



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 12, 2026 – 09:34 PM UTC

PDB ID : 6EKY / pdb_00006eky
Title : Direct-evolutioned unspecific peroxygenase from *Agrocybe aegerita*, in complex with 1-naphthol (II)
Authors : Ramirez-Escudero, M.; Sanz-Aparicio, J.
Deposited on : 2017-09-27
Resolution : 1.23 Å(reported)

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

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A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

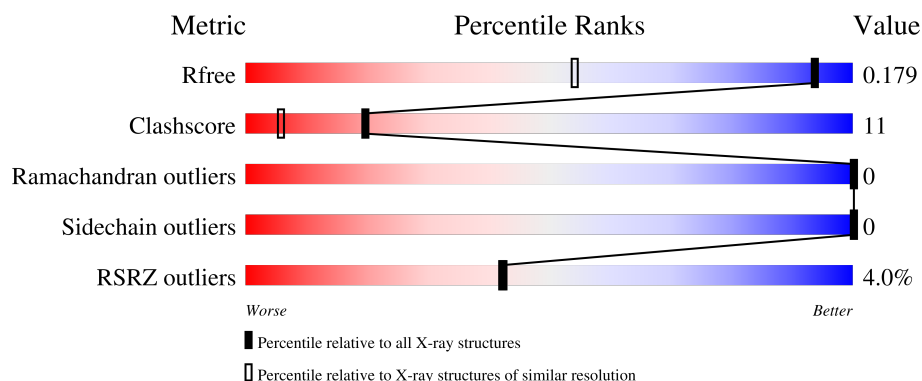
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.23 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	1630 (1.26-1.22)
Clashscore	190562	1668 (1.26-1.22)
Ramachandran outliers	187476	1635 (1.26-1.22)
Sidechain outliers	187428	1633 (1.26-1.22)
RSRZ outliers	180081	1630 (1.26-1.22)

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	328	

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
8	1NP	A	413[A]	-	-	X	-

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Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
8	1NP	A	413[B]	-	-	X	-
8	1NP	A	413[C]	-	-	X	-
8	1NP	A	413[D]	-	-	X	-
8	1NP	A	414[A]	-	-	X	-
8	1NP	A	414[B]	-	-	X	-
8	1NP	A	417	-	-	X	-
9	MOH	A	427	-	-	X	-

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

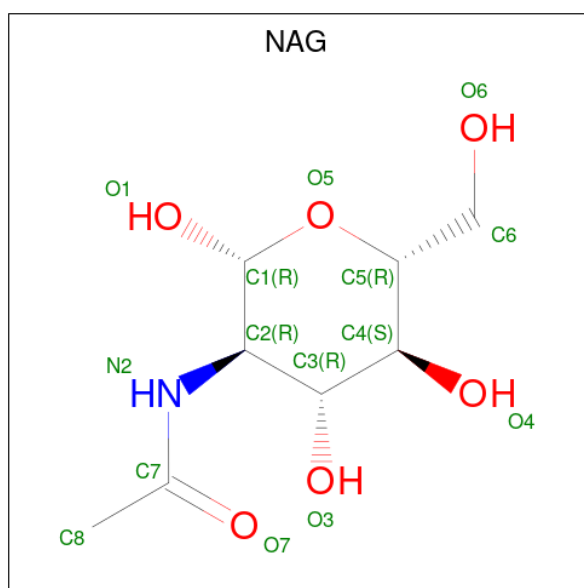
- Molecule 3 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Mg		
			1	1	0	0

- Molecule 4 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

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

- Molecule 5 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: C₈H₁₅NO₆).



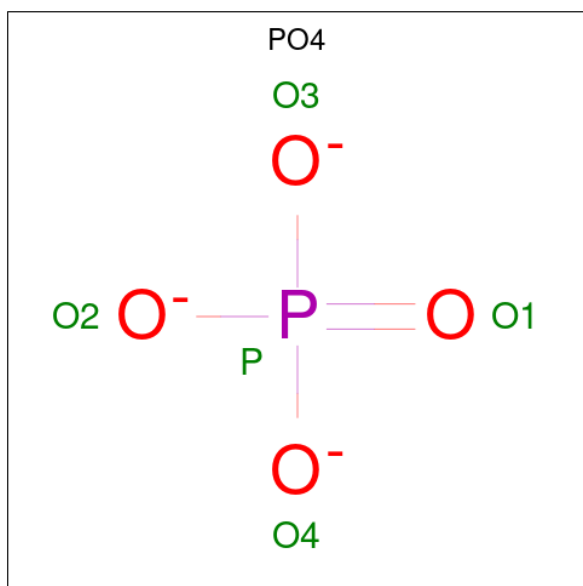
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	A	1	Total	C	N	O		
			14	8	1	5	0	0
5	A	1	Total	C	N	O		
			14	8	1	5	0	0
5	A	1	Total	C	N	O		
			14	8	1	5	0	0
5	A	1	Total	C	N	O		
			14	8	1	5	0	0

