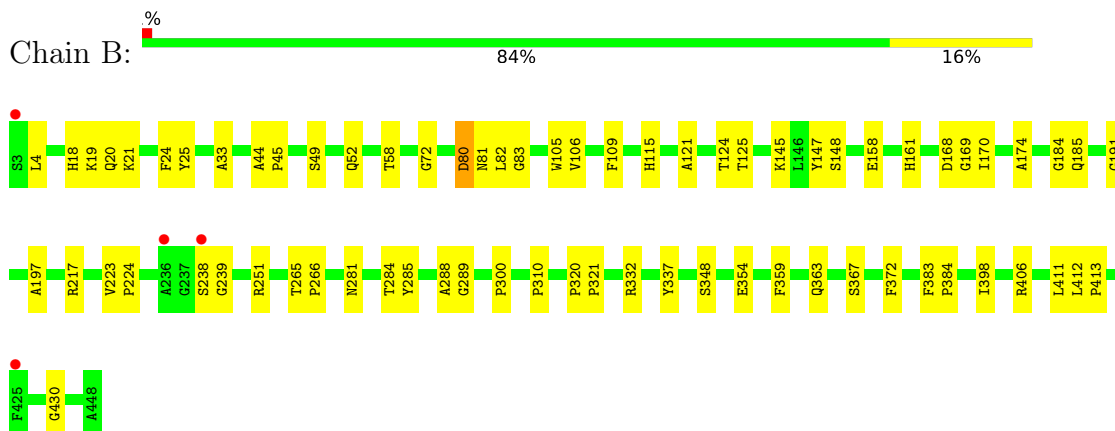
3 Residue-property plots

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 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: DYE-DECOLORIZING PEROXIDASE



- Molecule 1: DYE-DECOLORIZING PEROXIDASE



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



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





4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	67.33Å 45.71Å 140.47Å 90.00° 91.61° 90.00°	Depositor
Resolution (Å)	47.51 – 2.10 47.92 – 2.10	Depositor EDS
% Data completeness (in resolution range)	99.2 (47.51-2.10) 99.3 (47.92-2.10)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.96 (at 2.10Å)	Xtrriage
Refinement program	REFMAC 5.8.0073	Depositor
R, R_{free}	0.179 , 0.239 0.180 , 0.240	Depositor DCC
R_{free} test set	2523 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	20.9	Xtrriage
Anisotropy	0.156	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 43.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtrriage
Estimated twinning fraction	0.033 for h,-k,-l	Xtrriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	7583	wwPDB-VP
Average B, all atoms (Å ²)	22.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.13% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: HEM, GOA, K, IMD, MRD, MPD, EPE, NAG, OXD, ACT, BMA, FMT

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	0.27	0/3411	0.42	0/4653
1	B	0.25	0/3392	0.40	0/4627
All	All	0.26	0/6803	0.41	0/9280

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

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	288	ALA	Peptide

5.2 Too-close contacts [i](#)

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	3333	0	3229	51	0
1	B	3314	0	3213	54	0
2	C	28	0	25	0	0
3	D	39	0	34	0	0
4	A	43	0	30	1	0
4	B	43	0	30	2	0
5	A	28	0	26	1	0
5	B	28	0	26	1	0
6	A	15	0	13	4	0
6	B	15	0	14	4	0
7	A	8	0	14	0	0
8	A	16	0	28	5	0
8	B	40	0	70	11	0
9	A	3	0	1	1	0
10	A	30	0	36	14	0
10	B	30	0	36	7	0
11	A	12	0	9	0	0
11	B	12	0	9	0	0
12	A	6	0	0	0	0
13	A	1	0	0	0	0
13	B	1	0	0	0	0
14	B	5	0	0	0	0
15	A	318	0	0	5	0
15	B	215	0	0	7	0
All	All	7583	0	6843	127	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 9.

All (127) 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)
10:A:1460[A]:EPE:H32	10:A:1460[A]:EPE:C10	1.61	1.27
10:A:1460[A]:EPE:H32	10:A:1460[A]:EPE:H102	1.25	1.09
1:B:238:SER:HB2	1:B:239:GLY:HA2	1.29	1.09
10:A:1460[A]:EPE:H32	10:A:1460[A]:EPE:H101	1.37	1.03
10:A:1460[A]:EPE:C10	10:A:1460[A]:EPE:C3	2.35	1.00
10:A:1460[A]:EPE:H102	10:A:1460[A]:EPE:C3	1.87	0.90
1:A:409:ASN:ND2	1:A:419:SER:OG	2.06	0.88
10:A:1460[B]:EPE:C10	10:A:1460[B]:EPE:H32	2.04	0.87
10:A:1460[B]:EPE:H32	10:A:1460[B]:EPE:H101	1.60	0.83
1:B:125:THR:H	8:B:1461:MPD:HM1	1.46	0.79

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:147:TYR:OH	10:B:1462[B]:EPE:H51	1.88	0.74
1:B:238:SER:CB	1:B:239:GLY:HA2	2.07	0.72
1:B:168:ASP:OD2	6:B:1455[A]:IMD:H4	1.93	0.69
1:B:238:SER:HB2	1:B:239:GLY:CA	2.18	0.68
1:A:64:LEU:HD12	1:A:120:LEU:HD21	1.75	0.68
1:A:125:THR:H	8:B:1461:MPD:H53	1.58	0.68
10:A:1460[B]:EPE:C10	10:A:1460[B]:EPE:C3	2.66	0.67
1:A:145:LYS:HD3	10:A:1460[A]:EPE:H92	1.75	0.67
1:A:125:THR:H	8:B:1461:MPD:C5	2.10	0.65
1:A:168:ASP:OD2	6:A:1454[B]:IMD:H4	1.97	0.65
1:A:409:ASN:HD22	1:A:419:SER:HA	1.62	0.64
1:A:126:ASP:OD2	8:B:1461:MPD:H12	1.97	0.63
1:B:161:HIS:NE2	15:B:2087:HOH:O	2.29	0.63
1:A:265:THR:O	1:A:265:THR:HG23	1.99	0.62
1:B:284:THR:O	1:B:285:TYR:HB2	1.99	0.62
1:B:125:THR:H	8:B:1461:MPD:CM	2.10	0.62
1:B:124:THR:HA	8:B:1461:MPD:HM3	1.80	0.62
1:A:55:ASN:HB3	1:A:58:THR:OG1	2.01	0.61
1:A:404:GLN:CG	8:A:1457:MPD:H12	2.32	0.59
1:A:315:GLY:HA2	9:A:1459:FMT:C	2.33	0.58
1:B:223:VAL:HB	1:B:224:PRO:HD3	1.85	0.58
1:B:33:ALA:HB3	15:B:2023:HOH:O	2.03	0.58
1:A:439:ILE:HA	1:A:442:ILE:HD12	1.87	0.56
1:B:217:ARG:HB2	1:B:359:PHE:HB3	1.86	0.56
1:A:418:ALA:HA	5:A:1453:NAG:H61	1.87	0.56
1:B:310:PRO:HB3	15:B:2165:HOH:O	2.05	0.56
1:B:191:GLY:HA2	1:B:197:ALA:HB3	1.88	0.55
1:B:383:PHE:CG	1:B:384:PRO:HA	2.41	0.55
10:B:1462[B]:EPE:C7	10:B:1462[B]:EPE:N1	2.69	0.55
8:B:1460:MPD:H52	8:B:1460:MPD:O2	2.07	0.54
1:A:223:VAL:HB	1:A:224:PRO:HD3	1.88	0.54
1:A:175[A]:ILE:HB	1:A:178:PHE:HB3	1.89	0.54
1:B:105:TRP:CD2	1:B:430:GLY:HA3	2.42	0.54
1:B:44:ALA:HB3	1:B:45:PRO:HD3	1.91	0.53
10:A:1460[B]:EPE:C3	10:A:1460[B]:EPE:H102	2.32	0.53
1:A:284:THR:O	1:A:285:TYR:HB2	2.09	0.53
1:A:105:TRP:CD2	1:A:430:GLY:HA3	2.44	0.52
1:A:409:ASN:HD21	1:A:419:SER:HG	1.51	0.52
6:B:1455[B]:IMD:H4	15:B:2176:HOH:O	2.09	0.52
1:B:147:TYR:CZ	10:B:1462[B]:EPE:H51	2.45	0.51
1:B:20:GLN:NE2	15:B:2021:HOH:O	2.43	0.51

