



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 5, 2026 – 01:40 PM EDT

PDB ID : 4OEK / pdb_00004oek
Title : Crystal Structure of the Complex of goat Lactoperoxidase with Phenylethylamine at 2.47 Å Resolution
Authors : Kumar, M.; Singh, R.P.; Sinha, M.; Bhushan, A.; Kaur, P.; Sharma, S.; Singh, T.P.
Deposited on : 2014-01-13
Resolution : 2.47 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.015 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.50

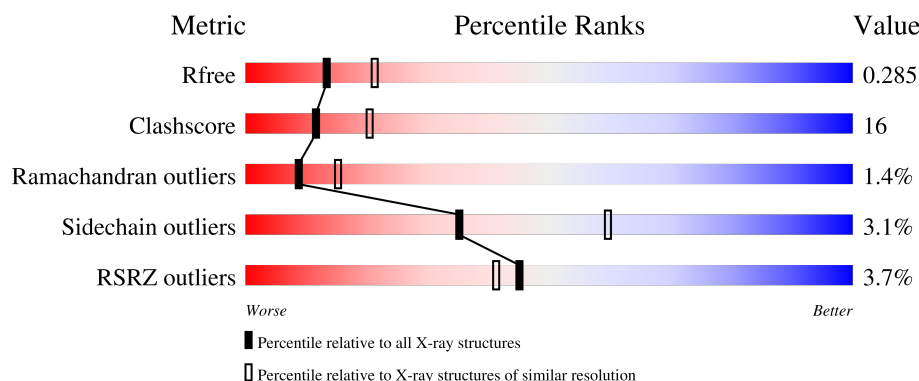
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.47 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	7589 (2.50-2.46)
Clashscore	190562	8295 (2.50-2.46)
Ramachandran outliers	187476	8164 (2.50-2.46)
Sidechain outliers	187428	8166 (2.50-2.46)
RSRZ outliers	180081	7593 (2.50-2.46)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	595	 4% 72% 26%
2	B	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	615	-	-	X	-
6	IOD	A	620	-	-	X	-
9	PEA	A	627	-	-	X	-

2 Entry composition [i](#)

There are 10 unique types of molecules in this entry. The entry contains 5166 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called Lactoperoxidase.

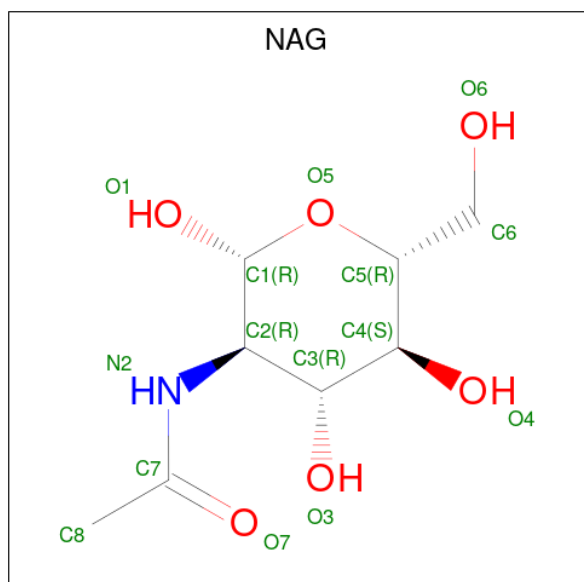
Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	595	Total	C	N	O	P	S	0	0	0
			4757	3021	844	865	1	26			

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



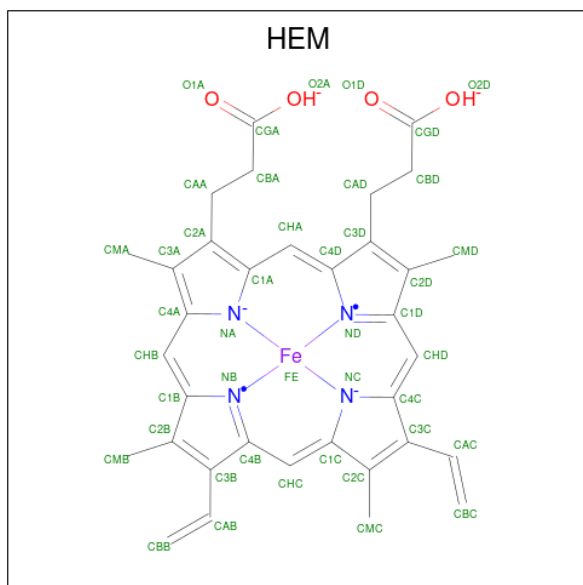
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	B	2	Total	C	N	O	0	0	0
			28	16	2	10			

- Molecule 3 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: C₈H₁₅NO₆).



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

- Molecule 4 is PROTOPORPHYRIN IX CONTAINING FE (CCD ID: HEM) (formula: $C_{34}H_{32}FeN_4O_4$).



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

- Molecule 5 is CALCIUM ION (CCD ID: CA) (formula: Ca).

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

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

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	13	Total	I	0	0
			13	13		

- Molecule 7 is THIOCYANATE ION (CCD ID: SCN) (formula: CNS).



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

- Molecule 8 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: C₂H₆O₂).



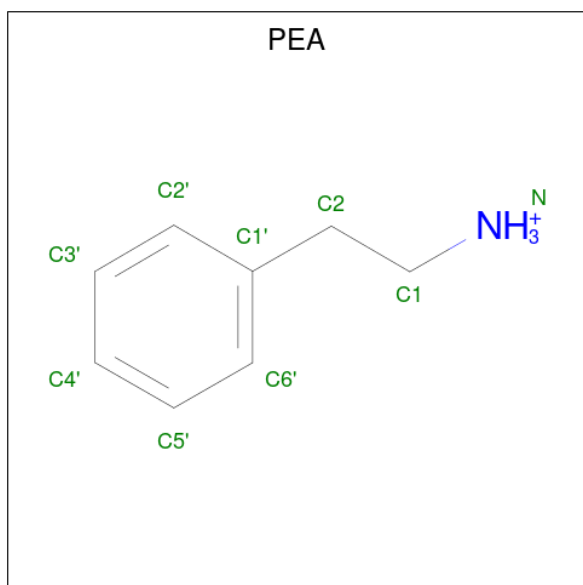
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
8	A	1	Total	C	O	0	0
			4	2	2		
8	A	1	Total	C	O	0	0
			4	2	2		
8	A	1	Total	C	O	0	0
			4	2	2		

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

- Molecule 9 is 2-PHENYLETHYLAMINE (CCD ID: PEA) (formula: $C_8H_{12}N$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
9	A	1	Total	C	N	0	0
			9	8	1		