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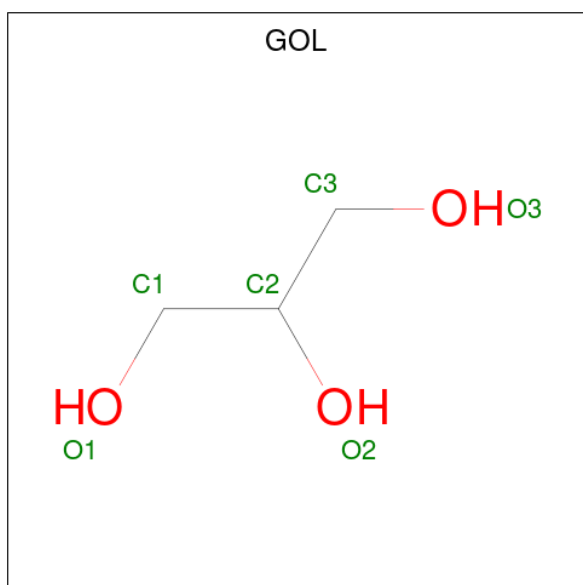
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	A	1	Total	C	N	O	0	0
			14	8	1	5		

- Molecule 6 is PHOSPHATE ION (CCD ID: PO4) (formula: O_4P).



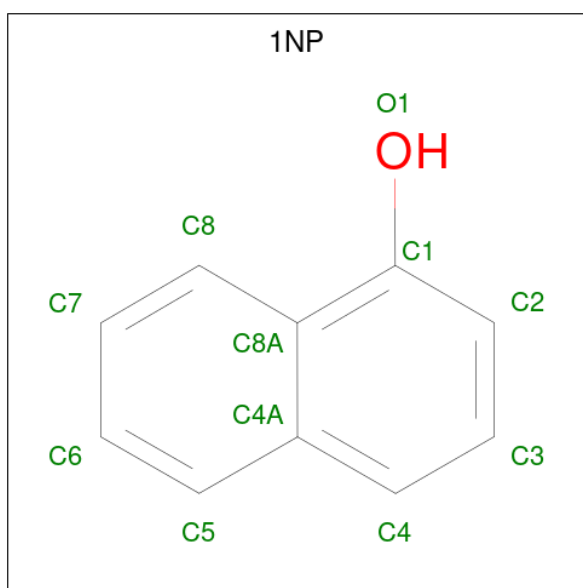
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	A	1	Total	O	P	0	0
			5	4	1		

- Molecule 7 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



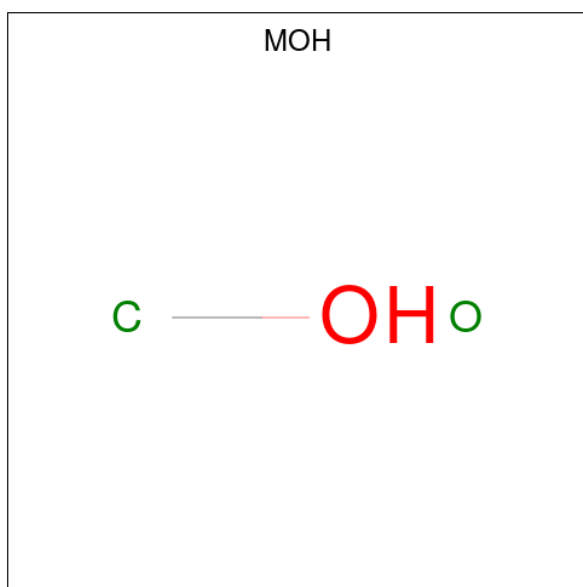
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	A	1	Total	C	O	0	0
			6	3	3		
7	A	1	Total	C	O	0	0
			6	3	3		
7	A	1	Total	C	O	0	0
			6	3	3		
7	A	1	Total	C	O	0	0
			6	3	3		

- Molecule 8 is 1-NAPHTHOL (CCD ID: 1NP) (formula: C₁₀H₈O).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
8	A	1	Total	C	O	0	1
			44	40	4		
8	A	1	Total	C	O	0	1
			22	20	2		
8	A	1	Total	C	O	0	0
			11	10	1		
8	A	1	Total	C	O	0	0
			11	10	1		
8	A	1	Total	C	O	0	0
			11	10	1		

- Molecule 9 is METHANOL (CCD ID: MOH) (formula: CH₄O).