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:151:ALA:HB1	1:A:335:ILE:HG12	1.93	0.50
1:A:335:ILE:HG13	1:A:336:PRO:HD2	1.91	0.50
8:B:1459:MPD:HM3	8:B:1459:MPD:O4	2.11	0.50
1:A:64:LEU:HD12	1:A:120:LEU:CD2	2.42	0.50
1:B:184:GLY:HA3	1:B:281:ASN:O	2.10	0.50
1:B:168:ASP:OD1	1:B:169:GLY:N	2.42	0.50
1:B:348:SER:HA	5:B:1453:NAG:H82	1.94	0.50
1:B:82:LEU:N	1:B:82:LEU:HD23	2.27	0.49
1:A:383:PHE:CG	1:A:384:PRO:HA	2.48	0.49
1:A:44:ALA:N	1:A:45:PRO:CD	2.75	0.49
1:B:251:ARG:NE	1:B:411:LEU:O	2.41	0.49
1:A:145:LYS:O	10:A:1460[A]:EPE:H51	2.13	0.49
10:A:1460[A]:EPE:H101	10:A:1460[A]:EPE:C3	2.20	0.48
1:A:404:GLN:HG3	8:A:1457:MPD:H12	1.95	0.48
1:A:335:ILE:HG13	1:A:336:PRO:CD	2.43	0.48
1:B:49:SER:OG	1:B:52[A]:GLN:HG3	2.14	0.48
1:A:389:PRO:HD2	15:A:2245:HOH:O	2.14	0.47
1:A:404:GLN:HG2	8:A:1457:MPD:C1	2.45	0.47
1:B:158:GLU:O	1:B:161:HIS:HB2	2.15	0.47
1:B:398:ILE:HG21	4:B:1449:HEM:CHC	2.44	0.47
10:B:1462[A]:EPE:H101	10:B:1462[A]:EPE:H62	1.41	0.47
1:A:367:SER:O	1:A:372:PHE:HB2	2.16	0.46
1:B:320:PRO:HA	1:B:321:PRO:HA	1.68	0.46
1:B:170:ILE:N	4:B:1449:HEM:O2A	2.48	0.46
1:A:235:PRO:O	1:A:244:ARG:HD2	2.15	0.46
1:A:72:GLY:HA3	1:A:115:HIS:O	2.16	0.46
1:B:19:LYS:HB2	1:B:121:ALA:HB1	1.98	0.46
1:B:265:THR:HA	1:B:266:PRO:HD3	1.74	0.46
1:B:106:VAL:HB	1:B:109:PHE:CD2	2.49	0.45
1:A:44:ALA:HB3	1:A:45:PRO:HD3	1.97	0.45
8:B:1461:MPD:H53	8:B:1461:MPD:O2	2.17	0.45
1:A:22:GLN:HG2	1:A:121:ALA:HB2	1.97	0.45
1:B:44:ALA:N	1:B:45:PRO:CD	2.80	0.45
1:A:124:THR:HA	8:B:1461:MPD:H52	1.97	0.45
10:A:1460[B]:EPE:H102	10:A:1460[B]:EPE:H22	1.49	0.45
1:B:284:THR:O	1:B:285:TYR:CB	2.64	0.45
1:B:223:VAL:N	1:B:224:PRO:CD	2.80	0.44
1:B:412:LEU:HA	1:B:413:PRO:HD3	1.81	0.44
1:A:47:VAL:HG21	1:A:442:ILE:HD13	2.00	0.44
8:B:1461:MPD:HM1	8:B:1461:MPD:O4	2.17	0.44
10:B:1462[B]:EPE:N1	10:B:1462[B]:EPE:H71	2.33	0.44

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:404:GLN:CG	8:A:1457:MPD:C1	2.95	0.44
1:B:18:HIS:CE1	15:B:2015:HOH:O	2.70	0.44
1:B:174:ALA:HB2	1:B:185:GLN:CD	2.38	0.43
1:A:359:PHE:HE2	6:A:1454[B]:IMD:H2	1.82	0.43
1:B:359:PHE:CE2	6:B:1455[A]:IMD:H2	2.53	0.43
1:B:337:TYR:HE2	1:B:354:GLU:HG2	1.83	0.43
10:A:1460[A]:EPE:H101	10:A:1460[A]:EPE:H52	2.01	0.43
1:B:25:TYR:CE2	1:B:145:LYS:HD3	2.54	0.43
1:B:265:THR:HG23	1:B:265:THR:O	2.19	0.43
1:B:80:ASP:N	1:B:80:ASP:OD1	2.52	0.42
1:A:359:PHE:CE2	6:A:1454[B]:IMD:H2	2.54	0.42
1:A:382:ASN:HD22	1:A:382:ASN:HA	1.71	0.42
1:A:233:ASN:ND2	15:A:2188:HOH:O	2.39	0.42
1:A:348:SER:O	1:A:349:ASN:HB2	2.20	0.42
1:A:126:ASP:OD2	1:B:21:LYS:NZ	2.40	0.42
1:A:145:LYS:HE3	15:A:2112:HOH:O	2.20	0.41
1:A:288:ALA:HA	1:A:289:GLY:HA2	1.86	0.41
1:B:81:ASN:OD1	1:B:83:GLY:N	2.52	0.41
1:B:367:SER:O	1:B:372:PHE:HB2	2.21	0.41
1:B:332:ARG:NH1	6:B:1455[B]:IMD:C2	2.83	0.41
1:B:4:LEU:HD12	1:B:4:LEU:HA	1.85	0.41
1:B:251:ARG:NH2	1:B:300:PRO:HG3	2.36	0.41
1:A:4:LEU:HD23	1:A:4:LEU:HA	1.90	0.41
1:A:409:ASN:ND2	1:A:419:SER:HA	2.31	0.41
6:A:1454[A]:IMD:H4	15:A:2258:HOH:O	2.20	0.41
8:A:1458:MPD:H13	15:A:2291:HOH:O	2.21	0.41
10:B:1462[A]:EPE:H61	15:B:2076:HOH:O	2.21	0.41
1:A:398:ILE:HG21	4:A:1449:HEM:CHC	2.51	0.41
1:A:217:ARG:HB2	1:A:359:PHE:HB3	2.03	0.40
1:A:432:GLU:HB3	1:A:434:PHE:CE1	2.56	0.40
1:B:147:TYR:CE1	10:B:1462[B]:EPE:H51	2.56	0.40
1:B:288:ALA:HA	1:B:289:GLY:HA2	1.55	0.40
1:B:72:GLY:HA3	1:B:115:HIS:O	2.21	0.40
1:A:175[A]:ILE:HG23	1:A:190:ALA:HA	2.02	0.40
1:B:24:PHE:O	1:B:148:SER:HA	2.22	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	448/446 (100%)	439 (98%)	9 (2%)	0	100	100
1	B	445/446 (100%)	429 (96%)	16 (4%)	0	100	100
All	All	893/892 (100%)	868 (97%)	25 (3%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains [i](#)