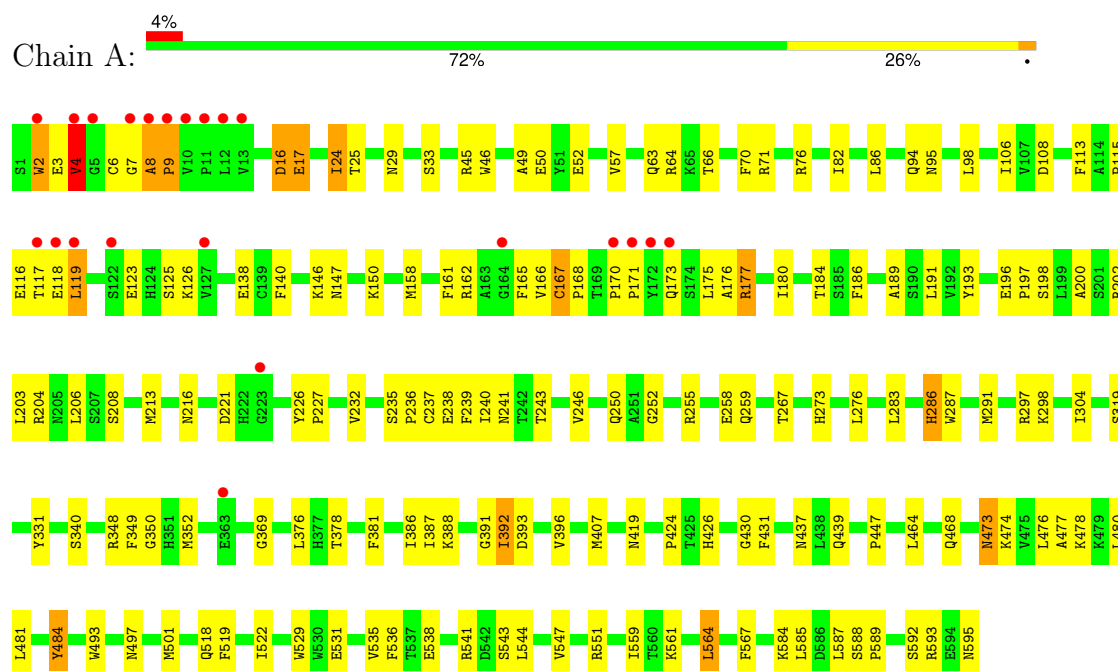
- Molecule 10 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
10	A	250	Total	O	0	0
			250	250		

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: Lactoperoxidase



• Molecule 2: 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	53.17Å 79.89Å 75.28Å 90.00° 101.80° 90.00°	Depositor
Resolution (Å)	73.69 – 2.47 73.69 – 2.47	Depositor EDS
% Data completeness (in resolution range)	88.0 (73.69-2.47) 88.0 (73.69-2.47)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.08	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.43 (at 2.48Å)	Xtriage
Refinement program	REFMAC 5.6.0117	Depositor
R, R_{free}	0.204 , 0.281 0.208 , 0.285	Depositor DCC
R_{free} test set	995 reflections (4.47%)	wwPDB-VP
Wilson B-factor (Å ²)	40.0	Xtriage
Anisotropy	0.340	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 26.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	5166	wwPDB-VP
Average B, all atoms (Å ²)	44.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.66% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: SCN, CA, IOD, EDO, NAG, PEA, HEM, SEP

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	1.01	2/4875 (0.0%)	0.96	11/6621 (0.2%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	240	ILE	C-N	13.77	1.53	1.33
1	A	241	ASN	C-N	-9.00	1.22	1.33

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	33	SER	CA-C-N	7.92	128.00	119.28
1	A	33	SER	C-N-CA	7.92	128.00	119.28
1	A	369	GLY	CA-C-N	6.10	127.46	119.84
1	A	369	GLY	C-N-CA	6.10	127.46	119.84
1	A	95	ASN	N-CA-C	5.50	117.92	110.88
1	A	484	TYR	N-CA-C	-5.34	100.59	109.46
1	A	240	ILE	O-C-N	-5.26	115.16	122.05
1	A	240	ILE	CA-C-N	5.20	129.72	121.72
1	A	240	ILE	C-N-CA	5.20	129.72	121.72
1	A	173	GLN	N-CA-C	5.16	117.06	110.61
1	A	94	GLN	N-CA-C	5.14	119.18	113.01

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	4757	0	4643	149	0
2	B	28	0	25	1	0
3	A	42	0	39	0	0
4	A	43	0	30	15	0
5	A	1	0	0	0	0
6	A	13	0	0	7	0
7	A	3	0	0	0	0
8	A	20	0	30	3	0
9	A	9	0	12	13	0
10	A	250	0	0	12	0
All	All	5166	0	4779	155	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 16.