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

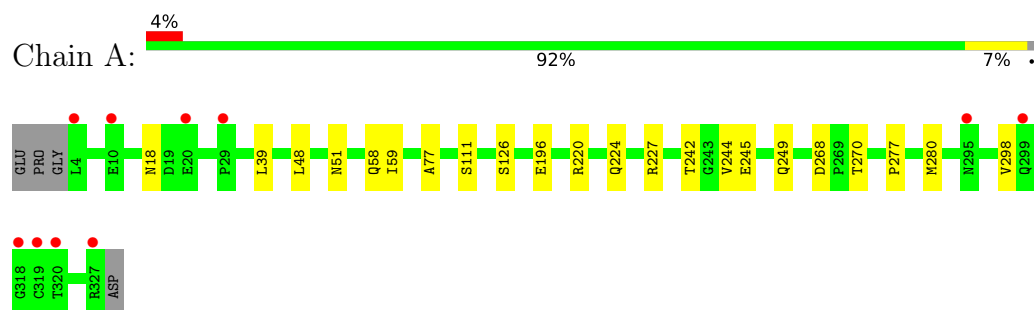
- Molecule 10 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
10	A	388	Total 388	O 388	0	0

3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

- Molecule 1: Aromatic peroxygenase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	51.18Å 57.97Å 60.98Å 90.00° 110.13° 90.00°	Depositor
Resolution (Å)	48.06 – 1.23 48.06 – 1.23	Depositor EDS
% Data completeness (in resolution range)	99.7 (48.06-1.23) 99.7 (48.06-1.23)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.35 (at 1.23Å)	Xtriage
Refinement program	REFMAC 5.8.0158	Depositor
R, R_{free}	0.160 , 0.169 0.170 , 0.179	Depositor DCC
R_{free} test set	4775 reflections (4.92%)	wwPDB-VP
Wilson B-factor (Å ²)	11.1	Xtriage
Anisotropy	0.457	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 38.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	3231	wwPDB-VP
Average B, all atoms (Å ²)	16.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.24% 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, PO4, MOH, MG, CL, NAG, GOL, 1NP

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	0.49	0/2642	0.72	1/3593 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	A	298	VAL	N-CA-C	5.42	116.05	110.36

There are no chirality outliers.

There are no planarity outliers.

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	2572	0	2454	34	0
2	A	43	0	30	10	0
3	A	1	0	0	0	0
4	A	1	0	0	0	0
5	A	70	0	65	1	0
6	A	5	0	0	0	0
7	A	24	0	32	3	0
8	A	99	0	63	39	0
9	A	28	0	0	6	0
10	A	388	0	0	13	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	3231	0	2644	63	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 11.

All (63) 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)
8:A:413[B]:1NP:C3	10:A:504:HOH:O	1.76	1.32
8:A:413[A]:1NP:H2	10:A:504:HOH:O	1.31	1.29
8:A:413[D]:1NP:C6	10:A:504:HOH:O	1.67	1.28
8:A:413[D]:1NP:H6	10:A:504:HOH:O	1.25	1.25
8:A:413[C]:1NP:C7	10:A:504:HOH:O	1.76	1.22
8:A:413[A]:1NP:C2	10:A:504:HOH:O	1.79	1.20
8:A:413[C]:1NP:H7	10:A:504:HOH:O	1.34	1.17
8:A:413[B]:1NP:H3	10:A:504:HOH:O	1.35	1.08
8:A:413[A]:1NP:C6	8:A:414[A]:1NP:H5	1.87	1.05
8:A:413[A]:1NP:H6	8:A:414[A]:1NP:C5	1.95	0.95
8:A:413[A]:1NP:C6	8:A:414[A]:1NP:C5	2.45	0.94
1:A:196:GLU:OE2	8:A:413[D]:1NP:H7	1.69	0.93
1:A:196:GLU:OE2	8:A:413[A]:1NP:H3	1.76	0.84
8:A:413[A]:1NP:C5	8:A:414[A]:1NP:H5	2.08	0.84
1:A:244:VAL:HG13	8:A:416:1NP:H8	1.66	0.78
1:A:196:GLU:OE2	8:A:413[C]:1NP:H6	1.84	0.77
8:A:413[B]:1NP:C7	8:A:414[B]:1NP:H8	2.15	0.77
8:A:413[B]:1NP:C2	10:A:504:HOH:O	2.17	0.76
1:A:280[A]:MET:HE3	8:A:417:1NP:C5	2.19	0.73
8:A:413[B]:1NP:H7	8:A:414[B]:1NP:C8	2.19	0.73
5:A:408:NAG:C7	10:A:509:HOH:O	2.37	0.72
2:A:401:HEM:HBB2	2:A:401:HEM:HHC	1.71	0.72
1:A:280[A]:MET:CE	8:A:417:1NP:C5	2.70	0.69
2:A:401:HEM:CMC	2:A:401:HEM:HBC2	2.22	0.69
2:A:401:HEM:HBC2	2:A:401:HEM:HMC2	1.76	0.68
8:A:413[B]:1NP:C8	8:A:414[B]:1NP:H8	2.25	0.66
8:A:413[B]:1NP:C7	8:A:414[B]:1NP:C8	2.72	0.66
2:A:401:HEM:HHA	2:A:401:HEM:HBA1	1.78	0.64
1:A:196:GLU:OE2	8:A:413[B]:1NP:H2	1.98	0.63
1:A:280[A]:MET:HE3	8:A:417:1NP:C6	2.28	0.62
8:A:413[C]:1NP:C6	10:A:504:HOH:O	2.25	0.61
1:A:196:GLU:CD	8:A:413[D]:1NP:H7	2.26	0.60
1:A:268:ASP:OD1	1:A:270[B]:THR:HG22	2.02	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:196:GLU:CG	8:A:413[D]:1NP:H7	2.34	0.57
1:A:111:SER:OG	10:A:501:HOH:O	2.17	0.57
1:A:18:ASN:HD21	9:A:427:MOH:C	2.19	0.55
1:A:227[A]:ARG:HE	9:A:427:MOH:C	2.20	0.54
1:A:220:ARG:HH21	7:A:412:GOL:C3	2.22	0.53
2:A:401:HEM:HMC2	2:A:401:HEM:CBC	2.41	0.49
1:A:220:ARG:HH21	7:A:412:GOL:H32	1.77	0.49
2:A:401:HEM:HH21	2:A:401:HEM:CBB	2.41	0.49
8:A:413[D]:1NP:C7	10:A:504:HOH:O	2.27	0.48
1:A:77:ALA:HB2	8:A:413[B]:1NP:C5	2.44	0.48
1:A:196:GLU:HG3	8:A:413[D]:1NP:C8	2.44	0.48
1:A:196:GLU:CG	8:A:413[D]:1NP:C7	2.91	0.47
1:A:280[A]:MET:CE	8:A:417:1NP:H5	2.44	0.47
1:A:196:GLU:HG3	8:A:413[D]:1NP:C7	2.44	0.47
2:A:401:HEM:CMC	2:A:401:HEM:CBC	2.93	0.47
1:A:58:GLN:HE21	9:A:425:MOH:C	2.28	0.46
1:A:244:VAL:CG1	8:A:416:1NP:H8	2.41	0.46
1:A:48:LEU:HD21	1:A:59:ILE:HA	1.98	0.46
1:A:277:PRO:O	8:A:417:1NP:H6	2.16	0.45
2:A:401:HEM:HBA1	2:A:401:HEM:CHA	2.44	0.45
1:A:77:ALA:HB2	8:A:413[C]:1NP:O1	2.17	0.45
9:A:427:MOH:C	9:A:428:MOH:C	2.97	0.43
1:A:126:SER:OG	2:A:401:HEM:O2A	2.37	0.42
1:A:39:LEU:HD11	2:A:401:HEM:C2B	2.54	0.42
1:A:242[A]:THR:HG21	9:A:429:MOH:C	2.50	0.41
1:A:245:GLU:O	1:A:249:GLN:HG3	2.20	0.41
1:A:224:GLN:HE21	7:A:411:GOL:C1	2.33	0.41
1:A:77:ALA:HB2	8:A:413[D]:1NP:H4	2.03	0.41
1:A:244:VAL:HG12	8:A:416:1NP:O1	2.21	0.41
1:A:51:ASN:OD1	9:A:430:MOH:C	2.70	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	329/328 (100%)	320 (97%)	9 (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	276/272 (102%)	276 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	18	ASN
1	A	71	ASN
1	A	224	GLN
1	A	254	GLN
1	A	307	ASN