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	357/353 (101%)	350 (98%)	7 (2%)	55	60
1	B	354/353 (100%)	350 (99%)	4 (1%)	73	79
All	All	711/706 (101%)	700 (98%)	11 (2%)	67	71

All (11) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	152[A]	SER
1	A	152[B]	SER
1	A	162	GLU
1	A	363	GLN
1	A	391	THR
1	A	406	ARG
1	A	421	SER
1	B	58	THR

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	B	80	ASP
1	B	363	GLN
1	B	406	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	382	ASN
1	A	409	ASN
1	B	18	HIS

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

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

5.5 Carbohydrates [i](#)

5 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	C	1	2,1	14,14,15	0.62	0	17,19,21	0.85	1 (5%)
2	NAG	C	2	2	14,14,15	0.55	0	17,19,21	1.01	0
3	NAG	D	1	3,1	14,14,15	0.55	0	17,19,21	0.96	1 (5%)
3	NAG	D	2	3	14,14,15	0.58	0	17,19,21	0.89	0
3	BMA	D	3	3	11,11,12	0.67	0	15,15,17	0.93	1 (6%)

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	C	1	2,1	-	0/6/23/26	0/1/1/1
2	NAG	C	2	2	-	0/6/23/26	0/1/1/1
3	NAG	D	1	3,1	-	0/6/23/26	0/1/1/1
3	NAG	D	2	3	-	2/6/23/26	0/1/1/1
3	BMA	D	3	3	-	1/2/19/22	0/1/1/1

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	3	BMA	C1-C2-C3	2.83	113.15	109.67
3	D	1	NAG	O5-C1-C2	-2.41	107.48	111.29
2	C	1	NAG	O5-C1-C2	-2.33	107.60	111.29

There are no chirality outliers.

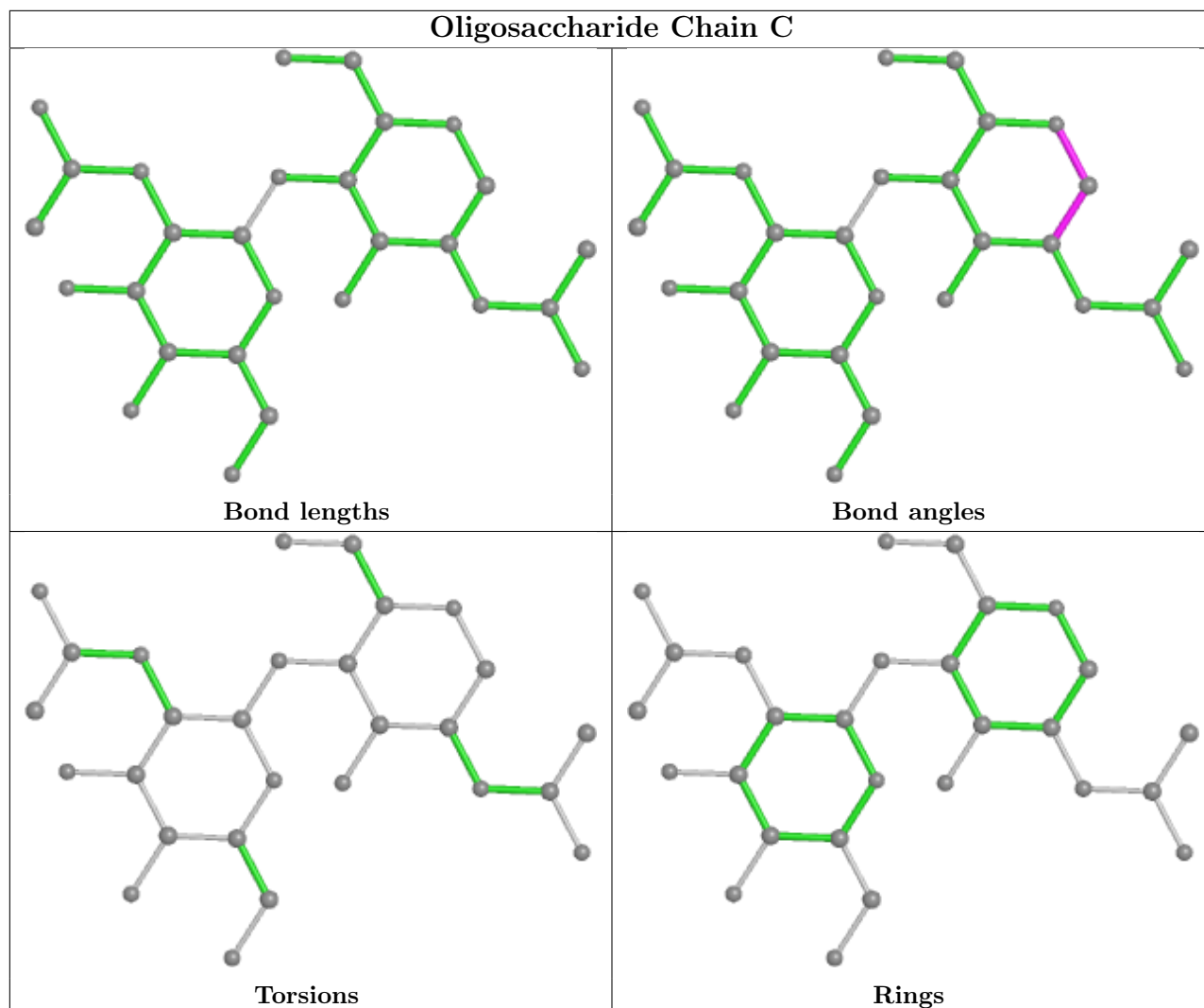
All (3) torsion outliers are listed below:

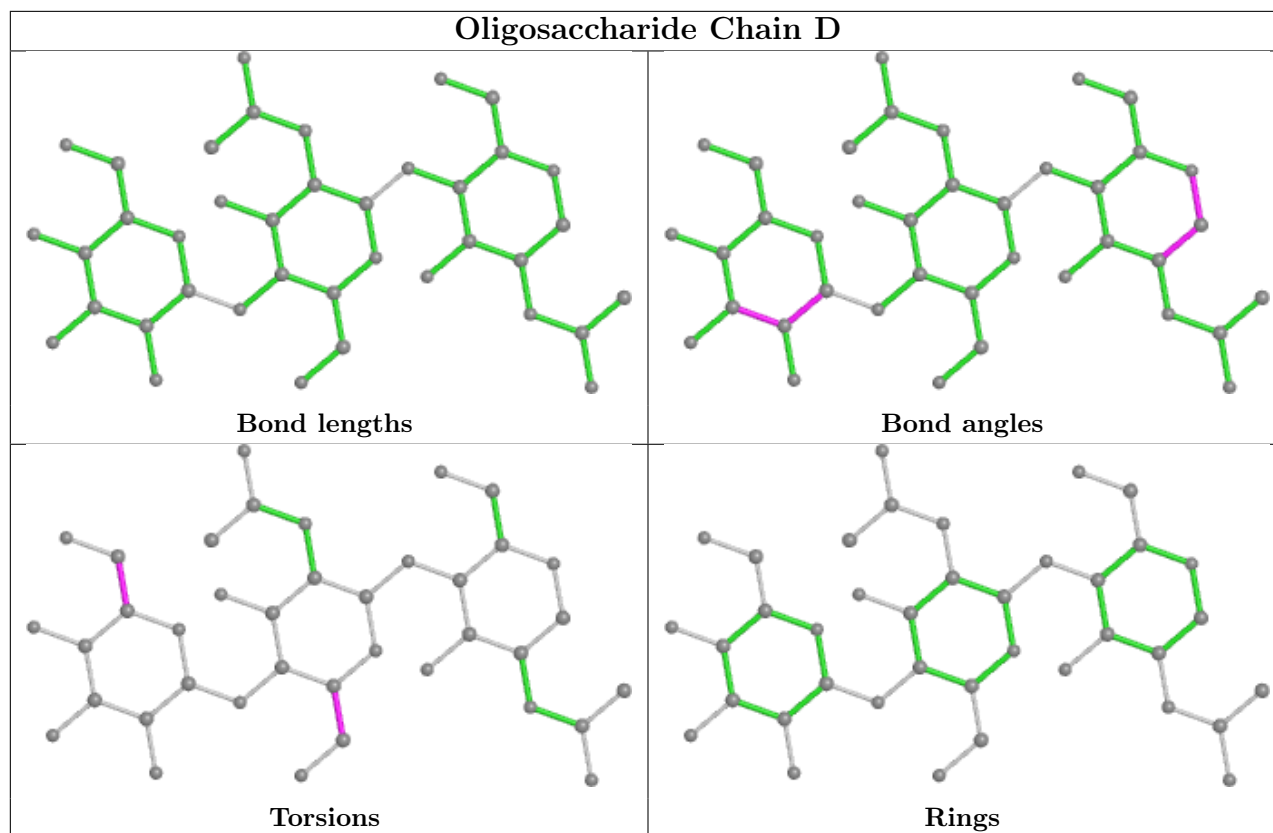
Mol	Chain	Res	Type	Atoms
3	D	2	NAG	C4-C5-C6-O6
3	D	2	NAG	O5-C5-C6-O6
3	D	3	BMA	O5-C5-C6-O6

There are no ring outliers.

No monomer is involved in short contacts.

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





5.6 Ligand geometry [i](#)

Of 35 ligands modelled in this entry, 2 are monoatomic - leaving 33 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	$\# Z > 2$
5	NAG	A	1453	1	14,14,15	0.56	0	17,19,21	1.07	1 (5%)
14	GOA	B	1466	-	4,4,4	0.98	0	4,4,4	0.98	0
5	NAG	B	1454	1	14,14,15	0.62	0	17,19,21	0.92	1 (5%)
5	NAG	B	1453	1	14,14,15	0.53	0	17,19,21	1.09	1 (5%)
10	EPE	B	1462[B]	-	15,15,15	1.97	1 (6%)	18,20,20	1.08	3 (16%)
8	MPD	B	1461	-	7,7,7	0.26	0	9,10,10	0.30	0
11	ACT	B	1464	-	3,3,3	0.78	0	3,3,3	0.68	0
11	ACT	B	1467	-	3,3,3	0.76	0	3,3,3	0.78	0
10	EPE	A	1460[B]	-	15,15,15	1.98	1 (6%)	18,20,20	1.04	2 (11%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
6	IMD	A	1455	-	3,5,5	0.36	0	4,5,5	0.55	0
6	IMD	B	1455[B]	4	3,5,5	0.36	0	4,5,5	0.57	0
9	FMT	A	1459	-	2,2,2	0.75	0	1,1,1	0.54	0
10	EPE	B	1462[A]	-	15,15,15	1.99	1 (6%)	18,20,20	1.10	3 (16%)
8	MPD	B	1459	-	7,7,7	0.25	0	9,10,10	0.26	0
11	ACT	A	1463	-	3,3,3	0.77	0	3,3,3	0.68	0
6	IMD	A	1454[B]	4	3,5,5	0.36	0	4,5,5	0.50	0
8	MPD	B	1458	-	7,7,7	0.24	0	9,10,10	0.26	0
8	MPD	B	1457	-	7,7,7	0.24	0	9,10,10	0.27	0
4	HEM	A	1449	1,6	41,50,50	1.30	4 (9%)	45,82,82	1.92	8 (17%)
4	HEM	B	1449	1,6	41,50,50	1.29	5 (12%)	45,82,82	1.89	7 (15%)
10	EPE	A	1460[A]	-	15,15,15	1.99	1 (6%)	18,20,20	1.06	3 (16%)
11	ACT	B	1463	-	3,3,3	0.77	0	3,3,3	0.77	0
5	NAG	A	1452	1	14,14,15	0.56	0	17,19,21	0.72	0
6	IMD	B	1455[A]	4	3,5,5	0.35	0	4,5,5	0.54	0
11	ACT	A	1462	-	3,3,3	0.77	0	3,3,3	0.75	0
8	MPD	A	1458	-	7,7,7	0.29	0	9,10,10	0.30	0
6	IMD	A	1454[A]	4	3,5,5	0.38	0	4,5,5	0.50	0
12	OXD	A	1464	-	5,5,5	1.78	1 (20%)	6,6,6	1.21	0
6	IMD	B	1456	-	3,5,5	0.32	0	4,5,5	0.57	0
11	ACT	A	1461	-	3,3,3	0.78	0	3,3,3	0.69	0
8	MPD	B	1460	-	7,7,7	0.24	0	9,10,10	0.26	0
7	MRD	A	1456	-	7,7,7	0.25	0	9,10,10	0.30	0
8	MPD	A	1457	-	7,7,7	0.26	0	9,10,10	0.29	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
5	NAG	A	1453	1	1/1/5/7	1/6/23/26	0/1/1/1
14	GOA	B	1466	-	-	2/2/2/2	-
5	NAG	B	1454	1	-	2/6/23/26	0/1/1/1
5	NAG	B	1453	1	-	2/6/23/26	0/1/1/1
10	EPE	B	1462[B]	-	-	6/9/19/19	0/1/1/1
8	MPD	B	1461	-	-	0/5/5/5	-
10	EPE	A	1460[B]	-	-	5/9/19/19	0/1/1/1
6	IMD	A	1455	-	-	-	0/1/1/1
6	IMD	B	1455[B]	4	-	-	0/1/1/1

Continued on next page...

Continued from previous page...

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
10	EPE	B	1462[A]	-	-	6/9/19/19	0/1/1/1
8	MPD	B	1459	-	-	0/5/5/5	-
6	IMD	A	1454[B]	4	-	-	0/1/1/1
8	MPD	B	1458	-	-	0/5/5/5	-
8	MPD	B	1457	-	-	0/5/5/5	-
4	HEM	A	1449	1,6	-	3/12/54/54	-
4	HEM	B	1449	1,6	-	2/12/54/54	-
10	EPE	A	1460[A]	-	-	5/9/19/19	0/1/1/1
5	NAG	A	1452	1	-	0/6/23/26	0/1/1/1
6	IMD	B	1455[A]	4	-	-	0/1/1/1
8	MPD	A	1458	-	-	2/5/5/5	-
6	IMD	A	1454[A]	4	-	-	0/1/1/1
12	OXD	A	1464	-	-	2/4/4/4	-
6	IMD	B	1456	-	-	-	0/1/1/1
8	MPD	B	1460	-	-	4/5/5/5	-
7	MRD	A	1456	-	-	1/5/5/5	-
8	MPD	A	1457	-	-	0/5/5/5	-