All (155) 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:258:GLU:OE2	4:A:606:HEM:CMB	1.73	1.36
1:A:108:ASP:OD2	4:A:606:HEM:CMD	1.76	1.32
6:A:616:IOD:I	10:A:719:HOH:O	2.36	1.13
1:A:258:GLU:OE2	4:A:606:HEM:HMB1	0.91	1.07
1:A:108:ASP:OD2	4:A:606:HEM:HMD1	0.82	0.99
1:A:108:ASP:CG	4:A:606:HEM:HMD1	1.86	0.98
1:A:551:ARG:HD3	1:A:584:LYS:HA	1.57	0.87
1:A:167:CYS:HB2	1:A:168:PRO:CD	2.09	0.82
1:A:158:MET:HE2	1:A:431:PHE:HA	1.64	0.80
10:A:761:HOH:O	2:B:2:NAG:O6	1.99	0.80
1:A:538:GLU:HG2	1:A:541:ARG:NH2	1.98	0.79
1:A:258:GLU:CD	4:A:606:HEM:HMB1	2.03	0.78
6:A:615:IOD:I	10:A:813:HOH:O	2.73	0.76
1:A:258:GLU:CD	4:A:606:HEM:CMB	2.59	0.74
1:A:258:GLU:HG3	9:A:627:PEA:H6'	1.69	0.73
1:A:203:LEU:HB3	1:A:213:MET:HE1	1.70	0.73
4:A:606:HEM:C4A	9:A:627:PEA:H12	2.25	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:255:ARG:HG2	9:A:627:PEA:H2'	1.70	0.72
4:A:606:HEM:C3A	9:A:627:PEA:H12	2.28	0.68
1:A:473:ASN:ND2	1:A:476:LEU:H	1.92	0.68
1:A:538:GLU:HG2	1:A:541:ARG:HH22	1.57	0.67
1:A:287:TRP:CZ2	1:A:291:MET:HE2	2.30	0.66
1:A:407:MET:HB3	1:A:501:MET:HE3	1.78	0.65
1:A:258:GLU:HG3	9:A:627:PEA:H21	1.77	0.65
1:A:167:CYS:HB2	1:A:168:PRO:HD3	1.79	0.64
1:A:108:ASP:CG	4:A:606:HEM:CMD	2.60	0.64
1:A:63:GLN:O	1:A:71:ARG:NH1	2.32	0.62
1:A:52:GLU:CG	1:A:57:VAL:HG12	2.29	0.62
1:A:123:GLU:HB3	1:A:126:LYS:HB2	1.81	0.62
1:A:167:CYS:CB	1:A:168:PRO:CD	2.79	0.61
1:A:255:ARG:CG	9:A:627:PEA:H22	2.31	0.61
1:A:203:LEU:HB3	1:A:213:MET:CE	2.30	0.61
1:A:255:ARG:HB3	9:A:627:PEA:H22	1.82	0.60
1:A:2:TRP:CD1	1:A:175:LEU:HD13	2.36	0.60
1:A:567:PHE:HB2	6:A:615:IOD:I	2.72	0.60
1:A:146:LYS:O	1:A:147:ASN:HB2	2.03	0.59
1:A:175:LEU:HD12	1:A:176:ALA:H	1.67	0.59
1:A:258:GLU:CG	9:A:627:PEA:H6'	2.33	0.58
1:A:206:LEU:C	10:A:861:HOH:O	2.47	0.57
1:A:239:PHE:HB2	10:A:852:HOH:O	2.03	0.57
1:A:387:ILE:HG22	1:A:388:LYS:HG3	1.86	0.57
1:A:8:ALA:H	1:A:9:PRO:HD3	1.70	0.56
1:A:16:ASP:OD1	1:A:16:ASP:C	2.47	0.56
1:A:501:MET:HE1	10:A:740:HOH:O	2.05	0.56
1:A:2:TRP:HB3	10:A:875:HOH:O	2.06	0.56
1:A:189:ALA:HB2	1:A:304:ILE:HD12	1.86	0.56
1:A:258:GLU:HG3	9:A:627:PEA:C6'	2.35	0.56
1:A:9:PRO:HG2	10:A:924:HOH:O	2.06	0.56
1:A:167:CYS:CB	1:A:168:PRO:HD3	2.36	0.56
1:A:348:ARG:HH11	1:A:437:ASN:HD22	1.53	0.55
1:A:167:CYS:HB2	1:A:168:PRO:HD2	1.88	0.55
1:A:197:PRO:HA	8:A:624:EDO:H12	1.89	0.55
1:A:200:ALA:O	1:A:204:ARG:HG3	2.05	0.55
1:A:221:ASP:HB2	1:A:226:TYR:CZ	2.42	0.55
1:A:588:SER:HB2	1:A:589:PRO:HD3	1.88	0.54
1:A:170:PRO:HB2	1:A:171:PRO:HD2	1.89	0.53
1:A:119:LEU:HD12	1:A:138:GLU:HG3	1.90	0.53
1:A:407:MET:HG3	10:A:723:HOH:O	2.08	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:237:CYS:HA	1:A:381:PHE:O	2.08	0.53
1:A:106:ILE:HG23	1:A:191:LEU:HD11	1.91	0.53
1:A:213:MET:HG2	1:A:273:HIS:CD2	2.44	0.52
1:A:349:PHE:C	1:A:349:PHE:CD1	2.87	0.52
1:A:66:THR:HB	1:A:70:PHE:N	2.25	0.51
1:A:63:GLN:HG3	1:A:71:ARG:HH12	1.74	0.51
1:A:7:GLY:O	1:A:8:ALA:HB2	2.10	0.51
1:A:386:ILE:CG2	1:A:392:ILE:HD13	2.40	0.50
1:A:473:ASN:C	1:A:473:ASN:HD22	2.18	0.50
1:A:518:GLN:O	1:A:522:ILE:HG23	2.11	0.50
1:A:25:THR:O	1:A:184:THR:HG22	2.10	0.50
1:A:348:ARG:HH11	1:A:437:ASN:ND2	2.09	0.50
1:A:298:LYS:NZ	1:A:535:VAL:O	2.28	0.49
1:A:113:PHE:CE2	1:A:115:PRO:HB3	2.48	0.49
4:A:606:HEM:HMC1	4:A:606:HEM:HBC2	1.93	0.49
1:A:45:ARG:CZ	1:A:49:ALA:HB2	2.42	0.49
1:A:255:ARG:CB	9:A:627:PEA:H22	2.42	0.49
1:A:287:TRP:CZ2	1:A:592:SER:HB2	2.48	0.49
1:A:165:PHE:CG	1:A:177:ARG:HD2	2.48	0.49
1:A:544:LEU:O	1:A:547:VAL:HG22	2.12	0.49
1:A:64:ARG:NH2	10:A:920:HOH:O	2.39	0.48
1:A:203:LEU:CB	1:A:213:MET:HE1	2.41	0.48
1:A:202:ARG:HE	1:A:250:GLN:CD	2.22	0.48
1:A:255:ARG:HG2	9:A:627:PEA:H22	1.95	0.48
1:A:216:ASN:HB2	1:A:227:PRO:O	2.14	0.47
1:A:376:LEU:HD13	10:A:942:HOH:O	2.13	0.47
1:A:393:ASP:HB3	1:A:559:ILE:HD11	1.94	0.47
1:A:246:VAL:HG22	8:A:625:EDO:H11	1.95	0.47
1:A:158:MET:HE2	1:A:430:GLY:O	2.15	0.47
1:A:3:GLU:O	1:A:4:VAL:C	2.58	0.47
1:A:46:TRP:CE2	1:A:340:SER:HB3	2.50	0.47
1:A:52:GLU:HG2	1:A:57:VAL:HG12	1.96	0.46
1:A:82:ILE:HD13	1:A:480:LEU:HD23	1.97	0.46
1:A:140:PHE:CE2	1:A:439:GLN:HG3	2.50	0.46
1:A:258:GLU:OE2	4:A:606:HEM:HMB2	1.98	0.46
1:A:276:LEU:HD21	1:A:585:LEU:HD23	1.96	0.46
1:A:424:PRO:C	1:A:426:HIS:H	2.24	0.46
1:A:170:PRO:CB	1:A:171:PRO:HD2	2.45	0.46
1:A:287:TRP:CE2	1:A:291:MET:HE2	2.50	0.46
1:A:349:PHE:CB	1:A:497:ASN:HD21	2.29	0.46
1:A:464:LEU:HA	1:A:481:LEU:HD12	1.98	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:286:HIS:ND1	1:A:286:HIS:N	2.60	0.46
1:A:593:ARG:NH1	1:A:593:ARG:HA	2.31	0.46
1:A:8:ALA:HB3	1:A:167:CYS:O	2.16	0.45
1:A:239:PHE:CD2	1:A:239:PHE:C	2.93	0.45
1:A:162:ARG:CZ	1:A:177:ARG:NH2	2.80	0.45
1:A:197:PRO:HD2	6:A:620:IOD:I	2.86	0.45
1:A:352:MET:HE3	1:A:493:TRP:CZ2	2.51	0.45
1:A:113:PHE:CE1	1:A:255:ARG:HD3	2.51	0.45
1:A:350:GLY:HA3	4:A:606:HEM:CBC	2.45	0.45
1:A:119:LEU:HD12	1:A:138:GLU:CG	2.47	0.45
1:A:193:TYR:OH	1:A:297:ARG:HA	2.16	0.45
1:A:117:THR:HG22	1:A:119:LEU:H	1.82	0.45
1:A:468:GLN:HG2	1:A:477:ALA:HB3	1.98	0.45
1:A:474:LYS:O	1:A:478:LYS:HG3	2.17	0.45
1:A:98:LEU:HD22	10:A:745:HOH:O	2.17	0.44
1:A:352:MET:CB	1:A:407:MET:HG2	2.46	0.44
1:A:76:ARG:HH22	1:A:419:ASN:ND2	2.16	0.44
1:A:276:LEU:CD2	1:A:587:LEU:HD11	2.47	0.44
1:A:378:THR:HG22	8:A:622:EDO:H12	1.98	0.44
1:A:481:LEU:HA	1:A:484:TYR:O	2.17	0.44
1:A:165:PHE:CD2	1:A:177:ARG:HD2	2.52	0.44
1:A:203:LEU:HD11	1:A:252:GLY:HA2	1.99	0.44
1:A:519:PHE:HA	1:A:522:ILE:HG12	2.00	0.44
1:A:8:ALA:N	1:A:9:PRO:HD3	2.32	0.44
1:A:259:GLN:HG2	4:A:606:HEM:CBB	2.48	0.44
1:A:29:ASN:CB	1:A:186:PHE:CZ	3.01	0.44
1:A:7:GLY:O	1:A:8:ALA:CB	2.66	0.43
1:A:236:PRO:HB3	6:A:618:IOD:I	2.89	0.43
1:A:258:GLU:HG3	9:A:627:PEA:C2	2.45	0.43
1:A:117:THR:HG22	1:A:119:LEU:N	2.34	0.43
1:A:227:PRO:HD3	1:A:267:THR:HG23	2.01	0.43
1:A:258:GLU:OE2	4:A:606:HEM:C2B	2.62	0.42
1:A:392:ILE:O	1:A:396:VAL:HG23	2.19	0.42
1:A:221:ASP:OD1	1:A:391:GLY:HA3	2.20	0.42
1:A:232:VAL:O	1:A:232:VAL:HG23	2.19	0.42
1:A:146:LYS:O	1:A:147:ASN:CB	2.66	0.42
1:A:24:ILE:HD13	1:A:24:ILE:HA	1.80	0.42
1:A:352:MET:CE	1:A:493:TRP:CZ2	3.02	0.42
1:A:202:ARG:HG2	1:A:250:GLN:OE1	2.20	0.42
1:A:116:GLU:HG2	1:A:161:PHE:O	2.19	0.41
1:A:235:SER:HB3	1:A:238:GLU:OE2	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:196:GLU:HB3	6:A:620:IOD:I	2.90	0.41
1:A:16:ASP:O	1:A:17:GLU:CB	2.69	0.41
1:A:50:GLU:OE2	1:A:447:PRO:HB3	2.20	0.41
1:A:150:LYS:HE2	1:A:419:ASN:HD22	1.85	0.41
1:A:283:LEU:HD13	1:A:587:LEU:O	2.20	0.41
1:A:331:TYR:HE2	6:A:619:IOD:I	2.74	0.41
1:A:29:ASN:HB3	1:A:186:PHE:CZ	2.56	0.41
1:A:298:LYS:HG2	1:A:536:PHE:CZ	2.56	0.41
1:A:595:ASN:HD22	1:A:595:ASN:N	2.19	0.41
1:A:529:TRP:CD1	1:A:531:GLU:H	2.39	0.41
1:A:564:LEU:HD12	1:A:564:LEU:HA	1.90	0.41
1:A:349:PHE:HA	1:A:497:ASN:HD21	1.86	0.40
1:A:193:TYR:O	1:A:200:ALA:HA	2.22	0.40
1:A:166:VAL:CG1	1:A:180:ILE:HG12	2.52	0.40
1:A:255:ARG:HB3	9:A:627:PEA:C2	2.49	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	592/595 (100%)	546 (92%)	38 (6%)	8 (1%)	9 15