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](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry

Of 36 ligands modelled in this entry, 2 are monoatomic - leaving 34 for Mogul analysis.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# $ Z > 2$	Counts	RMSZ	# $ Z > 2$
8	1NP	A	413[A]	-	12,12,12	1.50	2 (16%)	16,16,16	1.32	3 (18%)
9	MOH	A	428	-	1,1,1	0.19	0	-		
9	MOH	A	426	-	1,1,1	0.07	0	-		
7	GOL	A	410	-	5,5,5	0.36	0	5,5,5	0.85	0
8	1NP	A	414[B]	-	12,12,12	1.58	2 (16%)	16,16,16	0.90	0
5	NAG	A	407	1	14,14,15	0.42	0	17,19,21	0.84	1 (5%)
8	1NP	A	413[B]	-	12,12,12	1.37	1 (8%)	16,16,16	1.12	1 (6%)
5	NAG	A	405	1	14,14,15	0.23	0	17,19,21	0.77	0
9	MOH	A	431	-	1,1,1	0.11	0	-		
9	MOH	A	420	-	1,1,1	0.10	0	-		
7	GOL	A	411	-	5,5,5	0.28	0	5,5,5	0.25	0
9	MOH	A	425	-	1,1,1	0.11	0	-		
9	MOH	A	429	-	1,1,1	0.13	0	-		
5	NAG	A	404	1	14,14,15	0.25	0	17,19,21	0.74	0
9	MOH	A	424	-	1,1,1	0.15	0	-		
9	MOH	A	423	-	1,1,1	0.12	0	-		
5	NAG	A	408	1	14,14,15	0.26	0	17,19,21	1.47	2 (11%)
9	MOH	A	421	-	1,1,1	0.10	0	-		
9	MOH	A	427	-	1,1,1	0.11	0	-		
5	NAG	A	406	1	14,14,15	0.34	0	17,19,21	1.10	2 (11%)
8	1NP	A	413[C]	-	12,12,12	1.48	2 (16%)	16,16,16	1.35	3 (18%)
2	HEM	A	401	3,10,1	50,50,50	1.55	8 (16%)	67,82,82	1.60	12 (17%)
8	1NP	A	416	-	12,12,12	1.73	2 (16%)	16,16,16	1.07	1 (6%)
9	MOH	A	430	-	1,1,1	0.08	0	-		
7	GOL	A	412	-	5,5,5	0.29	0	5,5,5	0.51	0
8	1NP	A	415	-	12,12,12	1.73	2 (16%)	16,16,16	0.82	0
9	MOH	A	419	-	1,1,1	0.10	0	-		
6	PO4	A	409	-	4,4,4	0.94	0	6,6,6	0.46	0
9	MOH	A	422	-	1,1,1	0.07	0	-		
7	GOL	A	432	-	5,5,5	0.49	0	5,5,5	1.12	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
8	1NP	A	417	-	12,12,12	1.71	3 (25%)	16,16,16	1.03	1 (6%)
8	1NP	A	413[D]	-	12,12,12	1.66	2 (16%)	16,16,16	1.03	2 (12%)
9	MOH	A	418	-	1,1,1	0.09	0	-	-	-
8	1NP	A	414[A]	-	12,12,12	1.50	2 (16%)	16,16,16	0.94	2 (12%)