All (14) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
10	B	1462[A]	EPE	C10-S	-7.45	1.66	1.77
10	A	1460[A]	EPE	C10-S	-7.43	1.67	1.77
10	A	1460[B]	EPE	C10-S	-7.41	1.67	1.77
10	B	1462[B]	EPE	C10-S	-7.38	1.67	1.77
4	B	1449	HEM	C1B-NB	-3.64	1.34	1.40
4	A	1449	HEM	C1B-NB	-3.64	1.34	1.40
12	A	1464	OXD	C2-C1	-2.71	1.46	1.54
4	A	1449	HEM	C4D-ND	-2.68	1.35	1.40
4	A	1449	HEM	CHB-C1B	2.62	1.41	1.35
4	A	1449	HEM	FE-NB	2.59	2.09	1.96
4	B	1449	HEM	FE-NB	2.58	2.09	1.96
4	B	1449	HEM	C4D-ND	-2.48	1.36	1.40
4	B	1449	HEM	C3B-C4B	2.22	1.49	1.44
4	B	1449	HEM	CHB-C1B	2.22	1.40	1.35

All (29) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	1449	HEM	C1B-NB-C4B	5.52	110.77	105.07

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	1449	HEM	C1B-NB-C4B	5.14	110.39	105.07
4	A	1449	HEM	CHC-C4B-NB	5.02	129.88	124.43
4	B	1449	HEM	CHD-C1D-ND	4.89	129.75	124.43
4	A	1449	HEM	CHD-C1D-ND	4.54	129.36	124.43
4	B	1449	HEM	CHC-C4B-NB	4.29	129.09	124.43
4	B	1449	HEM	CHD-C1D-C2D	-3.93	118.84	124.98
4	B	1449	HEM	CHA-C4D-ND	3.92	129.22	124.38
4	A	1449	HEM	CHD-C1D-C2D	-3.89	118.90	124.98
4	A	1449	HEM	CHA-C4D-ND	3.18	128.31	124.38
5	A	1453	NAG	C2-N2-C7	-3.05	118.56	122.90
4	B	1449	HEM	CHA-C4D-C3D	-2.97	119.76	125.33
4	A	1449	HEM	CHA-C4D-C3D	-2.90	119.89	125.33
4	B	1449	HEM	CHB-C1B-NB	2.74	127.76	124.38
4	A	1449	HEM	CHB-C1B-NB	2.52	127.49	124.38
10	B	1462[B]	EPE	O2S-S-C10	2.44	109.86	106.92
10	A	1460[A]	EPE	O1S-S-C10	2.41	109.81	106.92
10	A	1460[B]	EPE	O1S-S-C10	2.40	109.81	106.92
4	A	1449	HEM	CAD-C3D-C4D	2.39	128.84	124.66
5	B	1453	NAG	C1-O5-C5	2.30	115.31	112.19
10	B	1462[A]	EPE	O3S-S-C10	2.25	109.40	105.77
10	B	1462[B]	EPE	O1S-S-C10	2.20	109.56	106.92
10	B	1462[A]	EPE	O1S-S-C10	2.17	109.53	106.92
10	B	1462[A]	EPE	O2S-S-C10	2.16	109.51	106.92
10	A	1460[A]	EPE	O3S-S-C10	2.13	109.21	105.77
10	A	1460[B]	EPE	O3S-S-C10	2.13	109.21	105.77
10	B	1462[B]	EPE	O3S-S-C10	2.10	109.17	105.77
5	B	1454	NAG	C3-C4-C5	2.08	113.95	110.24
10	A	1460[A]	EPE	O2S-S-C10	2.05	109.38	106.92

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
5	A	1453	NAG	C1

All (43) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
8	A	1458	MPD	C2-C3-C4-O4
8	A	1458	MPD	C2-C3-C4-C5
10	A	1460[A]	EPE	C10-C9-N1-C2
10	A	1460[B]	EPE	C10-C9-N1-C2
10	A	1460[B]	EPE	S-C10-C9-N1

Continued on next page...

Continued from previous page...

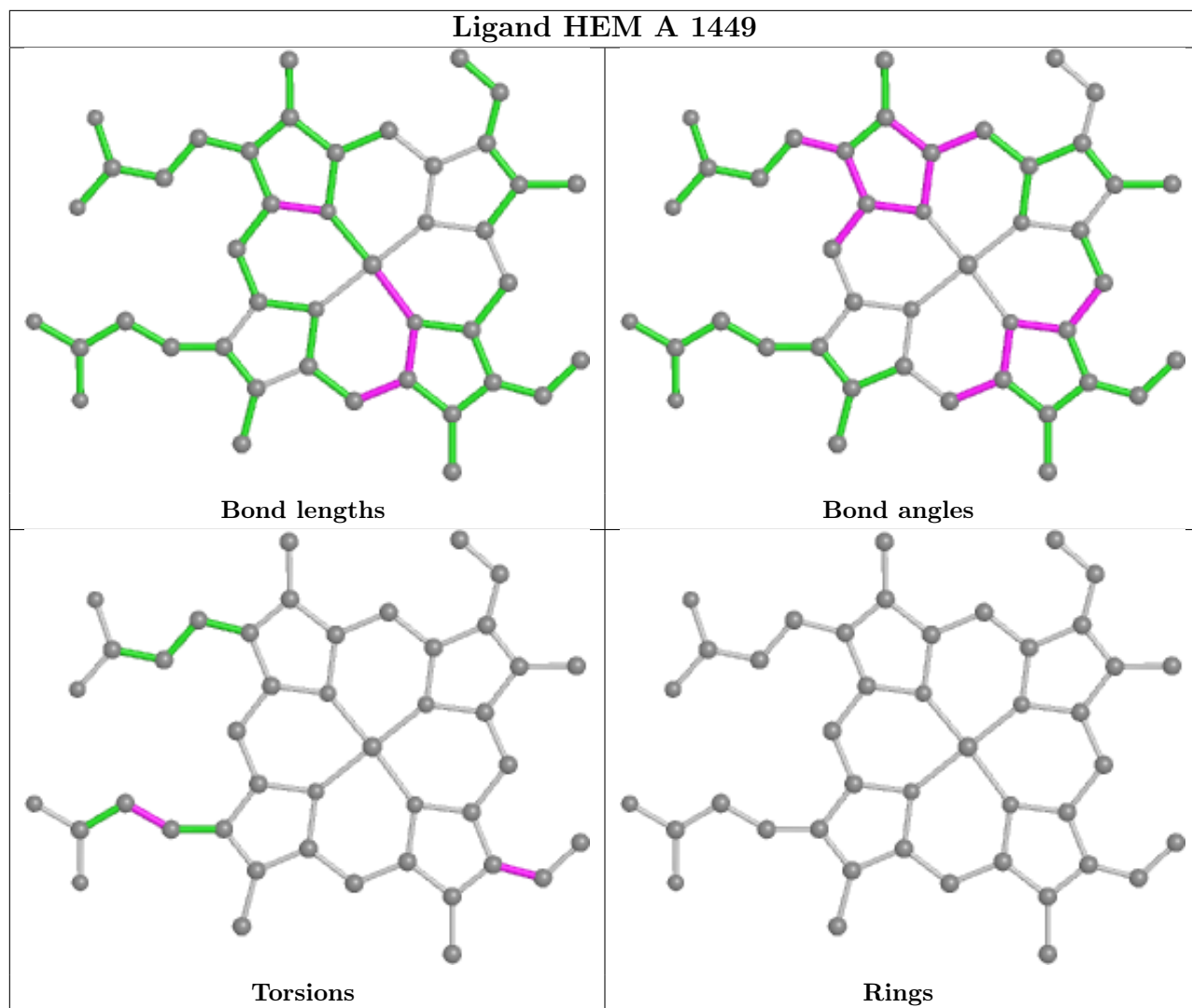
Mol	Chain	Res	Type	Atoms
10	B	1462[A]	EPE	C10-C9-N1-C6
10	B	1462[A]	EPE	S-C10-C9-N1
10	B	1462[B]	EPE	C10-C9-N1-C6
10	B	1462[B]	EPE	S-C10-C9-N1
14	B	1466	GOA	O-C-CA-O2
14	B	1466	GOA	OXT-C-CA-O2
5	B	1454	NAG	C4-C5-C6-O6
5	B	1453	NAG	O5-C5-C6-O6
10	B	1462[B]	EPE	N4-C7-C8-O8
5	B	1454	NAG	O5-C5-C6-O6
5	B	1453	NAG	C4-C5-C6-O6
4	A	1449	HEM	C2A-CAA-CBA-CGA
4	B	1449	HEM	C2A-CAA-CBA-CGA
10	A	1460[B]	EPE	C8-C7-N4-C3
10	B	1462[A]	EPE	C9-C10-S-O3S
5	A	1453	NAG	O5-C5-C6-O6
10	A	1460[B]	EPE	C8-C7-N4-C5
10	A	1460[B]	EPE	C10-C9-N1-C6
10	B	1462[B]	EPE	C10-C9-N1-C2
8	B	1460	MPD	O2-C2-C3-C4
4	A	1449	HEM	C2B-C3B-CAB-CBB
7	A	1456	MRD	C2-C3-C4-C5
10	A	1460[A]	EPE	N4-C7-C8-O8
10	B	1462[A]	EPE	C9-C10-S-O1S
10	B	1462[A]	EPE	C9-C10-S-O2S
8	B	1460	MPD	C2-C3-C4-O4
8	B	1460	MPD	C1-C2-C3-C4
10	B	1462[B]	EPE	C8-C7-N4-C5
10	A	1460[A]	EPE	C10-C9-N1-C6
10	B	1462[A]	EPE	C10-C9-N1-C2
10	B	1462[B]	EPE	C8-C7-N4-C3
12	A	1464	OXD	O3-C1-C2-O4
12	A	1464	OXD	O5-C1-C2-O6
4	B	1449	HEM	C2B-C3B-CAB-CBB
4	A	1449	HEM	C4B-C3B-CAB-CBB
10	A	1460[A]	EPE	C8-C7-N4-C3
8	B	1460	MPD	C2-C3-C4-C5
10	A	1460[A]	EPE	C8-C7-N4-C5