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	8	ALA
1	A	119	LEU
1	A	167	CYS
1	A	2	TRP
1	A	17	GLU
1	A	9	PRO

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Mol	Chain	Res	Type
1	A	319	SER
1	A	4	VAL

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
1	A	516/516 (100%)	500 (97%)	16 (3%)	35 60

All (16) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	4	VAL
1	A	6	CYS
1	A	16	ASP
1	A	24	ILE
1	A	86	LEU
1	A	118	GLU
1	A	125	SER
1	A	177	ARG
1	A	208	SER
1	A	243	THR
1	A	286	HIS
1	A	392	ILE
1	A	473	ASN
1	A	543	SER
1	A	561	LYS
1	A	564	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	147	ASN
1	A	245	HIS
1	A	377	HIS

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Mol	Chain	Res	Type
1	A	419	ASN
1	A	437	ASN
1	A	468	GLN
1	A	473	ASN
1	A	497	ASN
1	A	570	ASN
1	A	574	HIS
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 ⓘ

1 non-standard protein/DNA/RNA residue is modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# $ Z > 2$	Counts	RMSZ	# $ Z > 2$
1	SEP	A	198	1	8,9,10	0.86	0	7,12,14	2.68	5 (71%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
1	SEP	A	198	1	-	5/6/8/10	-

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	198	SEP	O2P-P-OG	3.81	116.61	106.67
1	A	198	SEP	OG-CB-CA	3.80	111.84	108.14
1	A	198	SEP	O3P-P-O1P	2.97	122.42	110.83
1	A	198	SEP	O3P-P-OG	-2.77	99.45	106.67
1	A	198	SEP	OG-P-O1P	-2.18	100.55	106.44

There are no chirality outliers.