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	1NP	A	413[A]	-	-	-	0/2/2/2
7	GOL	A	410	-	-	0/4/4/4	-
8	1NP	A	414[B]	-	-	-	0/2/2/2
5	NAG	A	407	1	-	0/6/23/26	0/1/1/1
8	1NP	A	413[B]	-	-	-	0/2/2/2
5	NAG	A	405	1	-	1/6/23/26	0/1/1/1
7	GOL	A	411	-	-	4/4/4/4	-
5	NAG	A	404	1	-	2/6/23/26	0/1/1/1
5	NAG	A	408	1	-	0/6/23/26	0/1/1/1
5	NAG	A	406	1	-	2/6/23/26	0/1/1/1
8	1NP	A	413[C]	-	-	-	0/2/2/2
2	HEM	A	401	3,10,1	-	4/14/54/54	-
8	1NP	A	416	-	-	-	0/2/2/2
7	GOL	A	412	-	-	1/4/4/4	-
8	1NP	A	415	-	-	-	0/2/2/2
7	GOL	A	432	-	-	2/4/4/4	-
8	1NP	A	417	-	-	-	0/2/2/2
8	1NP	A	413[D]	-	-	-	0/2/2/2
8	1NP	A	414[A]	-	-	-	0/2/2/2

All (26) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	401	HEM	FE-NB	5.12	2.10	1.94
2	A	401	HEM	FE-NC	4.11	2.08	1.95
8	A	416	1NP	C1-C8A	3.78	1.50	1.42
8	A	415	1NP	C8A-C4A	3.29	1.49	1.43
8	A	413[D]	1NP	C1-C8A	3.24	1.49	1.42
2	A	401	HEM	C4D-ND	-3.24	1.34	1.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	401	HEM	C1B-NB	-3.24	1.34	1.40
8	A	417	1NP	C1-C8A	3.13	1.48	1.42
8	A	413[D]	1NP	C8A-C4A	2.98	1.48	1.43
8	A	414[B]	1NP	C1-C8A	2.97	1.48	1.42
8	A	417	1NP	C8A-C4A	2.96	1.48	1.43
8	A	415	1NP	C1-C8A	2.92	1.48	1.42
8	A	416	1NP	C8A-C4A	2.77	1.48	1.43
8	A	414[B]	1NP	C8A-C4A	2.75	1.48	1.43
8	A	414[A]	1NP	C1-C8A	2.70	1.47	1.42
8	A	413[C]	1NP	C8A-C4A	2.67	1.48	1.43
8	A	414[A]	1NP	C8A-C4A	2.62	1.47	1.43
2	A	401	HEM	FE-ND	-2.58	1.87	1.94
8	A	413[C]	1NP	C1-C8A	2.50	1.47	1.42
8	A	413[A]	1NP	C8A-C4A	2.40	1.47	1.43
8	A	413[B]	1NP	C8A-C4A	2.36	1.47	1.43
2	A	401	HEM	C1C-C2C	-2.31	1.40	1.45
8	A	413[A]	1NP	C8-C8A	-2.25	1.37	1.42
2	A	401	HEM	C1D-ND	-2.14	1.34	1.38
8	A	417	1NP	C5-C4A	-2.11	1.37	1.42
2	A	401	HEM	C4B-NB	-2.07	1.34	1.38