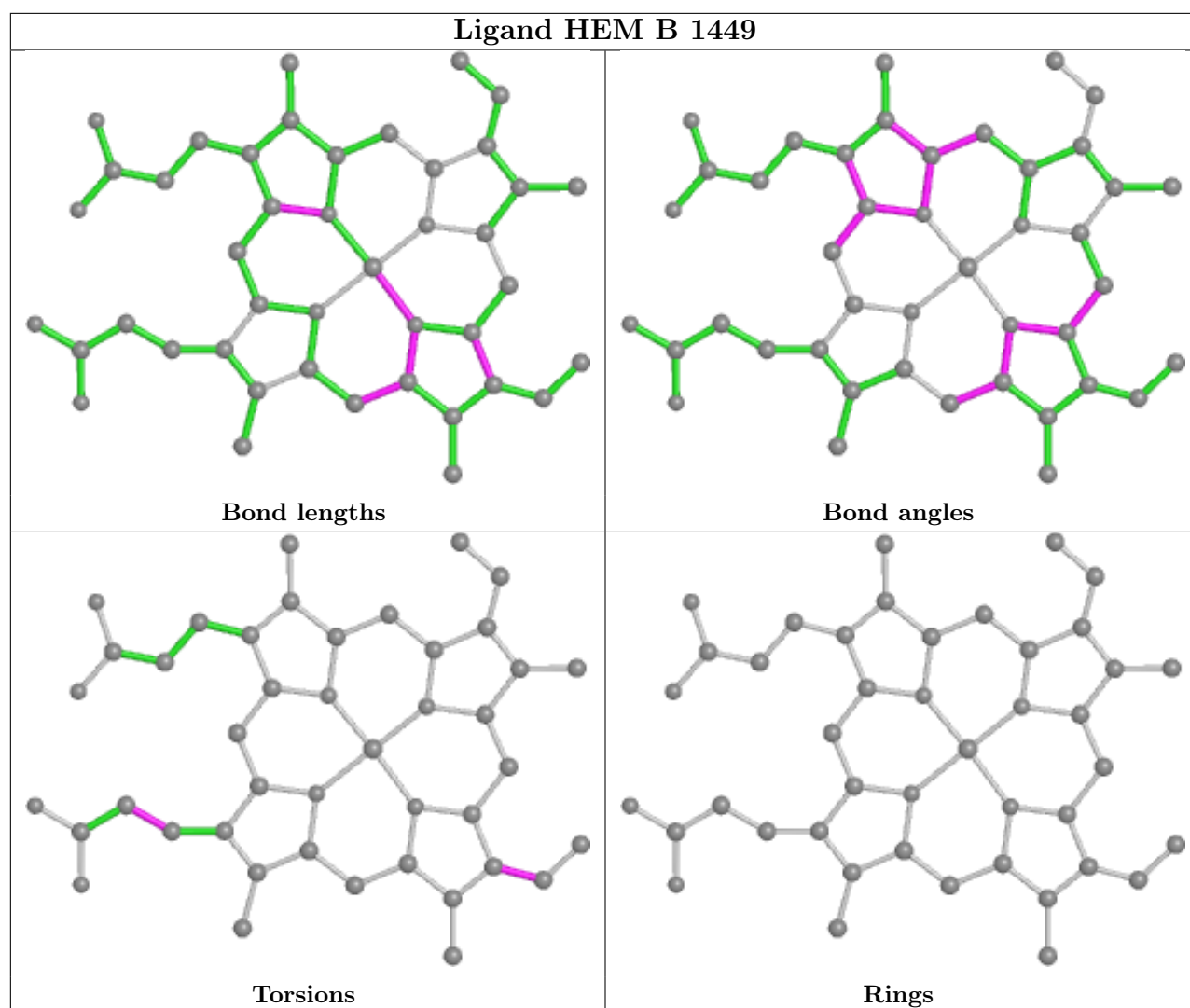
There are no ring outliers.