All (5) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	198	SEP	N-CA-CB-OG
1	A	198	SEP	C-CA-CB-OG
1	A	198	SEP	CB-OG-P-O2P
1	A	198	SEP	CB-OG-P-O1P
1	A	198	SEP	CB-OG-P-O3P

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates

2 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	1.08	1 (7%)	17,19,21	2.73	7 (41%)
2	NAG	B	2	2	14,14,15	0.96	1 (7%)	17,19,21	1.33	1 (5%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	NAG	B	1	1,2	-	2/6/23/26	0/1/1/1
2	NAG	B	2	2	-	0/6/23/26	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	C1-C2	2.92	1.56	1.52
2	B	1	NAG	O5-C1	-2.65	1.39	1.43

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	1	NAG	O5-C1-C2	-6.90	100.62	111.29
2	B	1	NAG	C2-N2-C7	4.71	129.21	122.90
2	B	1	NAG	C3-C4-C5	-4.21	102.60	110.23
2	B	1	NAG	C4-C3-C2	3.63	116.33	111.02
2	B	1	NAG	O7-C7-N2	2.87	127.06	121.98
2	B	2	NAG	O5-C1-C2	-2.51	107.41	111.29
2	B	1	NAG	O3-C3-C2	-2.47	104.27	109.40
2	B	1	NAG	O6-C6-C5	-2.02	104.46	111.33

There are no chirality outliers.

All (2) torsion outliers are listed below:

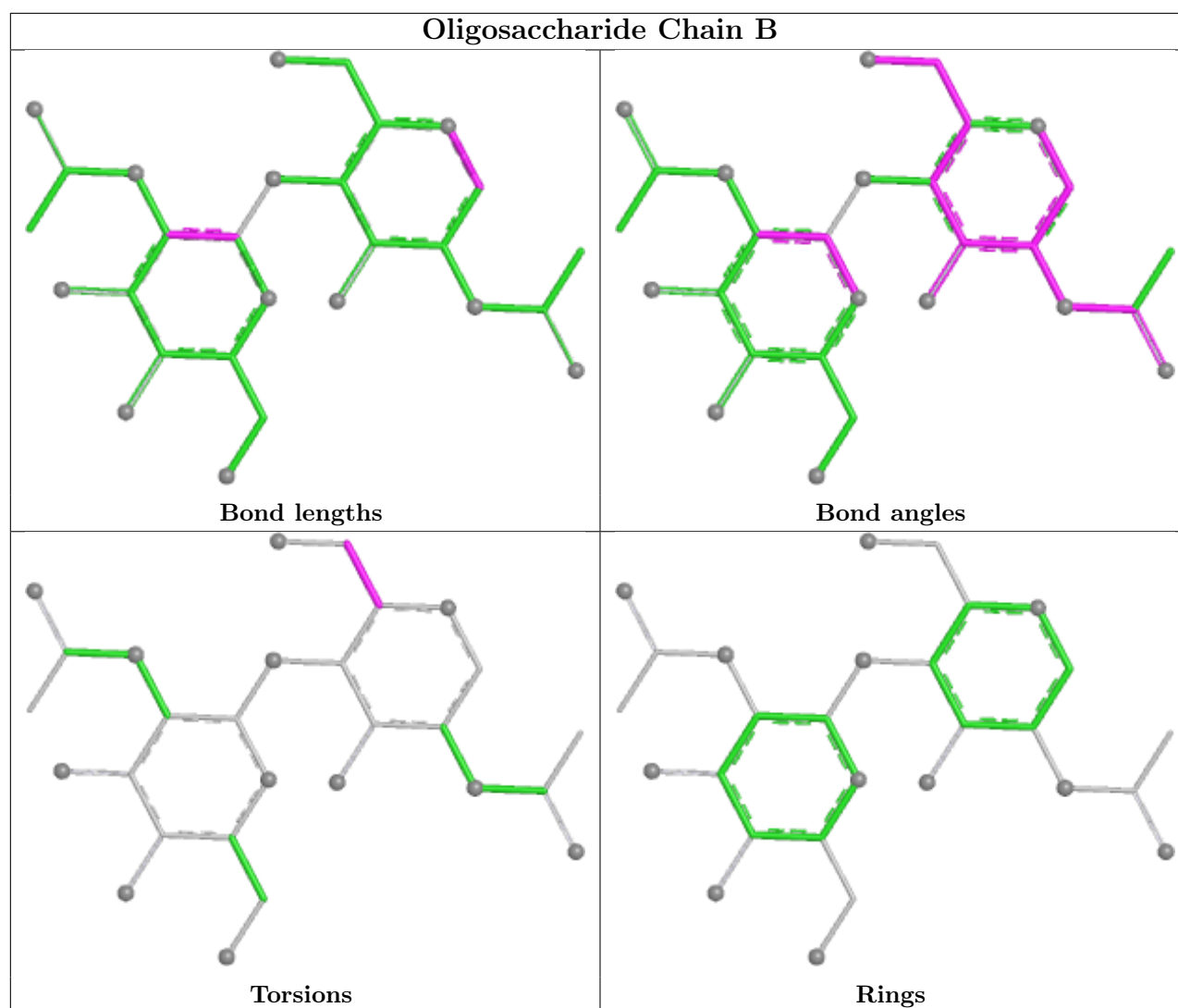
Mol	Chain	Res	Type	Atoms
2	B	1	NAG	C4-C5-C6-O6
2	B	1	NAG	O5-C5-C6-O6

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	2	NAG	1	0

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



5.6 Ligand geometry [i](#)

Of 25 ligands modelled in this entry, 14 are monoatomic - leaving 11 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	SCN	A	621	-	1,2,2	0.28	0	0,1,1	-	-
3	NAG	A	602	1	14,14,15	1.19	0	17,19,21	1.80	7 (41%)
8	EDO	A	624	-	3,3,3	0.55	0	2,2,2	0.19	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
8	EDO	A	622	-	3,3,3	0.13	0	2,2,2	0.91	0
8	EDO	A	625	-	3,3,3	0.26	0	2,2,2	0.97	0
4	HEM	A	606	1,10	50,50,50	2.56	23 (46%)	67,82,82	3.03	30 (44%)
3	NAG	A	605	1	14,14,15	0.45	0	17,19,21	2.45	6 (35%)
9	PEA	A	627	-	9,9,9	1.18	1 (11%)	10,10,10	1.21	1 (10%)
8	EDO	A	623	-	3,3,3	0.59	0	2,2,2	0.35	0
3	NAG	A	601	1	14,14,15	1.15	2 (14%)	17,19,21	1.79	2 (11%)
8	EDO	A	626	-	3,3,3	0.55	0	2,2,2	0.09	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
3	NAG	A	602	1	-	0/6/23/26	0/1/1/1
8	EDO	A	624	-	-	1/1/1/1	-
8	EDO	A	622	-	-	0/1/1/1	-
8	EDO	A	625	-	-	0/1/1/1	-
4	HEM	A	606	1,10	-	4/14/54/54	-
3	NAG	A	605	1	-	2/6/23/26	0/1/1/1
9	PEA	A	627	-	-	0/3/3/3	0/1/1/1
8	EDO	A	623	-	-	0/1/1/1	-
3	NAG	A	601	1	-	0/6/23/26	0/1/1/1
8	EDO	A	626	-	-	0/1/1/1	-