All (30) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	401	HEM	CHC-C4B-NB	4.33	129.08	124.42
5	A	408	NAG	C1-C2-N2	4.29	117.20	110.43
2	A	401	HEM	C1B-NB-C4B	3.93	109.86	105.21
2	A	401	HEM	CHD-C1D-ND	3.91	128.63	124.42
2	A	401	HEM	C3B-C4B-NB	-3.12	107.23	109.47
2	A	401	HEM	CHB-C1B-NB	3.00	128.07	124.37
2	A	401	HEM	CHA-C4D-ND	2.97	128.04	124.37
8	A	416	1NP	O1-C1-C8A	2.90	121.40	116.42
5	A	408	NAG	C1-O5-C5	2.83	115.98	112.19
2	A	401	HEM	C1A-CHA-C4D	-2.74	119.81	126.25
5	A	406	NAG	C2-N2-C7	2.70	126.52	122.90
2	A	401	HEM	C4D-ND-C1D	2.59	108.28	105.21
8	A	413[A]	1NP	C4-C3-C2	2.57	124.94	121.00
8	A	413[C]	1NP	C7-C6-C5	2.54	123.80	120.40
8	A	413[B]	1NP	C8-C8A-C4A	2.49	121.13	117.92
8	A	413[D]	1NP	O1-C1-C8A	2.48	120.67	116.42
8	A	413[C]	1NP	C8-C8A-C4A	2.42	121.05	117.92
2	A	401	HEM	C1C-CHC-C4B	-2.38	120.96	126.02

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	406	NAG	C1-C2-N2	2.35	114.13	110.43
8	A	413[A]	1NP	C8-C8A-C4A	2.32	120.92	117.92
8	A	417	1NP	C2-C1-C8A	-2.25	118.52	120.84
8	A	413[A]	1NP	O1-C1-C8A	2.23	120.25	116.42
2	A	401	HEM	CHD-C4C-NC	2.23	126.88	124.45
8	A	413[D]	1NP	C2-C1-C8A	-2.16	118.61	120.84
8	A	413[C]	1NP	O1-C1-C8A	2.13	120.06	116.42
2	A	401	HEM	C4C-CHD-C1D	-2.11	121.54	126.02
8	A	414[A]	1NP	O1-C1-C8A	2.04	119.92	116.42
2	A	401	HEM	C4C-NC-C1C	2.03	109.13	105.82
8	A	414[A]	1NP	C8-C8A-C4A	2.03	120.54	117.92
5	A	407	NAG	O5-C1-C2	-2.02	108.17	111.29

There are no chirality outliers.

All (16) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	A	406	NAG	C1-C2-N2-C7
7	A	411	GOL	C1-C2-C3-O3
7	A	411	GOL	O1-C1-C2-C3
7	A	432	GOL	O1-C1-C2-C3
7	A	412	GOL	O1-C1-C2-C3
2	A	401	HEM	C1A-C2A-CAA-CBA
7	A	411	GOL	O2-C2-C3-O3
7	A	432	GOL	O1-C1-C2-O2
7	A	411	GOL	O1-C1-C2-O2
5	A	404	NAG	C4-C5-C6-O6
5	A	404	NAG	O5-C5-C6-O6
2	A	401	HEM	CAA-CBA-CGA-O1A
2	A	401	HEM	C3A-C2A-CAA-CBA
5	A	405	NAG	C1-C2-N2-C7
2	A	401	HEM	CAA-CBA-CGA-O2A
5	A	406	NAG	C3-C2-N2-C7

There are no ring outliers.

17 monomers are involved in 59 short contacts:

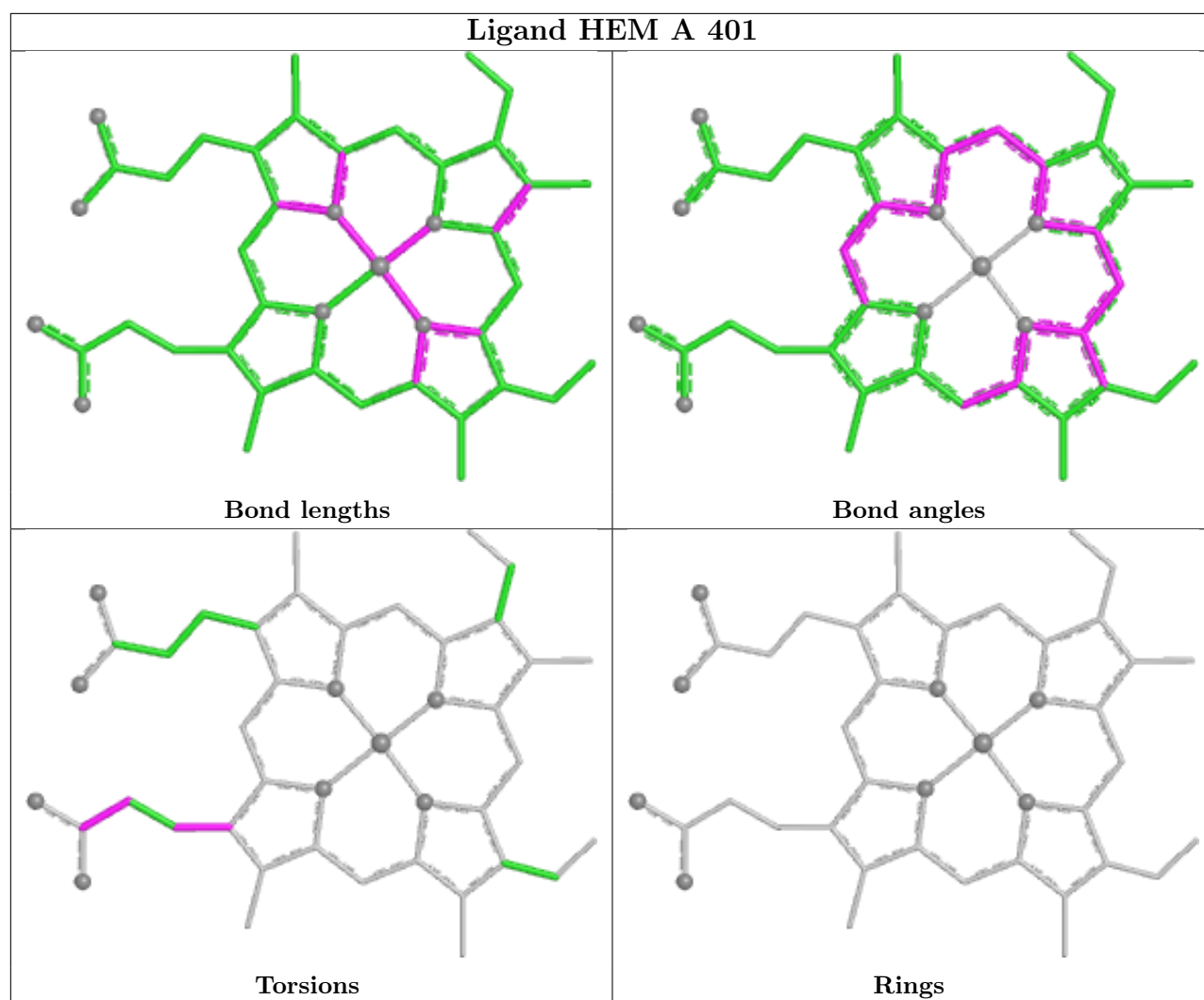
Mol	Chain	Res	Type	Clashes	Symm-Clashes
8	A	413[A]	1NP	7	0
9	A	428	MOH	1	0
8	A	414[B]	1NP	4	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
8	A	413[B]	1NP	9	0
7	A	411	GOL	1	0
9	A	425	MOH	1	0
9	A	429	MOH	1	0
5	A	408	NAG	1	0
9	A	427	MOH	3	0
8	A	413[C]	1NP	5	0
2	A	401	HEM	10	0
8	A	416	1NP	3	0
9	A	430	MOH	1	0
7	A	412	GOL	2	0
8	A	417	1NP	5	0
8	A	413[D]	1NP	10	0
8	A	414[A]	1NP	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	324/328 (98%)	0.28	13 (4%) 42 42	5, 13, 23, 32	10 (3%)

All (13) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	316	ALA	6.0
1	A	317	ALA	4.4
1	A	4	LEU	3.3
1	A	295	ASN	3.1
1	A	10	GLU	2.8
1	A	299	GLN	2.6
1	A	20	GLU	2.6
1	A	327	ARG	2.5
1	A	318	GLY	2.4
1	A	319	CYS	2.2
1	A	320	THR	2.2
1	A	302	LYS	2.1
1	A	29	PRO	2.0

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](#)