18 monomers are involved in 51 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	1453	NAG	1	0
5	B	1453	NAG	1	0
10	B	1462[B]	EPE	5	0
8	B	1461	MPD	9	0
10	A	1460[B]	EPE	5	0
6	B	1455[B]	IMD	2	0
9	A	1459	FMT	1	0
10	B	1462[A]	EPE	2	0
8	B	1459	MPD	1	0
6	A	1454[B]	IMD	3	0
4	A	1449	HEM	1	0
4	B	1449	HEM	2	0
10	A	1460[A]	EPE	9	0
6	B	1455[A]	IMD	2	0
8	A	1458	MPD	1	0
6	A	1454[A]	IMD	1	0
8	B	1460	MPD	1	0
8	A	1457	MPD	4	0

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





5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [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	446/446 (100%)	-0.35	0 100 100	10, 18, 26, 50	0
1	B	446/446 (100%)	-0.18	4 (0%) 84 86	13, 22, 35, 62	0
All	All	892/892 (100%)	-0.27	4 (0%) 92 93	10, 20, 32, 62	0

All (4) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	425	PHE	2.8
1	B	3	SER	2.3
1	B	238	SER	2.3
1	B	236	ALA	2.1

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

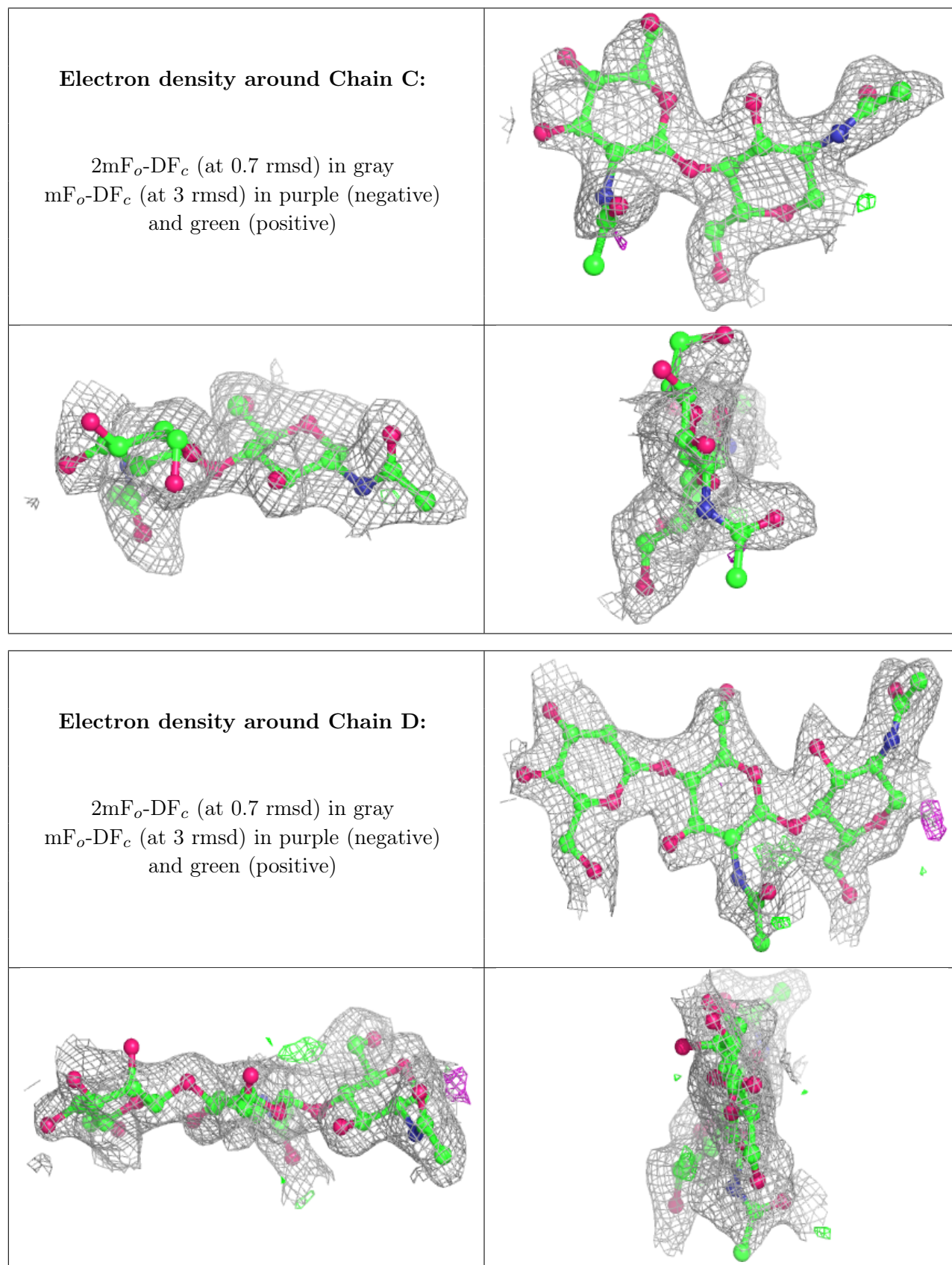
There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates [i](#)

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	NAG	C	2	14/15	0.76	0.22	46,54,63,75	0
3	NAG	D	2	14/15	0.80	0.20	46,50,53,53	0
3	BMA	D	3	11/12	0.80	0.24	48,52,55,58	0
2	NAG	C	1	14/15	0.91	0.10	25,31,35,41	0
3	NAG	D	1	14/15	0.91	0.11	25,34,38,44	0