All (26) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	606	HEM	FE-NB	5.56	2.12	1.94
4	A	606	HEM	C3C-C2C	5.08	1.47	1.37
4	A	606	HEM	FE-ND	5.05	2.10	1.94
4	A	606	HEM	C3B-C2B	4.53	1.46	1.37
4	A	606	HEM	CHC-C1C	4.25	1.46	1.38
4	A	606	HEM	C4A-NA	4.00	1.47	1.39
4	A	606	HEM	C4D-ND	-3.79	1.33	1.40
4	A	606	HEM	C3D-C2D	3.60	1.44	1.36
4	A	606	HEM	C1A-NA	3.60	1.46	1.39
4	A	606	HEM	C1C-NC	3.42	1.46	1.39
4	A	606	HEM	CHA-C4D	3.33	1.45	1.38
4	A	606	HEM	C4A-C3A	3.20	1.50	1.43

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	606	HEM	C2A-C3A	3.19	1.47	1.38
4	A	606	HEM	CHB-C1B	3.16	1.44	1.38
4	A	606	HEM	CHD-C4C	3.14	1.44	1.38
4	A	606	HEM	CHB-C4A	3.03	1.46	1.39
4	A	606	HEM	CHC-C4B	2.97	1.46	1.39
4	A	606	HEM	C1D-ND	-2.80	1.33	1.38
4	A	606	HEM	C4C-NC	2.70	1.44	1.39
4	A	606	HEM	C1B-NB	-2.69	1.35	1.40
3	A	601	NAG	O5-C1	-2.64	1.39	1.43
4	A	606	HEM	CHD-C1D	2.62	1.45	1.39
9	A	627	PEA	C2'-C1'	-2.52	1.34	1.38
4	A	606	HEM	C1B-C2B	2.27	1.49	1.44
4	A	606	HEM	C1C-C2C	2.25	1.49	1.45
3	A	601	NAG	C2-N2	-2.15	1.42	1.46

All (46) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	606	HEM	C3B-C4B-NB	9.18	116.06	109.47
4	A	606	HEM	C2A-C1A-NA	7.39	118.35	110.15
4	A	606	HEM	C3B-C2B-C1B	-7.02	101.14	106.41
3	A	605	NAG	C2-N2-C7	6.99	132.26	122.90
4	A	606	HEM	C2B-C1B-NB	6.77	117.62	109.84
4	A	606	HEM	C3D-C4D-ND	6.39	117.18	110.17
4	A	606	HEM	C2D-C1D-ND	6.20	117.07	109.90
3	A	601	NAG	C2-N2-C7	-5.63	115.36	122.90
4	A	606	HEM	C4A-NA-C1A	-5.44	96.96	105.82
4	A	606	HEM	C3C-C2C-C1C	-4.48	102.81	107.05
4	A	606	HEM	C1D-C2D-C3D	-4.36	102.39	106.98
3	A	605	NAG	C1-C2-N2	-3.92	104.26	110.43
4	A	606	HEM	C4C-NC-C1C	-3.81	99.60	105.82
4	A	606	HEM	C4B-C3B-C2B	-3.75	103.83	107.28
3	A	602	NAG	C2-N2-C7	-3.68	117.97	122.90
3	A	601	NAG	O5-C1-C2	-3.62	105.69	111.29
4	A	606	HEM	C3A-C4A-NA	3.59	115.90	110.14
4	A	606	HEM	CHD-C1D-C2D	-3.59	119.36	125.03
4	A	606	HEM	C4D-C3D-C2D	-3.46	101.85	106.89
4	A	606	HEM	CHA-C1A-C2A	-3.46	117.72	125.30
4	A	606	HEM	C2C-C1C-NC	3.45	116.02	109.64
4	A	606	HEM	CHA-C4D-C3D	-3.44	118.88	125.23
4	A	606	HEM	C1B-NB-C4B	-3.31	101.29	105.21
3	A	605	NAG	O5-C5-C6	3.21	113.92	107.66

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	606	HEM	C4D-ND-C1D	-3.14	101.49	105.21
4	A	606	HEM	CMA-C3A-C4A	3.13	130.19	125.42
4	A	606	HEM	CMD-C2D-C1D	2.70	129.26	125.03
4	A	606	HEM	CHC-C4B-C3B	-2.69	119.71	125.07
4	A	606	HEM	CAD-C3D-C4D	2.67	129.35	124.70
3	A	602	NAG	O3-C3-C2	-2.61	103.98	109.40
3	A	602	NAG	C3-C4-C5	-2.47	105.75	110.23
3	A	602	NAG	O5-C1-C2	-2.46	107.48	111.29
4	A	606	HEM	CBD-CAD-C3D	-2.46	105.73	112.53
4	A	606	HEM	CHB-C1B-C2B	-2.40	120.13	126.95
4	A	606	HEM	CHB-C4A-C3A	-2.39	120.48	127.43
4	A	606	HEM	CHC-C1C-C2C	-2.38	120.54	125.49
4	A	606	HEM	CAD-CBD-CGD	-2.30	107.56	113.67
3	A	602	NAG	C8-C7-N2	2.21	119.79	116.12
9	A	627	PEA	C2-C1-N	-2.18	106.73	112.71
3	A	605	NAG	O4-C4-C5	-2.12	104.10	109.32
3	A	602	NAG	C1-O5-C5	2.10	115.00	112.19
4	A	606	HEM	C1A-C2A-C3A	-2.08	103.64	106.87
3	A	605	NAG	O5-C1-C2	2.05	114.46	111.29
4	A	606	HEM	C4A-C3A-C2A	-2.02	104.50	106.82
3	A	605	NAG	C4-C3-C2	2.02	113.97	111.02
3	A	602	NAG	C1-C2-N2	2.01	113.61	110.43

There are no chirality outliers.