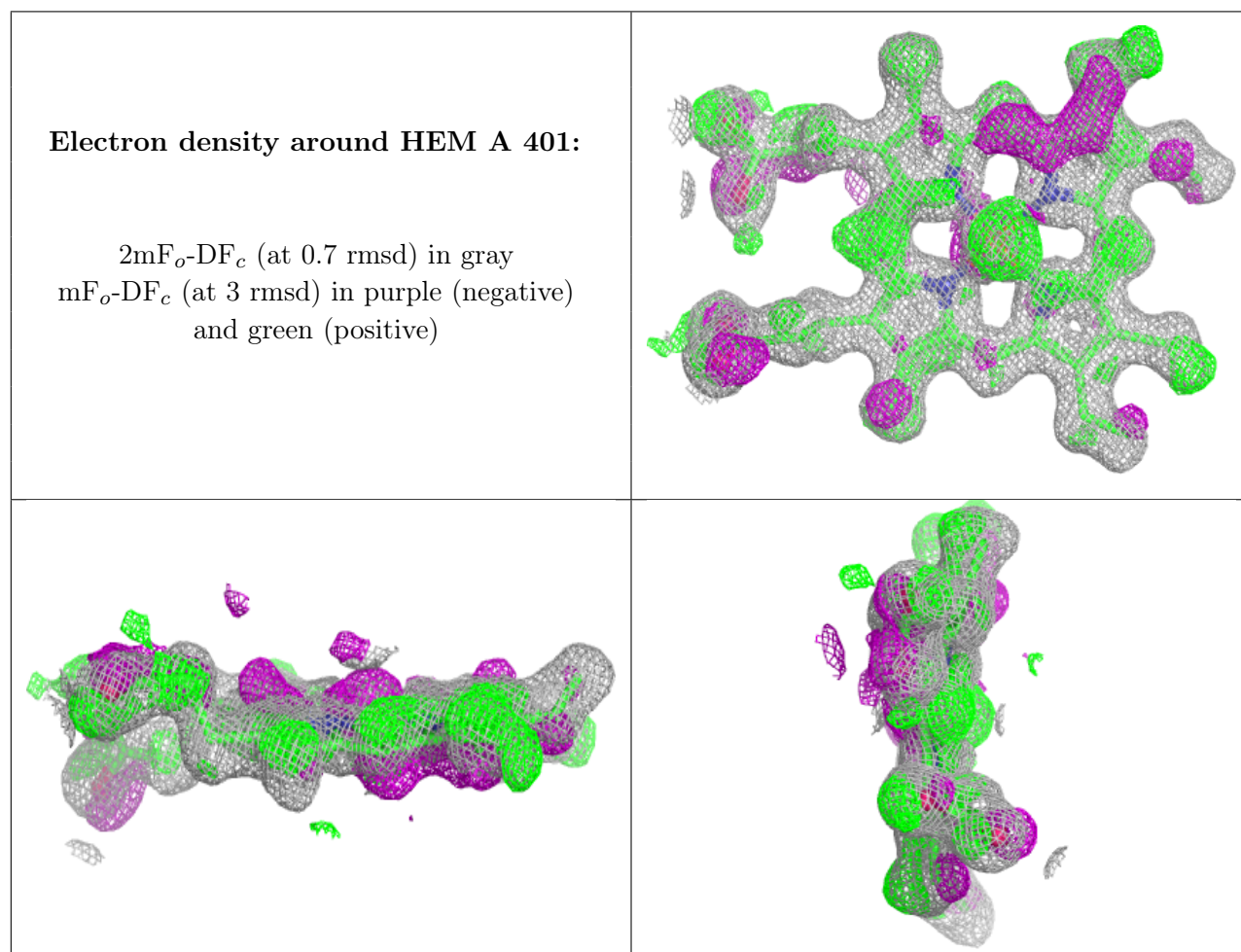
There are no oligosaccharides in this entry.

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
7	GOL	A	432	6/6	0.54	0.23	35,35,35,36	0
9	MOH	A	421	2/2	0.55	0.28	40,40,40,41	0
9	MOH	A	431	2/2	0.57	0.29	30,30,30,30	0
5	NAG	A	404	14/15	0.61	0.16	38,42,44,45	0
9	MOH	A	418	2/2	0.66	0.21	36,36,36,36	0
8	1NP	A	416	11/11	0.67	0.19	38,38,38,38	0
9	MOH	A	422	2/2	0.69	0.21	32,32,32,33	0
7	GOL	A	412	6/6	0.71	0.19	35,37,37,37	0
9	MOH	A	419	2/2	0.72	0.21	40,40,40,40	0
9	MOH	A	423	2/2	0.72	0.22	39,39,39,39	0
8	1NP	A	417	11/11	0.72	0.15	24,25,25,25	0
6	PO4	A	409	5/5	0.73	0.17	28,28,29,30	5
9	MOH	A	427	2/2	0.73	0.26	23,23,23,24	0
9	MOH	A	429	2/2	0.73	0.21	32,32,32,32	0
7	GOL	A	411	6/6	0.73	0.16	34,35,35,35	0
8	1NP	A	413[C]	11/11	0.76	0.17	18,20,20,20	11
8	1NP	A	413[D]	11/11	0.76	0.17	18,20,20,20	11
8	1NP	A	413[A]	11/11	0.76	0.17	22,22,23,23	11
8	1NP	A	413[B]	11/11	0.76	0.17	21,22,22,22	11
9	MOH	A	424	2/2	0.77	0.31	25,25,25,25	0
9	MOH	A	420	2/2	0.82	0.18	30,30,30,30	0
5	NAG	A	408	14/15	0.82	0.13	17,23,25,26	0
9	MOH	A	425	2/2	0.83	0.16	32,32,32,32	0
5	NAG	A	405	14/15	0.84	0.12	25,26,28,28	0
8	1NP	A	414[B]	11/11	0.86	0.12	21,22,22,22	11
8	1NP	A	414[A]	11/11	0.86	0.12	24,25,25,25	11
7	GOL	A	410	6/6	0.87	0.12	18,22,22,23	0
9	MOH	A	430	2/2	0.88	0.15	23,23,23,24	0
8	1NP	A	415	11/11	0.89	0.11	20,22,22,23	0
9	MOH	A	426	2/2	0.89	0.19	25,25,25,26	0
2	HEM	A	401	43/43	0.90	0.13	8,9,12,15	0
9	MOH	A	428	2/2	0.93	0.11	21,21,21,21	0
5	NAG	A	406	14/15	0.96	0.06	11,11,14,14	0
5	NAG	A	407	14/15	0.97	0.05	11,12,14,14	0
4	CL	A	403	1/1	0.99	0.13	30,30,30,30	0
3	MG	A	402	1/1	1.00	0.01	8,8,8,8	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.