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



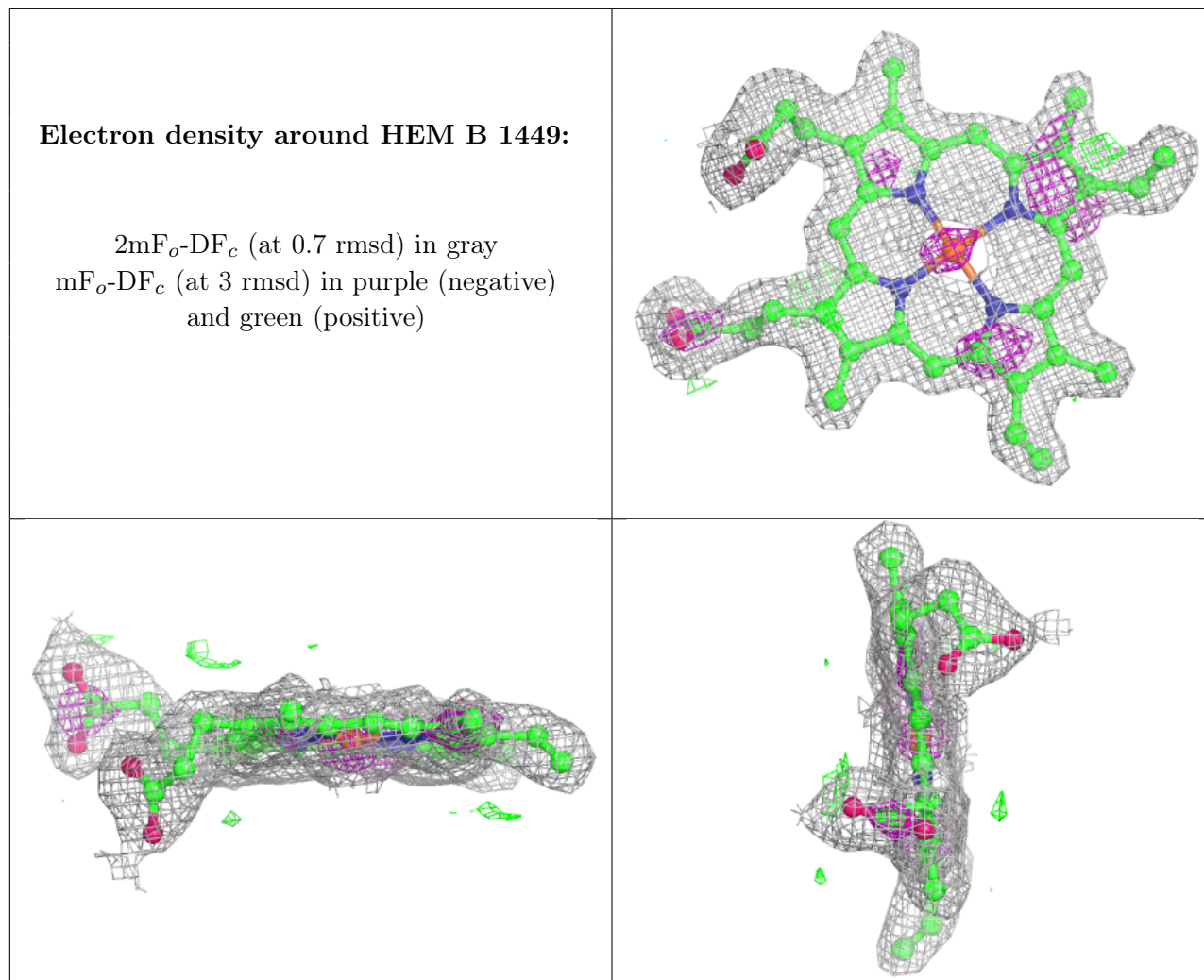
6.4 Ligands

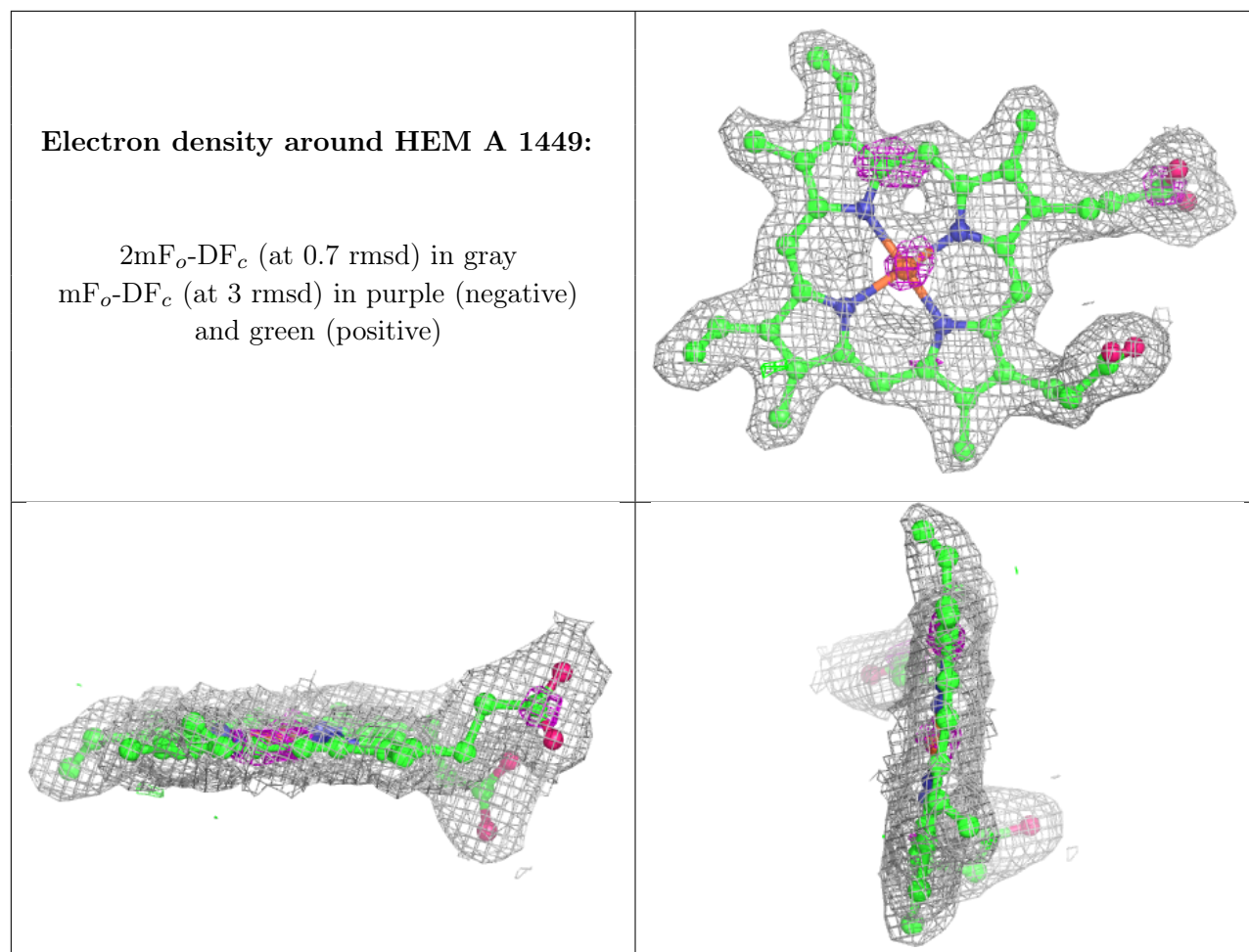
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
10	EPE	B	1462[A]	15/15	0.69	0.34	41,49,89,90	15
10	EPE	B	1462[B]	15/15	0.69	0.34	29,36,45,45	15
8	MPD	A	1458	8/8	0.75	0.28	41,46,52,57	0
5	NAG	B	1454	14/15	0.76	0.28	47,57,64,67	0
5	NAG	B	1453	14/15	0.77	0.25	40,51,65,72	0
8	MPD	B	1459	8/8	0.77	0.36	51,55,59,62	0
12	OXD	A	1464	6/6	0.77	0.24	44,50,51,52	0
14	GOA	B	1466	5/5	0.77	0.20	45,47,53,54	0
11	ACT	A	1461	4/4	0.78	0.36	34,39,41,45	0
11	ACT	B	1467	4/4	0.79	0.21	52,53,54,57	0
10	EPE	A	1460[A]	15/15	0.81	0.41	41,50,54,55	15
10	EPE	A	1460[B]	15/15	0.81	0.41	35,49,57,58	15
8	MPD	B	1460	8/8	0.82	0.22	43,48,50,53	0
5	NAG	A	1453	14/15	0.83	0.17	36,43,45,46	0
8	MPD	B	1458	8/8	0.84	0.20	42,44,47,50	0
11	ACT	A	1462	4/4	0.85	0.20	51,52,55,55	0
8	MPD	A	1457	8/8	0.85	0.21	48,54,58,59	0
7	MRD	A	1456	8/8	0.86	0.18	18,22,24,27	0
8	MPD	B	1461	8/8	0.86	0.19	19,27,30,33	0
11	ACT	B	1464	4/4	0.86	0.17	51,53,56,56	0
8	MPD	B	1457	8/8	0.87	0.16	21,24,27,31	0
9	FMT	A	1459	3/3	0.87	0.20	30,30,33,33	0
5	NAG	A	1452	14/15	0.88	0.13	26,32,36,38	0
11	ACT	A	1463	4/4	0.93	0.16	39,40,45,46	0
11	ACT	B	1463	4/4	0.93	0.29	43,43,46,50	0
6	IMD	B	1456	5/5	0.95	0.13	15,15,17,17	0
6	IMD	A	1454[A]	5/5	0.96	0.12	12,12,12,13	5
6	IMD	A	1454[B]	5/5	0.96	0.12	11,11,12,12	5
4	HEM	B	1449	43/43	0.96	0.12	13,15,20,30	0
6	IMD	B	1455[A]	5/5	0.97	0.09	15,15,16,16	5
6	IMD	B	1455[B]	5/5	0.97	0.09	14,14,15,15	5
4	HEM	A	1449	43/43	0.97	0.10	11,13,15,20	0
13	K	A	1465	1/1	0.97	0.08	27,27,27,27	1
13	K	B	1465	1/1	0.97	0.06	29,29,29,29	1
6	IMD	A	1455	5/5	0.97	0.10	21,22,23,23	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 [i](#)

There are no such residues in this entry.