All (7) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	605	NAG	C4-C5-C6-O6
3	A	605	NAG	O5-C5-C6-O6
8	A	624	EDO	O1-C1-C2-O2
4	A	606	HEM	CAA-CBA-CGA-O2A
4	A	606	HEM	CAA-CBA-CGA-O1A
4	A	606	HEM	CAD-CBD-CGD-O2D
4	A	606	HEM	CAD-CBD-CGD-O1D

There are no ring outliers.

5 monomers are involved in 29 short contacts:

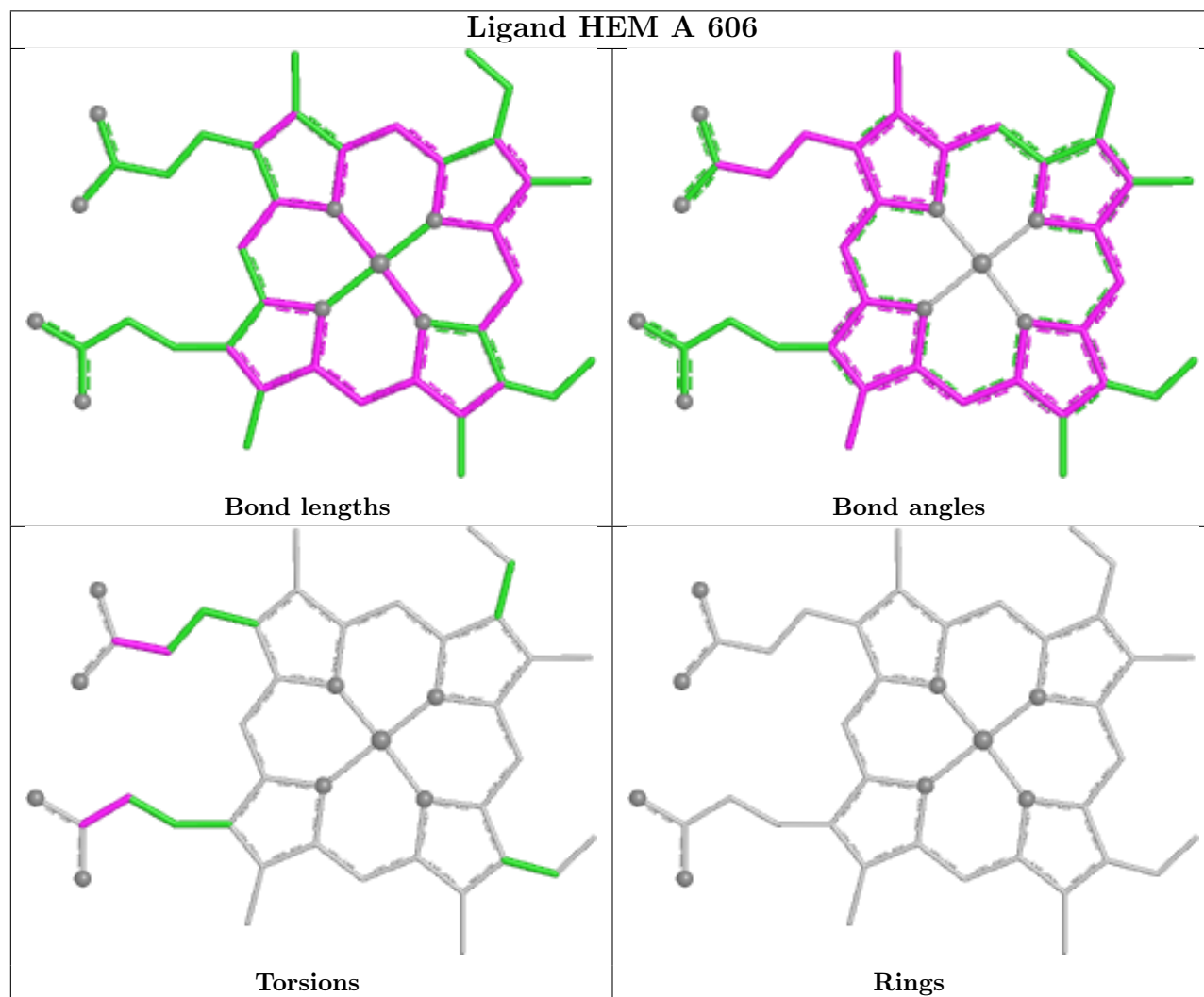
Mol	Chain	Res	Type	Clashes	Symm-Clashes
8	A	624	EDO	1	0
8	A	622	EDO	1	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
8	A	625	EDO	1	0
4	A	606	HEM	15	0
9	A	627	PEA	13	0

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



5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	594/595 (99%)	0.19	22 (3%) 45 41	24, 40, 83, 111	0

All (22) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	12	LEU	5.2
1	A	13	VAL	4.3
1	A	10	VAL	4.2
1	A	11	PRO	3.8
1	A	8	ALA	3.7
1	A	172	TYR	3.6
1	A	4	VAL	3.2
1	A	2	TRP	3.2
1	A	171	PRO	3.1
1	A	223	GLY	3.0
1	A	9	PRO	3.0
1	A	127	VAL	2.6
1	A	119	LEU	2.5
1	A	173	GLN	2.4
1	A	7	GLY	2.3
1	A	363	GLU	2.3
1	A	122	SER	2.2
1	A	164	GLY	2.2
1	A	118	GLU	2.2
1	A	5	GLY	2.1
1	A	117	THR	2.1
1	A	170	PRO	2.1

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

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

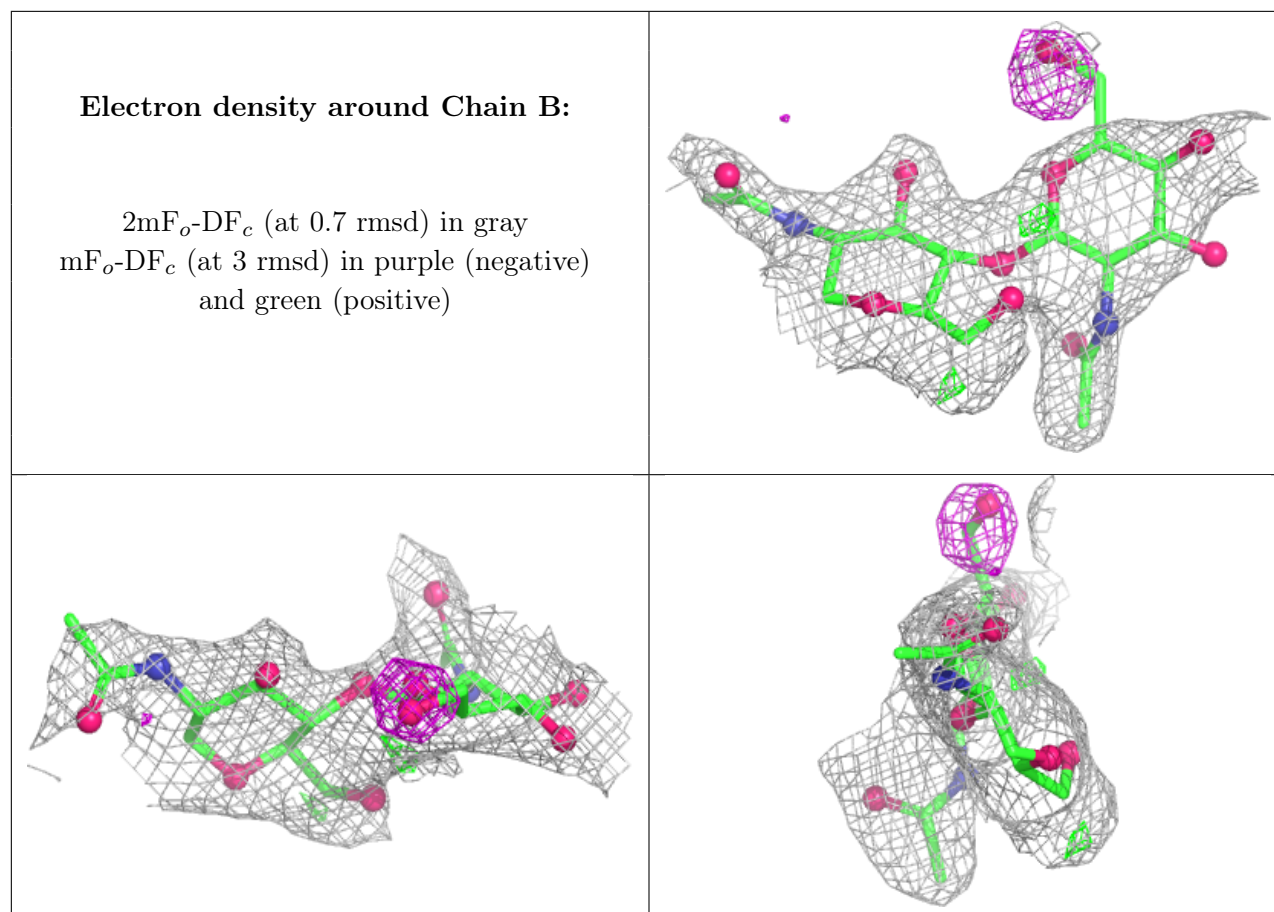
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
1	SEP	A	198	10/11	0.72	0.20	44,50,54,61	0

6.3 Carbohydrates [i](#)

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	NAG	B	2	14/15	0.57	0.17	52,56,59,60	14
2	NAG	B	1	14/15	0.87	0.12	49,52,55,56	0

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



6.4 Ligands [i](#)

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	NAG	A	605	14/15	0.53	0.17	51,55,57,57	14
3	NAG	A	601	14/15	0.73	0.12	45,48,51,51	0
7	SCN	A	621	3/3	0.76	0.16	42,42,43,49	0
8	EDO	A	625	4/4	0.76	0.16	32,34,36,37	0
8	EDO	A	623	4/4	0.80	0.21	34,39,42,43	0
9	PEA	A	627	9/9	0.81	0.20	27,32,37,37	0
8	EDO	A	622	4/4	0.82	0.23	22,22,24,28	0
8	EDO	A	626	4/4	0.85	0.16	34,35,38,38	0
3	NAG	A	602	14/15	0.85	0.08	46,49,51,52	0
6	IOD	A	612	1/1	0.90	0.12	53,53,53,53	1
8	EDO	A	624	4/4	0.93	0.10	35,36,36,37	0
6	IOD	A	619	1/1	0.95	0.11	54,54,54,54	1

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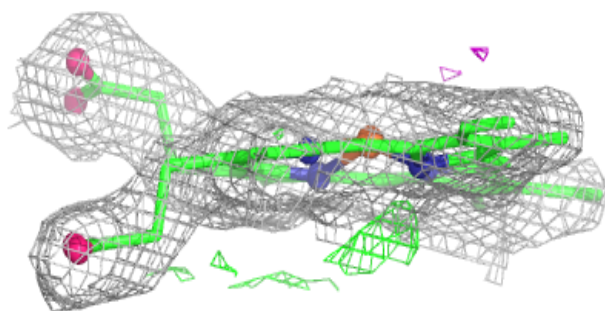
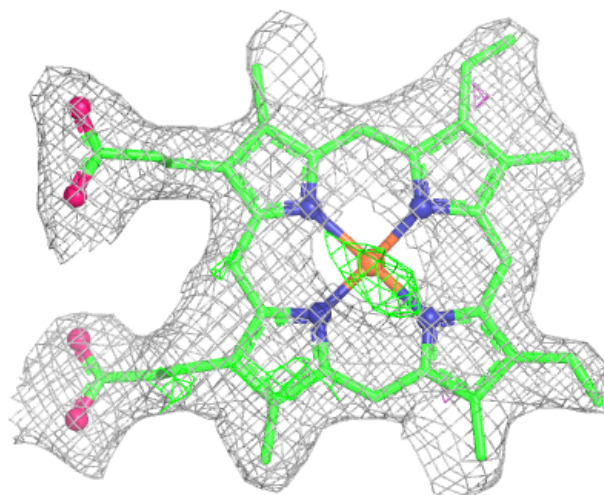
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
4	HEM	A	606	43/43	0.96	0.08	21,24,25,27	0
6	IOD	A	615	1/1	0.96	0.16	51,51,51,51	1
6	IOD	A	616	1/1	0.97	0.11	48,48,48,48	1
6	IOD	A	617	1/1	0.97	0.09	40,40,40,40	1
6	IOD	A	613	1/1	0.98	0.05	38,38,38,38	0
6	IOD	A	614	1/1	0.98	0.10	47,47,47,47	1
6	IOD	A	620	1/1	0.98	0.10	45,45,45,45	0
6	IOD	A	610	1/1	0.98	0.11	44,44,44,44	0
6	IOD	A	609	1/1	0.98	0.04	43,43,43,43	1
5	CA	A	607	1/1	0.99	0.03	27,27,27,27	0
6	IOD	A	611	1/1	0.99	0.09	43,43,43,43	0
6	IOD	A	618	1/1	0.99	0.04	45,45,45,45	1
6	IOD	A	608	1/1	1.00	0.03	34,34,34,34	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.

Electron density around HEM A 606:

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



6.5 Other polymers [i](#)

There are no such residues in this entry.